Department of Chemistry

Syllabus Module: CBCS

| Semeste | Course | Unit | Sub unit | No. of |
|---------|--------|----------------|-------------------------------------------------|----------|
| r | code – | | | lectures |
| | CEM-G | | | |
| | | Kinetic Theory | Concept of pressure and temperature: | 4 |
| | | of Gases and | Collision of gas molecules: Collision number | |
| Sem - I | CC1/GE | Real mases | and mean free path. Nature of distribution of | |
| | 1 | Keal gases | velocities Maxwell's distribution of speed | |
| | | | and kinetic energy: Average velocity root | |
| | | | mean square velocity and most probable | |
| | | | velocity: Principle of equipartition of energy | |
| | | | Deviation of real gases from ideal behavior: | |
| | | | compressibility factor: Poyle temperature: | |
| | | | Andrew's and Amaget's plate: yen der Waals | |
| | | | Andrew's and Anagat's plots, van der waars | |
| | | | state. Critical constants in terms of yan der | |
| | | | Waals constants: Law of corresponding states | |
| | | | waars constants, Law of corresponding states. | |
| | | Liquids | Definition of Surface tension, its dimension | 2 |
| | | | and principle of its determination using | |
| | | | stalagmometer; Viscosity of a liquid and | |
| | | | principle of determination of coefficient of | |
| | | | viscosity using Ostwald viscometer; Effect of | |
| | | | temperature on surface tension and coefficient | |
| | | | of viscosity of a liquid (qualitative treatment | |
| | | | only) | |
| | | Chemical | Introduction of rate law Order and | 5 |
| | | Kinetics | molecularity: Extent of reaction: rate | 5 |
| | | | constants: Rates of First second and nth order | |
| | | | reactions and their Differential and integrated | |
| | | | forms (with derivation): Pseudo first order | |
| | | | reactions; Determination of order of a reaction | |
| | | | by halfhalf-life and differential method. | |
| | | | Temperature dependence of rate constant: | |
| | | | Arrhenius equation, energy of activation: | |
| | | | | |

| 1 | 1 | 1 | |
|-------|-----------------|------------------------------------------------------------|---|
| | Atomic | Bohr's theory for hydrogen atom (simple | 4 |
| | Structure | mathematical treatment), atomic spectra of | |
| | | hydrogen and Bohr's model, Sommerfeld's | |
| | | model, quantum numbers and their | |
| | | significance, Pauli's exclusion principle, | |
| | | Hund's rule, electronic configuration of many- | |
| | | electron atoms, Aufbau principle and its | |
| | | limitations. | |
| | Chamical | Classification of classents on the basis of | F |
| | Cnemical | Classification of elements on the basis of | 5 |
| | Periodicity | electronic configuration: general | |
| | | characteristics of s-, p-, d- and f-block | |
| | | elements. Positions of hydrogen and noble | |
| | | gases. Atomic and ionic radii, ionization | |
| | | potential, electron affinity, and | |
| | | electronegativity; periodic and group-wise | |
| | | variation of above properties in respect of s- | |
| | | and p- block elements. | |
| | Acids and bases | Brönsted–Lowry concept, conjugate acids and | 4 |
| | | bases, relative strengths of acids and bases, | |
| | | effects of substituent and solvent, differentiating | |
| | | and leveling solvents. Lewis acid-base concept, | |
| | | classification of Lewis acids and bases, Lux- | |
| | | Flood concept and solvent system concept. Hard | |
| | | and soft acids and bases (HSAB concept), | |
| | | applications of HSAB process. | |
| | Fundamentals | Flectronic displacements: inductive effect | 3 |
| | of Organic | resonance and hyperconjugation: nucleophiles | 5 |
| | Chemistry | and electrophiles: reactive intermediates: | |
| | Chemistry | carbocations carbanions and free radicals | |
| | | care of caroanons, caroanons and free fudicals. | |
| | Stonocohomistr | Different types of isomerican association and | 0 |
| | Stereocnemistr | ontical isomerican concert of chirality or d | ð |
| | У | optical activity (up to two corbon stores); | |
| | | optical activity (up to two carbon atoms); | |
| | | asymmetric carbon atom; interconversion of | |
| | | Fischer and Newman representations; | |
| | | enantiomerism and diastereomerism, meso | |
| | | compounds; three and erythre, D and L, cis | |
| | | and <i>trans</i> nomenclature; CIP Rules: <i>R/S</i> (only | |
| | | | |

