

# WP2.4: Research infrastructures offering theoretical support for experiments

**Manuela Rodríguez-Gallardo**

On behalf Gert Aarts (ECT\*)

Universidad de Sevilla

Krakow, 10th October 2023

## → Leading Task 4: Gert Aarts

- ⇒ 4.1) ECT\*, European Centre for Theoretical Studies in Nuclear Physics and Related Areas  
Leading Gert Aarts



- ⇒ 4.2) Theo4Exp Virtual Access Infrastructure  
Leading Manuela Rodríguez-Gallardo



INSTYTUT FIZYKI JĄDROWEJ  
IM. HENRYKA NIEWODNICZAŃSKIEGO  
POLSKIEJ AKADEMII NAUK



UNIVERSITÀ  
DEGLI STUDI  
DI MILANO

# ECT\*, Trento (Italy)



# ECT\* mission



- ✓ to be a Centre at the frontline of research in theoretical nuclear physics
- ✓ to promote active contacts between theory and experiments, and to related areas of research
- ✓ to further the training of young researchers
- established in **1993 (30 years ago!)**
- Institutional member of ESF-Expert Committee NuPECC (Nuclear Physics European Collaboration Committee)
- community-driven, bottom-up approach

## SYMPOSIUM IN OCCASION OF THE 30TH ANNIVERSARY OF ECT\*

OCTOBER 4TH, 2023

ECT\* | Aula Leonardi

ECT\* | Villazzano, Trento,  
Strada delle Tabarelle 286

Info: [staff@ectstar.eu](mailto:staff@ectstar.eu)

### PROGRAMME



- 9.00 - 10.00 **Welcome and introduction by invited guests**
- 10.00 - 10.45 **Computing the Heart of Matter**  
Sonia Bacca (University of Mainz)
- 10:45 - 11:15 Coffee break
- 11:15 - 12:00 **Ab-initio Nuclear Physics in Trento**  
Francesco Pederiva (University of Trento)
- 12:00 - 14:00 Lunch at Villa Tambosi
- 14:00 - 14:45 **Color Glass Condensate in Collider Physics**  
Dionysios Triantafyllopoulos (ECT\*)
- 14:45 - 15:30 **Cosmic Laboratories for Nuclear Physics**  
Almudena Arcones (TU Darmstadt)
- 15:30-16:00 Coffee break
- 16:00 - 16:45 **Cosmic Thirty Years of Education and Research on Nuclear Many-Body Physics at the ECT\*; from traditional methods to quantum computing and machine learning**  
Morten Hjorth-Jensen (Michigan State University, USA and University of Oslo, Norway)
- 16:45 - 17:30 **Quantum Fractals**  
Ubirajara Van Kolck (JCLab Orsay & University of Arizona)
- 18:00 Dinner at Villa Tambosi

The ECT\* is part of the Fondazione Bruno Kessler. The Centre is funded by the Autonomous Province of Trento, funding agencies of EU Member and Associated states, and by INFN-TIFPA and has the support of the Department of Physics of the University of Trento.



7 OTTOBRE 2023

ORE 17.00

Aula Grande - Fondazione Bruno Kessler  
Via Santa Croce 77 - Trento

### LAURA FABBIIETTI IL LUNGO VIAGGIO DEGLI ANTINUCLEI

Gli antinuclei sono immagini speculari dei normali nuclei atomici, con la stessa massa ma carica opposta. Non esistono fonti naturali di antinuclei sulla Terra, ma possono essere prodotti in laboratorio presso grandi acceleratori di particelle. Gli antinuclei vengono cercati anche nello spazio, perché potrebbero essere la chiave di uno dei più grandi misteri della fisica: la materia oscura. La materia oscura è onnipresente e rappresenta cinque volte la massa di tutta la materia che possiamo osservare sotto forma di stelle nel cielo, pianeti e tutto il gas intermedio nelle galassie. Non è però possibile vedere o toccare la materia oscura perché non interagisce con la luce o con le forze elettriche. Gli antinuclei offrono un nuovo modo di guardare nello spazio per cercare la materia oscura in quanto essa può interagire per creare antinuclei altrimenti quasi assenti. Come possiamo trovare queste particelle nello spazio? Quali proprietà dobbiamo conoscere? E da dove possono provenire gli antinuclei nello spazio? Durante l'evento approfondiremo queste domande e seguiremo il lungo viaggio degli antinuclei dal centro della nostra galassia alla Stazione Spaziale Internazionale nello spazio.

EVENTO PUBBLICO IN OCCASIONE DEL TRESANTESIMO ANNIVERSARIO DI ECT\*

INFORMAZIONI  
[staff@ectstar.eu](mailto:staff@ectstar.eu)

ECT\* fa parte della Fondazione Bruno Kessler. Il Centro è finanziato dalla Provincia Autonoma di Trento, dalle agenzie di finanziamento degli Stati membri e associati dell'UE, dall'INFN-TIFPA e gode del supporto del Dipartimento di Fisica dell'Università di Trento.

La Prof.ssa LAURA FABBIIETTI si è laureata in fisica all'università Statale di Milano ed è docente di fisica nucleare presso l'Università Tecnica di Monaco (TUM). Nel 2007 ha diretto un gruppo di ricerca Helmholtz junior alla TUM in stretta collaborazione con la Gesellschaft für Schwerionenforschung (GSI) e, dal 2008, un gruppo di ricerca junior dell'Universe Cluster of Excellence. Nel 2011 ha vinto la cattedra presso la TUM, dove dirige la divisione di Density and Strange Hadronic Matter. Oggi è uno degli scienziati di punta del Cluster of Excellence ORIGINS e dell'Area Speciale di Ricerca 1258. La Prof.ssa Fabbietti conduce i suoi esperimenti all'LHC del CERN nell'ambito della collaborazione ALICE.



