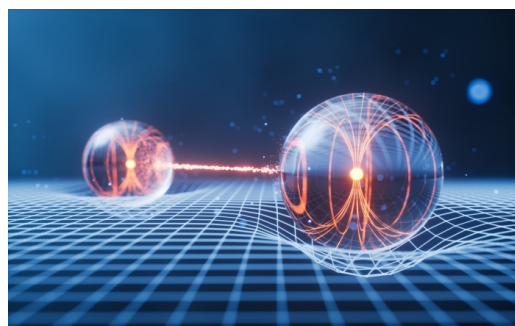




21st International Conference on Density Functional Theory and its Applications (DFT2026)



Abu. 24 - Abu. 27 2026

Kod. Z64-26

Mod.:

Aurrez aurrekoa

Edizioa

2026

Jarduera mota

Workshop

Data

Abu. 24 - Abu. 27 2026

Kokalekua

Miramar Jauregia

Hizkuntzak

Ingelesa

Balio akademikoa

40 ordu

Webgunea

<https://dft2026.com>

Antolakuntza Batzordea



Fundación
BBVA



Azalpena

The conference covers a broad range of topics in the field of density functional theory (DFT), from the latest theoretical developments to cutting-edge chemical and physical applications, bringing together the DFT community from around the world.

The scientific program spans four days and includes seven plenary talks and twelve invited lectures by internationally renowned speakers, who will present the current state of the art in DFT. In addition, 32 contributed talk slots will offer participants the opportunity to present their research at the world's largest platform for DFT developers and users.

Scientific Themes: New fundamental developments in density functional theory, Advances in exchange–correlation functionals, Ensemble DFT and multireference DFT, DFT and machine learning, Novel areas of application of DFT.

ORGANIZING COMMITTEE

- **Prof. Eduard Matito** - Donostia International Physics Center (DIPC)
- **Dr. Eloy Ramos-Cordoba** - Institute for Advanced Chemistry of Catalonia (IQAC-CSIC)
- **Dr. Irene Casademont Reig** - Donostia International Physics Center (DIPC)
- **Prof. Jesus M. Ugalde** - Euskal Herriko Unibertsitatea (EHU) and Donostia International Physics Center (DIPC)
- **Dr. Roberto Álvarez Boto** - Euskal Herriko Unibertsitatea (EHU)
- **Dr. Ishita Bhattacharjee** - Donostia International Physics Center (DIPC)

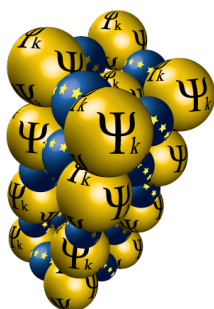
Helburuak

Bring together the scientific community working on density functional theory.

Bring leading experts in the field to Donostia and promote the scientific exchange with people at the DIPC, EHU and other BERCs

Place Donostia as an important hub working on computational chemistry and molecular physics

Ikastaroaren laguntzaile espezifikoak



Zuzendaritza



Eduard Matito Gras

DIPC

Dr. Eduard Matito is an Ikerbasque professor at the DIPC, where leads a group of 12 people. His works have contributed significantly to chemical bonding analysis and aromaticity. One of his most fascinating studies is the importance of the delocalization error in aromaticity, recognizing the spurious classification of recently synthesized molecules as aromatic [4,7]. He is part of the European Committee on Chemical Bonding (ECCB), chair of the DFT 2026 conference, co-organizer of ESCB2 and two editions of CTTC, and part of the scientific board of Aromaticity conference. He was awarded the Miguel A. Blanco Early Career International Prize for his works on chemical bonding and aromaticity. Since 2020, he is in the world's top 2% Scientists.

Matrikula prezioak

REGISTRATION FEES	2026-04-15 ARTE	2026-07-31 ARTE
Fee Waiver	0 EUR	0 EUR
Scientific Committee	250,00 EUR	250,00 EUR
Student Reduced Fee	300,00 EUR	400,00 EUR
Standard Fee	450,00 EUR	550,00 EUR

Kokalekua

Miramar Jauregia

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Gipuzkoa