| r- 11 | GE-2 | Chemical Thermodynami cs: | Intensive and extensive variables; state and path functions; isolated, closed and open systems; zeroth law of thermodynamics; | lectures |
|---------|------------------|-----------------------------------------------------------------|-------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----------|
| Somosto | Dapar | 115:4 | 6.Estimation of Fe(II) and Fe(III) in a given mixture using K ₂ Cr ₂ O ₇ solution. | No. of |
| | | | 4. Estimation of Fe (II) ions by titrating it with K2Cr2O7 using internal indicator. 5. Estimation of Cu (II) ions iodometrically using Na2S2O3. | |
| | | | 2. Estimation of oxalic acid by titrating it with KMnO4. 3. Estimation of water of crystallization in Mohr's salt by titrating with KMnO4. | |
| | CC1/GE 1 (PR) | | 1. Estimation of sodium carbonate and sodium hydrogen carbonate present in a mixture. | 40 |
| | | Nucleophilic Substitution and Elimination Reactions | <i>Nucleophilic substitutions</i> : SN1 and SN2 reactions; eliminations: E1 and E2 reactions (elementary mechanistic aspects); Saytzeff and Hofmann eliminations. | 3 |
| | | | one chiral carbon atoms) and <i>E/Z</i> nomenclature. | |

| | | statement of first law; enthalpy, H; relation | 2 |
|--|-----------------|-------------------------------------------------|---|
| | | between heat capacities, calculations of q, w, | |
| | | ΔU and ΔH for reversible, | |
| | | irreversible and free expansion of gases. | |
| | | Standard states; Heats of reaction; enthalpy of | 1 |
| | | formation of molecules and ions and enthalpy | |
| | | of combustion and its applications; Laws of | |
| | | thermochemistry, Kirchhoff's equations. | |
| | | Statement of the second law of | 3 |
| | | thermodynamics; Concept of heat reservoirs | |
| | | and heat engines; Carnot cycle; | |
| | | Physical concept of Entropy; Entropy change | |
| | | of systems and surroundings for various | |
| | | processes and transformations; Auxiliary state | |
| | | functions (G and A) and Criteria for | |
| | | spontaneity and equilibrium. | |
| | Chemical | Thermodynamic conditions for equilibrium, | 2 |
| | Equilibrium: | degree of advancement; Variation of free | |
| | | energy with degree of advancement; | |
| | | Equilibrium constant and standard Gibbs free | |
| | | energy change; Definitions of KP, KC and Kx | |
| | | and relation among them; | |
| | | van't Hoff's reaction isotherm, isobar and | 1 |
| | | isochore from different standard states; | |
| | | Shifting of equilibrium due to change in | |
| | | external parameters e.g. temperature and | |
| | | pressure; variation of equilibrium constant | |
| | | with addition to inert gas; Le Chatelier's | |
| | | principle | |
| | Redox reactions | Ion-electron method of balancing equation | 3 |
| | | of redox reaction. Elementary idea on | |
| | | standard | |
| | | redox potentials with sign conventions, | |
| | | Nernst equation (without derivation). | |
| | | Influence of | |
| | | | |