# ECT\* Scientific Board

membership suggested  
by ECT\* associates  
3-year term

[Almudena Arcones](#) | TU Darmstadt (D)

[Constantia Alexandrou](#) | University of Cyprus & The Cyprus Institute (CY)

[David Kaplan](#) | University of Washington (USA)

[Denis Lacroix](#) | CNRS/IN2P3 (F)

[Marek Lewitowicz](#) | NuPECC/GANIL (F)

[Alexandre Obertelli](#) | TU Darmstadt (D)

[Barbara Pasquini](#) | University of Pavia (I)

[Vittorio Somà](#) | CEA Saclay (F)

[Urs Wiedemann](#), Board Chair | CERN-TH (CH)

Ex officio: [Albino Perego](#) | University of Trento (I)

**2023 PROGRAMME OF ACTIVITIES**

<b>JANUARY</b> 30.1-2.2	Structure and topology of RNA in living systems L. TUBIANA (University of Trento), S. PASQUALI (University Paris Cité), A. BOZIC (JIS, Ljubljana)	10-28.7	Doctoral Training Program: Ab Initio Methods and Emerging Technologies for Nuclear Structure C. BARBIERI (University of Milan), A. ROGGERO (University of Trento)
<b>FEBRUARY</b> 20-24.2	LAVA Meeting C. BONANNO (INFN Firenze), M. P. LOMBARDO (INFN Firenze), M. PEARDON (Trinity College Dublin)	10-14.7	Tensor Spin Observables ** K. SLIFER (University of New Hampshire), D. HIGINBOTHAM (JLab), D. KELLER (University of Virginia), E. LONG (University of New Hampshire)
<b>MARCH</b> 13-17.3	Holographic Perspectives on Chiral Transport K. LANDSTEINER (IFT-UAM/CSIC Madrid), U. GURSOY (University of Utrecht), M. KAMINSKI (University of Alabama), D. KHARZEEV (Stony Brook)	17-21.7	Short-Distance Nuclear Structure and PDF * N. FOMIN (University of Tennessee), J. ARRINGTON (LBNL), W. COSYN (Florida International University), N. ROCCO (Fermi National Laboratory)
20-24.3	The Gradient Flow in QCD and other Strongly Coupled Field Theories C. MONAHAN (William & Mary), R. HARLANDER (University of Aachen), A. HASENFRATZ (University of Colorado, Boulder), O. WITZEL (Siegen University)	31.7-4.8	Quantum Sensing and Fundamental Physics with Levitated Mechanical Systems A. VINANTE (IFN-CNR), D. BUDKER (Johannes Gutenberg University Mainz), G. HETET (École Normale Supérieure Paris), H. ULBRICH (University of Southampton)
<b>MAY</b> 2-5.5	Quantum Science Generation   QSG D. DE BERNARDIS (INO-CNR BEC Center), V. AMITRANO (University of Trento - INO-CNR), A. BALDAZZI (University of Trento), A. BERTI (University of Trento - INO-CNR), I. CARUSOTTO (INO-CNR BEC Center), D. CONTESSI (University of Trento - INO-CNR), A. NARDIN (University of Trento - INO-CNR), L. PAVESI (University of Trento, Italy)	<b>AUGUST</b> 21-25.8	ECT*-APCTF Joint Workshop: Exploring Resonance Structure with Translucid CPDs * S. DIEHL (Justus Liebig University Giessen), V. BRAUN (University Regensburg), K. JOO (University of Connecticut), Y. OH (Kyungpook National University), C. VAN HULSE (University of Alcalá, Madrid), C. WEISS (Jlab)
15-19.5	Color Glass Condensate at the Electron-Ion Collider* D. TRIANTAFYLLOPOULOS (ECT*), N. ARMESTO (University of Santiago de Compostela), E. IANCU (University of Paris-Saclay, IPHT), T. LAPPI (University of Jyväskylä)	<b>SEPTEMBER</b> 4-8.9	Many-Body Quantum Physics with Machine Learning A. RIOS HUGUET (Institute of Cosmos Sciences, University of Barcelona), G. CARLEO (EPFL), E. INACK (PITP), A. LOVATO (ANL & TIFPA)
22-26.5	From First-Principles QCD to Experiments* M. HUBER (Giessen University), G. EICHMANN (LIP Lisbon), M. P. LOMBARDO (INFN Firenze), P. MARIS (Iowa State University), J. M. PALOWSKI (Heidelberg University)	11-15.9	MICRA2023: Microphysics in Computational Relativistic Astrophysics* E. O'CONNOR (Stockholm University), C. FROHLICH (Carolina State University), A. PEREGO (University of Trento)
29.5-1.6	2nd CMS Heavy Ion Workshop: Bringing Together the LHC Heavy Ion Community G. KRINTIRAS (The University of Kansas), Y.J. LEE (MIT), W. LI (Rice University), C. LOURENCO (CERN), A. STAHL (CERN)	18-22.9	Parton Distribution Functions at a Crossroad * M. DING (Helmholtz Zentrum Dresden Rosendorf), J. PAPAVALSILIOU (University of Valencia), C. QUINTANS (LIP, Lisbon), C. ROBERTS (Nanjing University)
<b>JUNE</b> 5-9.6	Nuclear and Particle Physics on a Quantum Computer: Where do we stand now? A. BAZAVON (Michigan State University), M. PAVONDI (University of Maryland), R. DE MICHIELLI (SI), T. JAROLE (University of Bonn), B. KUBIS (HISKP Bonn), D. LERSCH (FSU Tallahassee)	25-29.9	Strongly Interacting Matter in Extreme Magnetic Fields * S. VARESE (UNICAMP), A. AYALA (UNAM), D. BLASCHKE (University of Wrocław), G. ENDRODI (University of Bielefeld), R. FARIAS (Universidade Federal de Santa Maria)
12-16.6	Quantum Simulation of Gravitational Problems on Condensed Matter Analog Models I. CARUSOTTO (INO-CNR BEC Center), R. BALBINOT (University of Bologna), G. FERRARI (University of Trento), M. RINALDI (University of Trento)	<b>OCTOBER</b> 9-13.10	ROCKSTAR: Towards a Roadmap of the Crucial measurements of Key observables in Strangeness reactions for neutron $\alpha$ TARA equation of state ** A. SCORDO (INFN-INFN), D. BOSNAR (University of Zagreb), C. CURCEANU (INFN-INFN), A. RAMOS (Institut de Ciències del Cosmos, Barcelona), F. SAKUMA (RIKEN), O. VAZQUEZ-DOCE (INFN-INFN), I. VIDANA (INFN Catania)
26-30.6	Machins Learning for Lattice Field Theory and Beyond D. HACKETT (MIT), G. AARTIS (Swansea University & ECT*), D. BACHTIS (Swansea University), B. LUCINI (Swansea University), P. SHANAHAN (MIT)	23-27.10	Critical Stability of Few-Body Quantum Systems * A. KHEVTSKY (INFN Pisa), T. FREDERICO (Instituto Tecnológico de Aeronáutica), O. FYNBO (Aarhus University), J.M. RICHARD (Institut de Physique des 2 Infinis de Lyon)
<b>JULY</b> 3-7.7	COLMO: Quantum Collapse Models Investigated with Particle, Nuclear, Atomic and Macro systems C. CURCEANU (INFN-INFN), A. BASSI (University and INFN Trieste), M. DERAKHSHANI (Rutgers University), L. DIOSI (University Budapest), S. DONADI (INFN Trieste), K. PISCICCHIA (CREP)	<b>NOVEMBER</b> 20-24.11	ALPACA: modern Algorithms in machine learning and data analysis: from medical Physics to research with Accelerators and in underground laboratories ** F. NAPOLITANO (INFN Frascati), R. DEL GRANDE (TU Munich), F. GROSA (CERN), M. SKURZOK (Jagiellonian University Krakow)

