

## 。 GDAŃSK UNIVERSITY OF TECHNOLOGY

## Subject card

Subject name and code	MATHEMATICAL METHODS IN CHEMISTRY, PG_00038882								
Field of study	Chemistry								
Date of commencement of studies	February 2025		Academic year of realisation of subject			2024/2025			
Education level	second-cycle studies		Subject group			Obligatory subject group in the field of study			
Mode of study	Full-time studies		Mode of delivery			at the university			
Year of study	1	Language of instruction		า	Polish				
Semester of study	1	ECTS credits		2.0					
Learning profile	general academic profile		Assessment form			exam			
Conducting unit	Department of Physical Chemistry -> Faculty of Chemistry								
Name and surname	Subject supervisor prof. dr hab. inż. Jacek Czub								
of lecturer (lecturers)	Teachers		prof. dr hab. i	ab. inż. Jacek Czub					
			dr inż. Mateu	sz Kogut					
Lesson types and methods	Lesson type	Lecture	Tutorial	Laboratory	Projec	Project Seminar		SUM	
of instruction	Number of study hours	15.0	15.0	0.0	0.0	0.0		30	
	E-learning hours included: 0.0								
Learning activity and number of study hours	Learning activity	Participation in didactic classes included in study plan		Participation in consultation hours		Self-study		SUM	
	Number of study hours	30		5.0		15.0		50	
Subject objectives	The aim of the course is to provide students with general knowledge and skills in advanced mathematical methods used in chemistry and related fields, with a special emphasis on linear algebra, vector analysis, Fourier series and transform, probability theory and numerical methods.								
Learning outcomes	Course out	utcome Subject outcome Method of verifica			rification				

	Lecture					
Subject contents	<ul> <li>Lecture:</li> <li>Linear algebra (operations on vectors and matrices, inner product axioms, systems of linear equations, multiple linear regression, four fundamental subspaces, orthogonal bases, orthogonal (unitary) matrices, eigenvalue problem, diagonalization, properties of symmetric (hermitian) matrices, similarity transformation, spectral decomposition of a matrix, positive-definite matrices, singular value decomposition, introduction to LCAO-MO method, Hueckel method, normal mode analysis, principal component analysis (parametric curves and surfaces, vectors tangent and normal to curves and surfaces, arc length, total differential and total derivative, chain rule, gradient and gradient operator, Lagrange multipliers, directional derivative, scalar and vector fields, integration change of variables, Jacobian, line and surface integrals, properties of a gradient field, curl of a vector field, flux of a vector field, Stokes theorem, Green theorems, divergence of a vector field, continuity equation, Gauss theorem, laplacian, diffusion equation, Poisson equation, Laplace equation, Poisson-Boltzmann equation, vector identities)</li> <li>Fourier analysis (brief introduction to Sturm-Liouville theory, representing functions in orthogonal function bases, orthogonal polynomials, Fourier series and its convergence, Fourier basis, complex form of Fourier series, Fourier transform, inverse Fourier transform, Dirac delta-function, application of Fourier transform in spectroscopy and crystallography, discrete Fourier transform, fast Fourier transform, convolution theorem and its applications, Nyquist-Shannon sampling theorem)</li> <li>Stochastic processes theory (probability density, cumulative distribution function, stochastic processes and their properties, stationary stochastic processes, autocorrelation, white noise, Monte Carlo methods, Langevin equation and Brownian dynamics, introduction to Markov processes, Markov matrices and their spectral decomposition, Markov state models in chemical kinetics and</li></ul>					
	<ol> <li>Introduction to Octave program representation, operations on</li> <li>Numerical approach to solving processes, simple models de cooperativity at the molecular</li> <li>Eigenvalue problem: principal aromaticity, effect of heteroate</li> <li>Eigenvalue problem: normal r spectra</li> <li>Basics of vector analysis: nun optimization of the structure o</li> <li>Selected topics in Fourier ana</li> </ol>	g ordinary differential equations and to scribing the kinetics of bistable reactio level. component analysis & Hueckel metho om on energy levels, fundamentals of e node analysis in the context of IR spec nerical approach to optimization, minim	Markov modeling of (bio)chemical ns, dimerization thermodynamics, id, quantum-chemical basis of electronic excitations troscopy, numerical prediction of IR nization of multivariable function, nsform to analysis of time series, in			
	transform of images	nsform of free induction decay data fro	im an NMR spectrometer; Fourier			
Prerequisites and co-requisites	Basic background in calculus and	linear algebra.				
Assessment methods	Subject passing criteria	Passing threshold	Percentage of the final grade			
and criteria	Practicals	50.0%	70.0%			
	Lecture	50.0%	30.0%			
Recommended reading	Basic literature	<ol> <li>Donald A. McQuarrie "Mathematical Methods for Scientists and Engineers"</li> <li>Erich Steiner "The Chemistry Maths Book"</li> <li>Henry Margenau "The Mathematics of Physics and Chemistry"</li> <li>Zbigniew Skoczylas, "Elementy analizy wektorowej: teoria, przykłady, zadania" (in polish only)</li> </ol>				
	Supplementary literature	<ol> <li>Gilbert Strang "LInear Algebra and Its Applications", 4th ed.</li> <li>George B. Arfken, Hans J. Weber "Mathematics for Physicists", 7th ed.</li> </ol>				
	eResources addresses	Adresy na platformie eNauczanie: Metody matematyczne w chemii (lato 2024_25, Chemia II sem.1) - Moodle ID: 38736 https://enauczanie.pg.edu.pl/moodle/course/view.php?id=38736				
Example issues/ example questions/ tasks being completed	<ol> <li>Create a routine z(v) that will function z = f(x,y), e.g. z = x<sup>2</sup> the vector of coordinates v = program which will do the follomove point P by a given step in value of z = f(x,y), (5) repeated. Use the toeplitz() function to conthe Hückel theory. Calculate the Hückel theory.</li> </ol>	d in the MATLAB/Octave environment: take a vector of arguments, $v = [x y]$ , a + 4y <sup>2</sup> . Based on this routine, create and [x y] and return the normalized gradien owng: (1) read the starting point P, (2) along the direction of negative gradien at steps 2-4 until the value of the function create a tridiagonal matrix describing a he electronic energy of the conjugated oresent benzene instead of hexatriene.	nd return the value of an arbitrary other routine grad(v) that will take t, grad(v) = [gx gy]. Then, write a calculate the gradient at point P, (3) t, (4) compute the resulting change on ceases to decrease. hexatriene molecule according to $\pi$ -system. Repeat the procedure			

Work placement
----------------

Document generated electronically. Does not require a seal or signature.