| | complex formation, precipitation and | |
|-----------------|---------------------------------------------------------------------------|---|
| | change of pH on redox potentials; formal | |
| | potential. | |
| | 1 | |
| | Feasibility of a redox titration, redox | 1 |
| | potential at the equivalence point, redox | |
| | indicators | |
| | Indicators | |
| Error Analysis | <i>Error analysis</i> : accuracy and precision of | 3 |
| and Computer | quantitative analysis determinate | |
| A pplications | indeterminate, systematic and random errors: | |
| representations | methods of losst squares and standard | |
| | devictions | |
| | deviations. | |
| | <i>Computer applications</i> : general introduction | |
| | to computers different components of a | |
| | computer: hardware and software: input and | |
| | output devices: binery numbers and | |
| | suitherstice later destion to | |
| | arithmetic; introduction to | |
| | computer languages | |
| | | |
| Solutions | Ideal solutions and Raoult's law, deviations | 2 |
| | from Raoult's law – non-ideal solutions; | |
| | Vapour pressure-composition and | |
| | temperature-composition curves of ideal and | |
| | non-ideal solutions: Distillation of solutions: | |
| | Lever rule: A zeotropes Nernst distribution | |
| | law and its applications, solvent extraction | |
| | law and its applications, solvent extraction | |
| Phase | Phases, components and degrees of freedom | 2 |
| Equilibria | of a system, criteria of phase equilibrium: | |
| | Gibbs Phase Rule: Derivation of Clausius – | |
| | Clapevron equation and its importance in | |
| | phase aquilibria: Dhase diagrams of one | |
| | phase equilibria, i hase diagrams of one- | |
| | component systems (water and CO ₂) | |
| Solids | Forms of solids crystal systems unit cells | 2 |
| | Bravais lattice types Symmetry elements: | |
| | Laws of Crystallography Law of constancy | |
| | of interfacial angles. Law of rational indicase | |
| | or internatian angles, Law or rational indices; | |
| | Millor indices of different planes and | |
| | Miller indices of different planes and | |
| | Miller indices of different planes and interplanar distance, Bragg's law; | |

| | Aliphatic | Functional group approach for the following | 3 |
|------|--------------|--------------------------------------------------------|----|
| | Hydrocarbons | reactions (preparations & reactions) to be | |
| | | studied in context to their structures. | |
| | | Alkanes: (up to 5 Carbons). Preparation: | |
| | | catalytic hydrogenation Wurtz reaction | |
| | | Kolhe's synthesis | |
| | | Koloe 5 Syllelesis. | |
| | | Alkenes: (up to 5 Carbons). Preparation: | |
| | | elimination reactions: dehydration of alcohols | |
| | | and dehydrohalogenation of alkyl halides; cis | |
| | | alkenes (partial catalytic hydrogenation) and | |
| | | trans alkenes (Birch reduction). Reactions: | |
| | | addition of bromine, addition of HX | |
| | | [Markownikoff's (with | |
| | | mechanism) and anti-Markownikoff's | |
| | | addition by dration ozonolysis | |
| | | | |
| | | Alkynes: (up to 5 Carbons). Preparation: | |
| | | acetylene from CaC ₂ ; by dehalogenation of | |
| | | tetra halides and dehydrohalogenation of | |
| | | vicinal dihalides. | |
| GE-2 | Practical | Experiment 1: Study of kinetics of acid- | 24 |
| | | catalyzed hydrolysis of methyl acetate | |
| | | Experiment 2: Study of kinetics of | |
| | | decomposition of H_2O_2 (Clock Reaction) | |
| | | | |
| | | Experiment 3: Study of viscosity of | |
| | | unknown liquid (glycerol, sugar) with | |
| | | respect to water. | |
| | | Experiment 4: Determination of solubility | |
| | | of sparingly soluble salt in water, in | |
| | | electrolyte with common ions and in neutral | |
| | | electrolyte (using common indicator) | |
| | | Experiment 5: Preparation of buffer solutions | |
| | | and find the pH of an unknown buffer | |
| | | | |
| | | solution by colour matching method | |