\*STRONG-2020 supported workshop \*\*EUROLABS supported workshop

The ECT\* is part of the Bruno Kessler Foundation. The Centre is funded by the Autonomous Province of Trento, funding agencies of EU Member and Associated States, INFN-TIFPA, and has the support of the Department of Physics of the University of Trento. The Director of the ECT\* is Prof. Gert Aarts (Trento and Swansea University)  
For information: [staff@ectstar.eu](mailto:staff@ectstar.eu) | [www.ectstar.eu](http://www.ectstar.eu)



# 2023 Activities

full programme:  
24 workshops and collaboration meetings  
from January to November  
selected by Scientific Board

## DTP: Ab Initio Methods and Emerging Technologies for Nuclear Structure

# From hybrid to the new(er) normal

- aim is to stimulate scientific discussion in an informal and creative environment
- hence, ECT\* returns to in-person participation as the preferred way of interaction
- to build on the past experiences and to support an inclusive global research community, presentations at workshops may be broadcast via zoom





# EURO-LABS supported meetings

- EXOTICO: EXOTIc atoms meet nuclear COLLisions for a new frontier precision era in low-energy strangeness nuclear physics 17-21/10 2022, **49** participants, **8** participants supported
- Tensor Spin Observables 10-14/7 2023, **24** participants, **4** participants supported
- DTP Ab Initio Methods and Emerging Technologies for Nuclear Structure 10-28/7 2023, **33** participants, **15** participants supported

# DTP 2023: Ab Initio Methods and Emerging Technologies for Nuclear Structure

- 10-28 July 2023

organizers:

- Carlo Barbieri  
(University of Milan, INFN)
- Alessandro Roggero  
(University of Trento)
- 33 participants



supported by

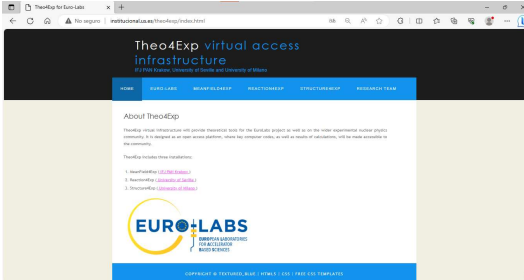


# EURO-LABS supported meetings: comments

- ECT\* Scientific Board is selection panel
- only 3 meetings supported so far: preserve financial support until after STRONG-2020 ends (May 2024)
- 2024: 31 workshop proposals submitted, 22 selected, EURO-LABS assignment to follow

# Theo4Exp

[institucional.us.es/theo4exp](http://institucional.us.es/theo4exp)



The screenshot shows a web browser displaying the Theo4Exp website. The page title is "Theo4Exp virtual access infrastructure". Below the title, it says "EURO-LABS" and "UNIVERSITY OF GRANADA AND UNIVERSITY OF MADRID". The navigation menu includes "HOME", "EURO-LABS", "ABOUT THEO4EXP", "REACTOR4EXP", "STRUCTURE4EXP", and "RESEARCH TEAM". The main content area is titled "About Theo4Exp" and contains the following text: "Theo4Exp virtual infrastructure will provide theoretical tools for the EuroLab project as well as on the wider experimental nuclear physics community. It is designed as an open access platform, where key computer codes, as well as results of calculations, will be made accessible to the community." Below this text, it states "Theo4Exp includes three installations:" followed by a numbered list: "1. Accelerator ([Link to the system](#))", "2. Reactor4Exp ([Link to the system](#))", and "3. Structure4Exp ([Link to the system](#))". At the bottom of the page, there is a logo for "EURO-LABS" and the text "EUROPEAN LABORATORIES FOR ACCELERATOR BASED SCIENCES". The footer contains the text "COPYRIGHT © TEXTURED, BLISS / HTML © 2011 / THEO4EXP TEMPLATE".

# Theo4Exp Virtual Access Infrastructure

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

- Theo4Exp virtual access infrastructure will provide theoretical tools for the EURO-LABS project as well as for the wider experimental nuclear physics comm.
- 3 installations:
  - ⇒ MeanField4Exp (Kraków): will provide access to mean-field theory service in the domain of **nuclear structure** physics. Deputies: Jerzy Dudek (IHPC/CNRS & Strasbourg U.)/ Piotr Bednarczyk
  - ⇒ Reaction4Exp (Sevilla): will provide codes used for **nuclear reaction** calculations. Coordinator: Manuela Rodríguez-Gallardo
  - ⇒ Structure4Exp (Milano): will provide virtual access to other codes that use advanced tools of **nuclear structure** theory. Deputy: Gianluca Coló

# Contracted personnel

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- **MeanField4Exp**(IFJ PAN, Kraków): 2-year contract
  - ⇒ Dr. Abdelghafar Gaamouchi, from 02/2023.
- **Reaction4Exp** (U. Sevilla): 2-year contract
  - ⇒ Carla Muñoz (Master), from 09/2023.
- **Structure4Exp** (U. Milano): 1-year contract
  - ⇒ Dr. Imane Moumene, from 03/2023.



# International Review Panel (IRP)

WP2.4: Research infrastructures offering theoretical support for experiments

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- **Members:** Piotr Bednarczyk (Chairperson, IFJ PAN), Krzysztof Rusek (UW), Antonio M. Moro (USE), Ian J. Thompson (LLNL), Enrico Vigezzi (UMIL) and Angela Gargano (INFN)
- **Second meeting** took place (online) on **20th September 2023** with participation of the Coordinating Team (Manuela Rodríguez-Gallardo, Jerzy Dudek and Gianluca Colò) and the Coordinator of WP2 (Adam Maj).
- Two points to **highlight**:
  - ⇒ The **user authentication** has to be the simplest as possible and open to all potential users.
  - ⇒ It is necessary to make **institutions** see that this new infrastructure is beneficial to them and therefore it will be necessary to invest on it for maintenance.