| | | | Experiment 6: Determination of surface | |
|----------|-------|---------------------------------------------------|--------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----------|
| | | | tension of a liquid using Stalagmometer | |
| | | | | N 6 |
| Semes | Paper | Unit | Sub unit | NO. Of |
| ter- III | | | | lectures |
| | GE-3 | Chemical Bonding and Molecular Structure | <i>Ionic Bonding:</i> General characteristics of ionic bonding. Energy considerations in ionic bonding, lattice energy and solvation energy and their importance in the context of stability and solubility of ionic compounds. Statement of Born-Landé equation for calculation of lattice energy, Born-Haber cycle and its applications, polarizing power and polarizability. Fajan's rules, ionic character in covalent compounds, bond moment, dipole moment and percentage ionic character. | 8 |
| | | | <i>Covalent bonding:</i> VB Approach: Shapes of some inorganic molecules and ions on the basis of VSEPR and hybridization with suitable examples of linear, trigonal planar, square planar, tetrahedral, trigonal bipyramidal and octahedral arrangements. Concept of resonance and resonating structures in various inorganic and organic compounds. | 3 |
| | | | MO Approach: Rules for the LCAO method, bonding and antibonding MOs and their characteristics for <i>s-s</i> , <i>s-p</i> and <i>p-p</i> combinations of atomic orbitals, nonbonding combination of orbitals, MO treatment of homonuclear diatomic molecules of 1st and 2nd periods. (including idea of <i>s- p</i> mixing) and heteronuclear diatomic molecules such as CO, NO and NO+. Comparison of VB and MO approaches. | 4 |
| | | Comparative study of p- block elements: | Group trends in electronic configuration, modification of pure elements, common oxidation states, inert pair effect, and their important compounds in respect of the following groups of | 5 |

| | | elements: | |
|--|-----------------------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|---|
| | | i) B-Al-Ga-In-Tl | |
| | | ii) C-Si-Ge-Sn-Pb | |
| | | iii) N-P-As-Sb-Bi | |
| | | iv) O-S-Se-Te | |
| | | v) F-Cl-Br-I | |
| | Transition Elements (<i>3d</i> series) | General group trends with special reference to electronic configuration, variable valency, colour, magnetic and catalytic properties, ability to form complexes and stability of various oxidation states (Latimer diagrams) for Mn, Fe and Cu. | 4 |
| | | Lanthanoids and actinoids: Electronic configurations, oxidation states, colour, magnetic properties, lanthanide contraction, separation of lanthanides (ion exchange method only). | 2 |
| | Coordination Chemistry | Werner's coordination theory, Valence Bond Theory (VBT): Inner and outer orbital complexes of Cr, Fe, Co, Ni and Cu (coordination numbers 4 and 6). Structural and stereoisomerism in complexes with coordination numbers 4 and 6. Drawbacks of VBT. IUPAC system of nomenclature | 5 |
| | ELECTROCH EMISTRY | 1) Ionic Equilibria Strong, moderate and weak electrolytes, degree of ionization, factors affecting degree of ionization, ionization constant and ionic product of water; Ionization of weak acids and bases, pH scale, common ion effect; Salt hydrolysis-calculation of hydrolysis constant, degree of hydrolysis and pH for different salts; Buffer solutions; Solubility and solubility product of sparingly soluble | 4 |

| | | salts - applications of solubility product | |
|--|--------------------------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|---|
| | | principle | |
| | | 2) Conductance | 4 |
| | | Conductance, cell constant, specific conductance and molar conductance; Variation of specific and equivalent conductance with dilution for strong and weak electrolytes; Kohlrausch's law of independent migration of ions; Equivalent and molar conductance at infinite dilution and their determination for strong and weak electrolytes; Ostwald's dilution law; Application of conductance measurement (determination of solubility product and ionic product of water); Conductometric titrations | |
| | | (acid-base) Transport Number and principles Moving-boundary method | |
| | Flootromotivo | Foreday's laws of algotrolysis rules of | E |
| | force | oxidation/reduction of ions based on half-cell potentials, applications of electrolysis in metallurgy and industry; Chemical cells, reversible and irreversible cells with examples; Electromotive force of a cell and its measurement, Nernst equation; Standard electrode (reduction) potential; Electrochemical series; Concentration cells with and without transference, liquid junction potential; pH Determination using hydrogen electrode and quinhydrone; Qualitative discussion of potentiometric titrations (acid- base, redox, precipitation) | J |
| | Aromatic Hydrocarbons | <i>Benzene: Preparation</i> : from phenol, by decarboxylation, from acetylene. <i>Reactions</i> : electrophilic substitution reaction (general mechanism); nitration (with mechanism), halogenations (chlorination and bromination), | 2 |