# User authentication

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Manuela Rodríguez Gallardo

- We have an application developed by WP5.2 EURO-LABS to control the user access:  
<https://iam-eurolabs.ijclab.in2p3.fr>



INDIGO - DataCloud

Welcome to **indigo-dc**

Sign in with

eduGAIN 



# User authentication

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<https://iam-eurolabs.ijclab.in2p3.fr>



INDIGO - DataCloud

Welcome to **indigo-dc**

Sign in with

eduGAIN 

- We hope soon to include other identity networks like ORCID

# At the moment, access via institutional identity provider (IdP)

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INDIGO - DataCloud

Welcome to **indigo-dc**

Sign in with



INDIGO - DataCloud

Sign in with your IdP

You will be redirected for authentication to:

**Universidad de Sevilla**

Proceed?



INDIGO - DataCloud

Sign in with your IdP

**Universidad de Sevilla**

[Back to login page](#)



**Identificación de usuario**

**Autenticación centralizada**

¿Problemas con su contraseña de acceso o sobre factor de autenticación (2FA)? Siga las instrucciones indicadas en <https://iso.us.es>.

Introduzca su UVUSU o su correo

**USUARIO**

**CONTRASEÑA**

**ACEPTAR**

Quiero recuperar mi contraseña

**OTROS MEDIOS DE AUTENTICACIÓN**

**Confirmar**

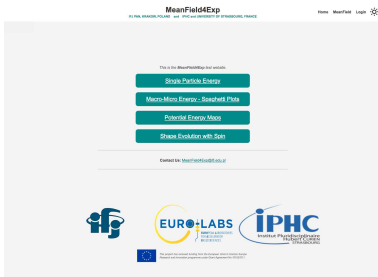
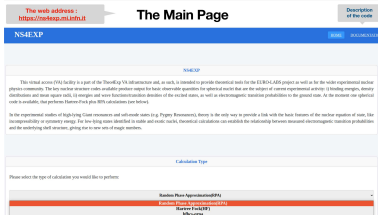
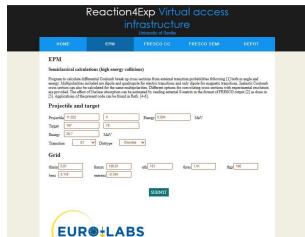
Una vez que se haya autenticado no será necesario identificarse de nuevo para acceder a otros recursos.  
Para desconectarse, recomendamos que cierre su navegador (cerrando todas las ventanitas).

ISIS ISO v1.9.3 - 4/30/21

# What about the platform for each installation?

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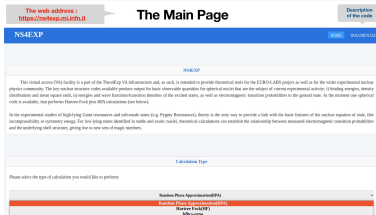
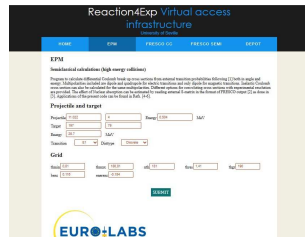
Manuela Rodríguez Gallardo

# What about the platform for each installation?

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Let's see!

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

## MeanField4Exp

IFJ PAN, KRAKÓW, POLAND and IPHC and UNIVERSITY OF STRASBOURG, FRANCE

[Home](#) [MeanField](#) [Login](#) 

This is the MeanField4Exp test website.

[Single Particle Energy](#)

[Macro-Micro Energy - Spaghetti Plots](#)

[Potential Energy Maps](#)

[Shape Evolution with Spin](#)

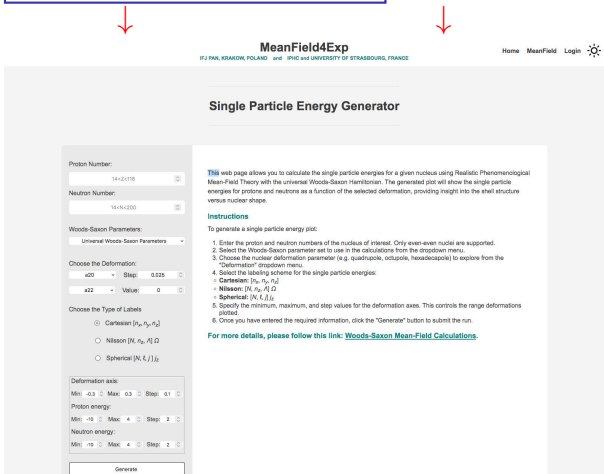
Contact Us: [MeanField4Exp@ifj.edu.pl](mailto:MeanField4Exp@ifj.edu.pl)



This project has received funding from the European Union's Horizon Europe Research and Innovation programme under Grant Agreement No 101027511.

List of propositions for the user choice

Guidelines



**MeanField4Exp**  
IFJ PAN, KRAKOW, POLAND and IPHC and UNIVERSITY OF STRASBOURG, FRANCE

Home MeanField Login

## Single Particle Energy Generator

**Proton Number:**  
14+Z+118

**Neutron Number:**  
16+N+200

**Woods-Saxon Parameters:**  
Universal Woods-Saxon Parameters

**Choose the Deformation:**  
a20 Step: 0.025  
a22 Value: 0

**Choose the Type of Labels:**  
 Cartesian  $(n_x, n_y, n_z)$   
 Nilsson  $[N, n_x, A] \Delta$   
 Spherical  $[N, l, j] l_x$

**Deformation axis:**  
Min: -0.3 Max: 0.3 Step: 0.1  
**Proton energy:**  
Min: -18 Max: 4 Step: 2  
**Neutron energy:**  
Min: -18 Max: 4 Step: 2

Generate

**This** web page allows you to calculate the single particle energies for a given nucleus using Realistic Phenomenological Mean-Field Theory with the universal Woods-Saxon Hamiltonian. The generated plot will show the single particle energies for protons and neutrons as a function of the selected deformation, providing insight into the shell structure versus nuclear shape.

**Instructions**  
To generate a single particle energy plot:

1. Enter the proton and neutron numbers of the nucleus of interest. Only even-even nuclei are supported.
2. Select the Woods-Saxon parameter set to use in the calculations from the dropdown menu.
3. Choose the nuclear deformation parameter (e.g. quadrupole, octupole, hexadecapole) to explore from the "Deformation" dropdown menu.
4. Select the labeling scheme for the single particle energies:
  - Cartesian:  $(n_x, n_y, n_z)$
  - Nilsson:  $[N, n_x, A] \Delta$
  - Spherical:  $[N, l, j] l_x$
5. Specify the minimum, maximum, and step values for the deformation axes. This controls the range deformations plotted.
6. Once you have entered the required information, click the "Generate" button to submit the run.