| | | | and Friedel-Crafts reaction (alkylation and acylation) (up to 4 carbons on benzene). | |
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| | | Organometallic Compounds | Introduction; <i>Grignard reagents</i> : <i>Preparations</i> (from alkyl and aryl halide); Reformatsky reaction. | 1 |
| | | Aryl Halides | <i>Preparation:</i> (chloro- and bromobenzene): from phenol, Sandmeyer reactionand effect of nitro substituent (activated nucleophilic substitution | 1 |
| | GE-3 (PR) | Qualitative semimicro analysis of mixtures containing two radicals. | Cation Radicals: Na ⁺ ,K ⁺ , Ca ^{2+,} Sr ^{2+,} Ba ²⁺ , Al ³⁺ , Cr ³⁺ , Mn ²⁺ /Mn ⁴⁺ , Fe ³⁺ , Co ²⁺ /Co ³⁺ , Ni2+, Cu ²⁺ , Zn ²⁺ , Pb ²⁺ , Sn ²⁺ /Sn ⁴⁺ , NH ⁴⁺ , Anion Radicals: F ⁻ , Cl ⁻ , Br ⁻ , BrO ³⁻ , I ⁻ , IO ³⁻ , SCN ⁻ , S ²⁻ , SO4 ²⁻ , NO ³⁻ , NO ²⁻ , PO4 ³⁻ , AsO4 ³⁻ 'BO ₃ ³⁻ , CrO4 ²⁻ /Cr2O7 ²⁻ | 40 |
| Semes ter- IV | Paper | Unit | Sub unit | No. of lectures |
| | GE-4 | Alcohols, Phenols and Ethers | Alcohols: (up to 5 Carbons). Preparation: 1°-, 2°- and 3°- alcohols: using Grignard reagent, reduction of aldehydes, ketones, carboxylic acid and esters; Reactions: With sodium, oxidation (alkaline KMnO4, acidic dichromate). Diols: Pinacol- pinacolone rearrangement (with mechanism) (with symmetrical diols only). | 3 |
| | | | <i>Phenols: Preparation:</i> cumene hydroperoxide method, from diazonium salts; acidic nature | 2 |

| | | substitution: nitration and halogenations; | |
|---|--------------|-----------------------------------------------|---|
| | | Reimer – Tiemann reaction, Schotten – | |
| | | Baumann reaction, Fries rearrangementand | |
| | | Claisen rearrangement. | |
| | | | |
| | | Ethers: Preparation: Williamson's ether | |
| | | synthesis; Reaction: cleavage of ethers with | |
| | | HI. | |
| | | | |
| | Carbonyl | Aldehydes and Ketones (aliphatic and | 5 |
| | Compounds | <i>aromatic):</i> (Formaldehye, acetaldehyde, | |
| | | acetone and benzaldehyde): Preparation: | |
| | | from acid chlorides, from nitriles and from | |
| | | Grignard reagents; general properties of | |
| | | aldehydes and ketones; Reactions: with HCN, | |
| | | NaHSO3, NH2-G derivatives and with | |
| | | Tollens' and Fehling's reagents; iodoform | |
| | | test: aldol condensation (with mechanism): | |
| | | | |
| | | Cannizzaro reaction (with mechanism), Wittig | |
| | | reaction, benzoin condensation; Clemmensen | |
| | | | |
| | | reduction, Wolff- Kishner reduction | |
| | Quantum | Spectroscopy and its importance in | 8 |
| | Chemistry & | chemistry Wave-particle duality Link | _ |
| | Snectroscony | between spectroscopy and quantum | |
| | Specifoscopy | shemistry. Electromegnetic rediction and its | |
| | | chemistry. Electromagnetic radiation and its | |
| | | interaction with matter. | |
| | | Types of spectroscopy Difference between | |
| | | atomic and molecular spectra | |
| | | atomic and molecular spectra | |
| | | Postulates of quantum mechanics, quantum | |
| | | mechanical operators | |
| | | r | |
| | | Free particle. Particle in a 1-D box | |
| | | (complete solution), quantization, | |
| | | normalization of wave functions, concept of | |
| | | zero-point energy | |
| | | Zero point energy. | |
| | | Rotational Motion: Schrödinger equation of | |
| | | a rigid rotator and brief discussion of its | |
| | | results (solution not required). Quantization | |
| 1 | | results (solution not required). Quantization | |