For more details, please follow this link: [Woods-Saxon Mean-Field Calculations](#).

# Meanfield4Exp: Single particle drawings

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Manuela Rodríguez Gallardo

Proton Number:

Neutron Number:

Woods-Saxon Parameters:  
 Universal Woods-Saxon Parameters

Choose the Deformation:  
 a32 - Step: 0.025  
 a22 - Value: 0

Choose the Type of Labels  
 Cartesian [ $n_x, n_y, n_z$ ]  
 Nilsson [ $N, n_x, l$ ]  $\Omega$   
 Spherical [ $N, l, j$ ] $l_z$

Deformation axis:  
 Min: -0.3 Max: 0.3 Step: 0.1

Proton energy:  
 Min: -10 Max: 4 Step: 2

Neutron energy:  
 Min: -18 Max: -2 Step: 2

This web page allows you to calculate the single particle energies for a given nucleus using Realistic Phenomenological Mean-Field Theory with the universal Woods-Saxon Hamiltonian. The generated plot will show the single particle energies for protons and neutrons as a function of the selected deformation, providing insight into the shell structure versus nuclear shape.

### Instructions

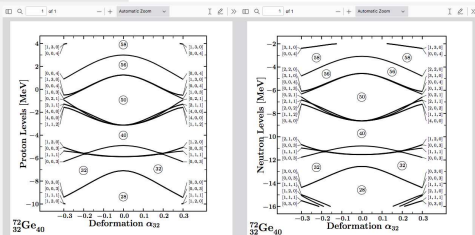
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4. Select the labeling scheme for the single particle energies:
  - Cartesian: [ $n_x, n_y, n_z$ ]
  - Nilsson: [ $N, n_x, l$ ]  $\Omega$
  - Spherical: [ $N, l, j$ ] $l_z$
5. Specify the minimum, maximum, and step values for the deformation axes. This controls the range deformations plotted.
6. Once you have entered the required information, click the "Generate" button to submit the run.

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
Example

$Z = 32, N = 40$ , oct. def.



## MeanField4Exp

IFJ PAN, KRAKOW, POLAND and IPHC and UNIVERSITY OF STRASBOURG, FRANCE

Home MeanField Login 

---

### Macro-Micro Energy - Spaghetti Plots Generator

Proton Number:

Neutron Number:

$\Delta Z$ :   $\Delta N$ :

Woods-Saxon Parameters:

Choose the Type of Energy:  
Total energy = E(FYU) + Shell(E) + Correlation

Choose the Deformation:  
 $a_{12}$   Step:   
 $a_{20}$   Value:

Deformation axis:  
Min: -3  Max: 0.3  Step: 0.1   
Isotopes energy:  
Min: -4  Max: 12  Step: 2   
Isotone energy:  
Min: -4  Max: 12  Step: 2

This form allows you to generate spaghetti plots of the macro-micro energy for a chain of isotopes and isotones using

**Realistic Phenomenological Nuclear Mean-Field Theory Calculations**  
with the  
Universal Woods-Saxon Hamiltonian.

**Instructions for Formatting the Plot**

- 1. Specification of the Nucleus and Parametrisation**  
Choose a central nucleus by entering the proton and neutron numbers.  
Select the range of nuclei by adjusting  $\Delta Z$  and  $\Delta N$ .  
"Woods-Saxon Parameters" allows you to select a parameter set for the Woods-Saxon Hamiltonian from the dropdown menu.
- 2. Energy Specifications**  
Choose the variant of the total energy formula according to which  
$$E_{tot} = E_{Macro} + E_{Shell} + E_{pairing}$$
The macroscopic energy,  $E_{Macro}$ , appears in two variants, the so-called Yukawa-folded [E(FYU)] and the Lubin-Strasbourg Drop [E(LSD)] liquid drop models. The standard Strutinsky shell-correction energy is denoted  $E_{shell}$ , whereas the pairing term,  $E_{pair}$ , can be chosen in the form of the so-called pairing correlation energy with a simple (BCS) approximation or a more advanced Particle Number Projection (PNP) approximation.  
Those interested can find the original references clicking "[Reference List](#)".
- 3. Description of Nuclear Deformation**  
"Choose the Deformation" Allows you to select the nuclear deformation from the dropdown menu.  
Deformation axis, isotope energy, and isotone energy allow you to adjust the x and y axes.
- 4. Execution**  
Once all input data is specified, click "Generate" to submit the run.  
For more details, please follow the link: [Woods-Saxon Mean-Field Calculations](#).



# Meanfield4Exp: Total energy drawings

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

Proton Number:

Neutron Number:

$\Delta Z$ :   $\Delta N$ :

Woods-Saxon Parameters:

Choose the Type of Energy:   
 Total energy =  $E_{PYU} + \text{Shell}(x) + \text{Correlation}$

Choose the Deformation:   
 $a32$  Step:    
 $a20$  Value:

Deformation axis:   
 Min:  Max:  Step:

Isotopes energy:   
 Min:  Max:  Step:

Isotones energy:   
 Min:  Max:  Step:

This form allows you to generate spaghetti plots of the macro-micro energy for a chain of isotopes and isotones using

Realistic Phenomenological Nuclear Mean-Field Theory Calculations  
with the  
Universal Woods-Saxon Hamiltonian.

## Instructions for Formatting the Plot

### 1. Specification of the Nucleus and Parametrisation

Choose a central nucleus by entering the proton and neutron numbers.

Select the range of nuclei by adjusting  $\Delta Z$  and  $\Delta N$ .

"Woods-Saxon Parameters" allows you to select a parameter set for the Woods-Saxon Hamiltonian from the dropdown menu.

### 2. Energy Specifications

Choose the variant of the total energy formula according to which

$$E_{\text{tot}} = E_{\text{macro}} + E_{\text{pair}} + E_{\text{spin}}$$

The macroscopic energy,  $E_{\text{macro}}$ , appears in two variants, the so-called "Yukawa-folded" [E]PYU [ ] and the Lubin-Strasbourg Drop [E]LSD [ ] liquid drop models. The standard Strutinsky shell-correction energy is denoted  $E_{\text{shell}}$ , whereas the pairing term,  $E_{\text{pair}}$ , can be chosen in the form of the so-called pairing correlation energy with a simple (BCS) approximation or a more advanced Pairing Number Projection (PNP) approximation.

Those interested can find the original references clicking "[Reference List](#)".

### 3. Description of Nuclear Deformation

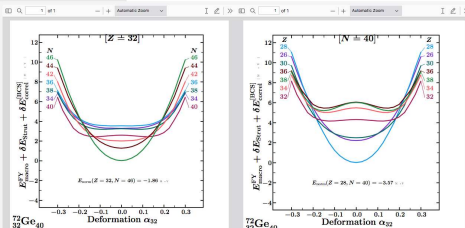
"Choose the Deformation" Allows you to select the nuclear deformation from the dropdown menu.

Deformation axis, isotope energy, and isotone energy allow you to adjust the x and y axes.

### 4. Execution

Once all input data is specified, click "Generate" to submit the run.

For more details, please follow the link: [Woods-Saxon Mean-Field Calculations](#).



## Nuclear Potential-Energy Maps Generator

Proton Number:

Neutron Number:

Choose the Type of Energy:  
Total energy = E(TFU) + ShellE + Correlation

Step:  Smoothing:

Deformation:  
List of 3D/4D Deformation Spaces:

X axis:  Y axis:  Min1:  Min2:

Choose a range of deformation:

X-axis: Min:  Max:  Step:

Y-axis: Min:  Max:  Step:

Table of the Energy Minima

Welcome to the Nuclear Potential-Energy Maps Generator allowing you to explore the elementary shape properties of atomic nuclei using

**Realistic Phenomenological Nuclear Mean-Field Theory Calculations**  
with the  
**Universal Woods-Saxon Hamiltonian.**

To design the appearance of the potential-energy map you are interested in, please employ the formatting table located on the left side of the page. This formatting table will allow you to choose the parameters needed by the algorithms to generate a map of the nuclear potential energy surface for the atomic nucleus which you define to start with.

### Instructions for formatting the drawing

- 1. Specification of the nucleus**  
Select the proton and neutron numbers of the nucleus needed. For information: our data base contains the results pre-calculated only for even-even nuclei.
- 2. Energy specifications**  
Choose the variant of the total energy formula according to which
 
$$E_{\text{tot}} = E_{\text{macro}} + E_{\text{shell}} + E_{\text{pairing}}$$

The macroscopic energy,  $E_{\text{macro}}$ , appears in three variants, the so-called Yukawa-folded [E(FYU)], Lublin-Strasbourg Drop [E(LSD)], and "traditional" Myers-Swiatecki [E(M-S)] liquid drop models. The standard Strutinsky shell-correction energy is denoted  $E_{\text{shell}}$ , whereas the pairing term,  $E_{\text{pair}}$ , can be chosen in the form of the so-called pairing correlation energy with a simple (RCG) approximation or a more advanced Particle Number Projection (PNP) approximation.

Those interested can find the original references clicking "[Reference List](#)".
- 3. Map appearance and format specifications**  
The size of the colour steps [in MeV] for the energy contour plotting can be adjusted by using the input box Step. The user can influence the aesthetic aspects such as the smoothness of the contour lines via the input field Smoothing.
- 4. Description of nuclear deformation**  
The user can select among the potential energy surfaces calculated using multiprocessor computer systems within an ensemble of 3D and 4D (3-dimensional and 4-dimensional original deformation space). The graphical representation corresponds to choosing two deformation variables as x- and y-coordinates. The total energies are minimised over the remaining variables at the disposal. In particular, in the case of a 3D space of deformations d1, d2 and d3 the user can select projections (d1,d2), alternatively (d1,d3) or (d2,d3). In the case of the 4D spaces the user will have 6 independent choices at her/his disposal.

Our system is being intensively developed — but at present we offer only one illustrative variant, probably the most often found in the literature, namely  $a_{32}$ ,  $a_{22}$  and  $a_{42}$  (also known under the notation (beta, gamma) and beta<sub>42</sub>).

By default, our algorithm displays the whole range of the pre-calculated deformation space. However, the user can customise the deformation ranges by selecting the option "Choose the range of deformation". By clicking the user displays self-explanatory input fields for the minimum, maximum, and step values for the x- and y-axes.
- 5. Execution**  
Once all the input data are specified, click the "Generate" button to submit the run.

For more details, please follow the link: [Woods-Saxon Mean-Field Calculations](#).

# Meanfield4Exp: Total energy map drawings

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

Potential energy contour

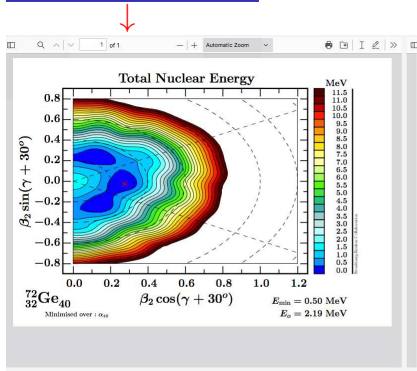


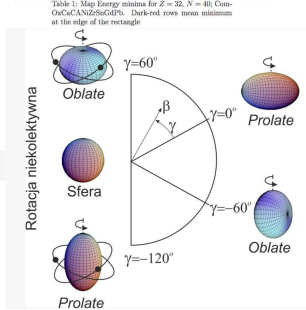
Table of minima

Strasbourg-Krakow Collaboration

Nucleus  $^{72}\text{Ge}_{40}$

- Total Energy:  $E(\text{FYU}) + \text{Shell}[e] + \text{Correlation}[BCS]$
- Minimum Energy:  $E_{\min} = 0.50 \text{ MeV}$
- Spherical-point Energy:  $E_s = 2.19 \text{ MeV}$

Energy [MeV]	$\beta_2 \cos(\gamma + 30^\circ)$	$\beta_2 \sin(\gamma + 30^\circ)$	$\alpha_{40}$	$\beta$	$\gamma(^{\circ})$
0.0000	0.275	-0.025	-0.025	0.276	-35
0.0015	0.125	0.250	-0.025	0.280	33
0.0718	0.150	-0.225	0.050	0.270	-86



## At high spin, microscopic effects are negligible

### MeanField4Exp

IFJ PAN, KRAKOW, POLAND and IPHC and UNIVERSITY OF STRASBOURG, FRANCE

Home MeanField Login

## Shape Evolution with Spin

In this page you can explore the Shape Evolution with Spin using the table located on the left side of the page. This table will allow you to choose the parameters needed by the algorithms to generate a map of the nuclear potential energy surface at a specific Spin of rotation for the atomic nucleus which you define to start with.