| | | of rotational energy levels. Microwave (pure rotational) spectra of diatomic molecules. Selection rules. Structural information derived from rotational spectroscopy. <i>Vibrational Motion:</i> Schrödinger equation of a linear harmonic oscillator and brief discussion of its results (solution not | |
|--|-------------------------------------------------|-----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|---|
| | | energy levels. Selection rules, IR spectra of diatomic molecules. | |
| | Carboxylic Acids and Their Derivatives | Carboxylic acids (aliphatic and aromatic): strength of organic acids: comparative study with emphasis on factors affecting pK values; <i>Preparation:</i> acidic and alkaline hydrolysis of esters (<i>B</i> Ac2 and <i>A</i> Ac2 mechanisms only) and from Grignard reagents. <i>Carboxylic acid derivatives</i> (aliphatic): (up to 5 carbons). <i>Preparation:</i> acid chlorides, anhydrides, esters and amides from acids; <i>Reactions:</i> Interconversion among acid derivatives. <i>Reactions:</i> Claisen condensation; Perkin reaction. | 2 |
| | Amines and Diazonium Salts | Amines (aliphatic and aromatic): strength of organic bases; Preparation: from alkyl halides, Hofmann degradation; Reactions: with HNO2 (distinction of 1°-, 2°- and 3°- amines), Schotten – Baumann reaction , Diazo coupling reaction (with mechanism). Diazonium salts: Preparation: from aromatic amines; Reactions: conversion to benzene, phenol, benzoic acid and nitrobenzene. | 3 |

| | | <i>Nitro compounds</i> (aromatic): reduction under different conditions (acidic, neutral and alkaline). | 1 |
|--------------|------------------------------------------------------------------------|------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|----|
| | Amino Acids | <i>Amino Acids: Preparations</i> (glycine and alanine only): Strecker synthesis, Gabriel's phthalimide synthesis; general properties; zwitterion, isoelectric point | 2 |
| | Carbohydrates | <i>Carbohydrates:</i> classificationand general properties; glucose and fructose: constitution; osazoneformation; oxidation-reduction reactions; ascending (Kiliani –Fischer method) and descending (Ruff's method) in monosaccharides (aldoses only); mutarotation | 4 |
| | Crystal Field Theory | Crystal field effect, octahedral symmetry. Crystal field stabilization energy (CFSE), Crystal field effects for weak and strong fields. Tetrahedral symmetry. Factors affecting the magnitude of D. Spectrochemical series. Comparison of CFSE for <i>O_h</i> and <i>T_d</i> complexes, Tetragonal distortion of octahedral geometry. Jahn-Teller distortion, Square planar coordination | 4 |
| GE-4 (PR) | 1.Qualitative Analysis of Single Solid Organic Compound(s) | Experiment A: Detection of special elements (N, Cl, and S) in organic compounds. Experiment B: Solubility and Classification (solvents: H2O, dil. HCl, dil. NaOH) Experiment C: Detection of functional groups: Aromatic-NO2, Aromatic -NH2, - COOH, carbonyl (no distinction of –CHO and >C=O needed), - OH (phenolic) in solid organic compounds. Experiments A - C with unknown (at least 6) solid samples containing not more than two of the | 40 |

| | above type of functional groups should be done. | |
|------------------------------------------------------|----------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------------|--|
| 2.Identification of a pure organic compound | <i>Solid compounds</i> : oxalic acid, tartaric acid, succinic acid, resorcinol, urea, glucose, benzoic acid and salicylic acid. <i>Liquid Compounds</i> :methyl alcohol, ethyl alcohol, acetone, aniline, dimethylaniline, benzaldehyde, chloroform and nitrobenzene | |