### Instructions for formatting the drawing

- 1. Specification of the nucleus**  
Select the proton and neutron numbers of the nucleus needed. For information: our data base contains the results pre-calculated only for even-even nuclei.
- 2. Energy specifications**  
Choose the variant of the total energy formula according to which
 
$$E_{tot} = E_{Macro} + E_{rotational}$$

The macroscopic energy,  $E_{Macro}$ , appears in two variants, the so-called Yukawa-folded [E(FYU)], Lublin-Strasbourg Drop [E(LSD)]. The second part ...

Those interested can find the original references clicking "[Reference List](#)".
- 3. Map appearance and format specifications**  
The size of the colour stripes [in MeV] for the energy contour plotting can be adjusted by using the input box Step. The user can influence the aesthetic aspects such as the smoothness of the contour lines via the input field Smoothing.
- 4. Description of nuclear deformation**  
The user can select among the potential energy surfaces calculated using multiprocessor computer systems within an ensemble of 3D and 4D (3-dimensional and 4-dimensional original deformation spaces). The graphical representation corresponds to choosing two deformation variables as x, and y-coordinates. The total energies are minimised over the remaining variables at the disposal. In particular, in the case of a 3D space of deformations d1, d2 and d3 the user can select projections (d1,d2), alternatively (d1,d3) or (d2,d3). In the case of the 4D spaces the user will have 6 independent choices at her/his disposal.

Our system is being intensively developed — but at present we offer only one illustrative variant, probably the most often found in the literature, namely  $\alpha_{20}$ ,  $\alpha_{22}$  and  $\alpha_{40}$  (also known under the notation (beta, gamma) and beta<sub>4</sub>).

By default, our algorithm displays the whole range of the pre-calculated deformation spaces. However, the user can customise the deformation ranges by selecting the option "Choose the range of deformation". By clicking the user displays self-explanatory input fields for the minimum, maximum, and step values for the x- and y-axis.

n Number:

on Number:

se the Type of Energy:  
if energy = E(FYU) + Rotational Energy ▾

0.5 ▾ Smoothing: 2 ▾

62 ▾ WhichSpin: ly ▾

mation:

of 3D/4D Deformation Spaces:  
bet, gam, a40 ▾

axis	Y axis	Min1	Min2
x	gam	a40	

Choose a range of deformation

es:  
0 ▾ Max: 1.2 ▾ Step: 0.2 ▾

es:  
-0.8 ▾ Max: 0.8 ▾ Step: 0.2 ▾

Generate

017. hiliar

Printed

- All My
- Android
- App
- Desktop
- Down
- Tor
- Print

Devices

Shared

- Beta
- Sigma

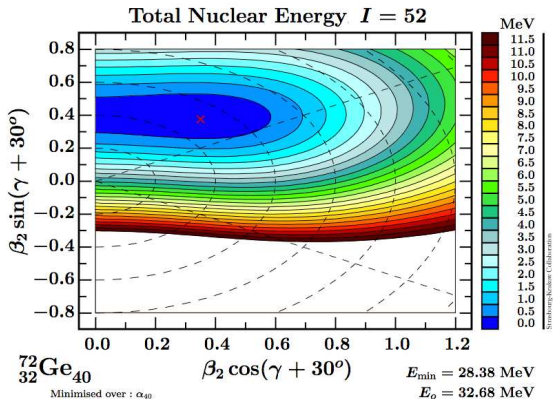
# Meanfield4Exp: Drawing at spin= $56\hbar$

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

Preliminary

Energy minimized with respect  $\beta_4$



# Meanfield4Exp: New options being considered/elaborated

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

01. *Shape coexistence and evolution with spin and temperature;*
02. *Pairing and its evolution with spin and temperature;*
03. *Axial symmetry and related isomers: K-isomers, yrast traps;*
04. *Many-particle many hole excited configurations – degeneracies;*
05. *Electromagnetic transitions;*
06. *Cranking, 3D cranking and chirality;*
07. *Nuclear point group symmetries;*
08. *Band crossings, Band termination;*
09. *Jacobi and Poincare shape transitions;*
10. *Giant-Dipole Resonances – spin & temperature dependence;*
11. *Configuration controlled shape evolution;*
12. *Mass parameters, mass tensor, collective motion;*

## Reaction4Exp Virtual access infrastructure

University of Sevilla

HOME
EPM
FRESKO CC
FRESKO SEMI
DEPOT

### EPM

**Semiclassical calculations (high energy collisions)**

Program to calculate differential Coulomb break up cross sections from external transition probabilities following [1] both in angle and energy. Multipolarities included are dipole and quadrupole for electric transitions and only dipole for magnetic transitions. Inelastic Coulomb cross section can also be calculated for the same multipolarities. Different options for convoluting cross sections with experimental resolution are provided. The effect of Nuclear absorption can be estimated by reading external S-matrix in the format of FRESKO output [2] as done in [3]. Applications of the present code can be found in Refs. [4-6].

**Projectile and target**

Projectile:   Energy:  MeV

Target:


Energy:  MeV

Transition:  Distype:

**Grid**

thimin:  thimax:  ntb:  thres:  thp:

benc:  enerenc:

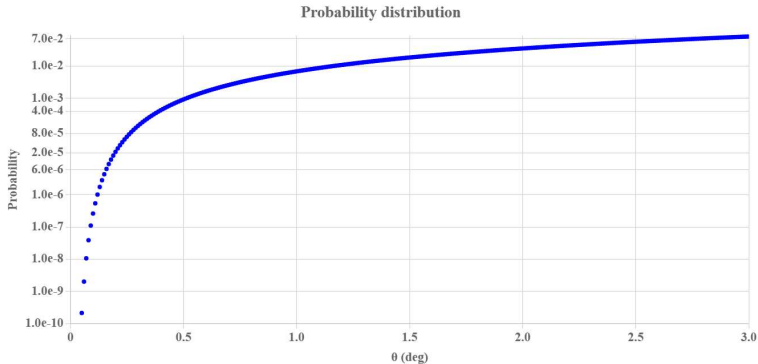


# Reaction4Exp: Breakup probability

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Manuela Rodríguez Gallardo

$^{11}\text{Li} + ^{208}\text{Pb}$  at 759 MeV





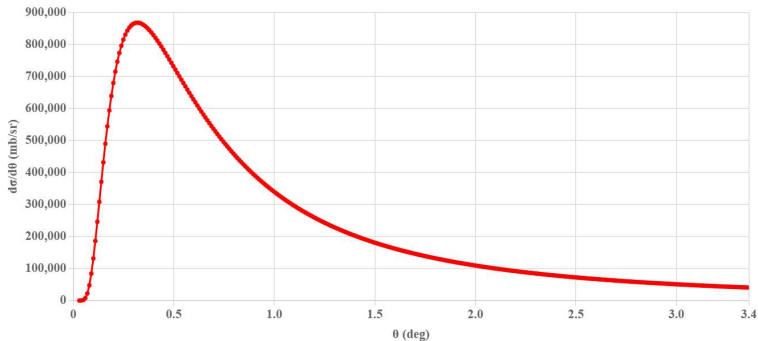
# Reaction4Exp: BU angular distribution

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

$^{11}\text{Li} + ^{208}\text{Pb}$  at 759 MeV

Angular distribution of differential cross section



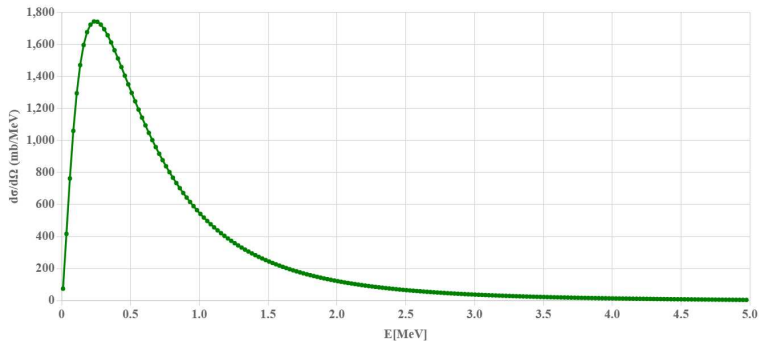
# Reaction4Exp: BU energy distribution

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

$^{11}\text{Li} + ^{208}\text{Pb}$  at 759 MeV

Energy distribution of cross section



- 1 Optical Model calculations: **FRESCO**  
(<http://www.fresco.org.uk/>)
- 2 Coupled-Channels calculations: **FRESCO**
- 3 Semiclassical calculations (high energy collisions):  
**EPM\_SEV**  
*Example: V. Pesudo et al., Phys. Rev. Lett. 118 (2017) 152502*
- 4 Double folding potentials from density distributions:
  - ⇒ **DFPOT**  
*J. Cook, Comp. Phys. Comm. 25(2), 125-139 (1982)*
  - ⇒ **SPP**  
*L.C. Chamon, B.V. Carlson and L.R. Gasques, Comp. Phys. Comm. 267 (2021) 108061*

The web address :  
<https://ns4exp.mi.infn.it>

## The Main Page

Description of the code

NS4EXP

[HOME](#)
[DOCUMENTATION](#)

**NS4EXP**

This virtual access (VA) facility is a part of the Theo4Exp VA infrastructure and, as such, is intended to provide theoretical tools for the EURO-LABS project as well as for the wider experimental nuclear physics community. The key nuclear structure codes available produce output for basic observable quantities for spherical nuclei that are the subject of current experimental activity: i) binding energies, density distributions and mean square radii, ii) energies and wave functions/transition densities of the excited states, as well as electromagnetic transition probabilities to the ground state. At the moment one spherical code is available, that performs Hartree-Fock plus RPA calculations (see below).

In the experimental studies of high-lying Giant resonances and soft-mode states (e.g. Pygmy Resonances), theory is the only way to provide a link with the basic features of the nuclear equation of state, like incompressibility or symmetry energy. For low-lying states identified in stable and exotic nuclei, theoretical calculations can establish the relationship between measured electromagnetic transition probabilities and the underlying shell structure, giving rise to new sets of magic numbers.

**Calculation Type**

Please select the type of calculation you would like to perform:

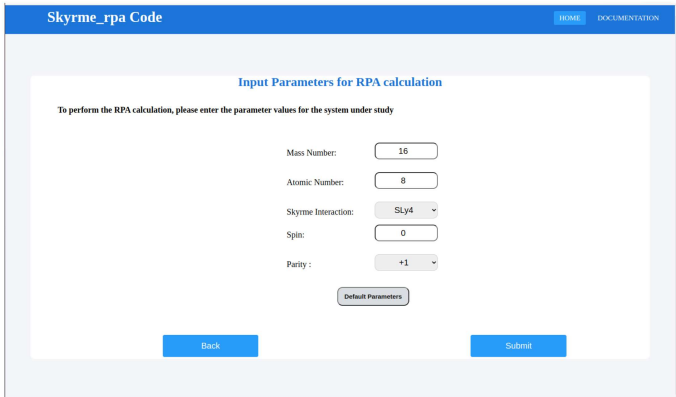
Random Phase Approximation(RPA) ▾

Random Phase Approximation(RPA)

Hartree Fock(HF)

hfcs-qrpa

## Input Form Page



The screenshot shows a web interface for the Skyrme\_rpa Code. At the top, there is a blue header with the text "Skyrme\_rpa Code" on the left and "HOME" and "DOCUMENTATION" on the right. Below the header, the main content area is titled "Input Parameters for RPA calculation" in blue. A sub-header reads: "To perform the RPA calculation, please enter the parameter values for the system under study". The form contains several input fields: "Mass Number" with a text box containing "16", "Atomic Number" with a text box containing "8", "Skyrme Interaction" with a dropdown menu showing "SLy4", "Spin" with a text box containing "0", and "Parity" with a dropdown menu showing "+1". Below these fields is a button labeled "Default Parameters". At the bottom of the form area, there are two blue buttons: "Back" on the left and "Submit" on the right.

## Results Page

Skyrme\_rpa Code
HOME DOCUMENTATION

### Results

Calculations completed successfully!

[View input file](#)

File name	File size	Download	Display	Plot
Plot_Bel_EM.dat	15.82 KB	<a href="#">Download</a>	<a href="#">Display</a>	<a href="#">Plot</a>
Plot_Bel_IS.dat	15.82 KB	<a href="#">Download</a>	<a href="#">Display</a>	<a href="#">Plot</a>
Plot_Bel_IV.dat	15.82 KB	<a href="#">Download</a>	<a href="#">Display</a>	<a href="#">Plot</a>
density.out	6.31 KB	<a href="#">Download</a>	<a href="#">Display</a>	<a href="#">Plot</a>
skyrme_rpa.out	10.74 KB	<a href="#">Download</a>	<a href="#">Display</a>	

[Generate combined Plots](#)

Email with results sent successfully to the same address you signed up with.

Example:

$^{16}\text{O}(0^+)$

EM, IS and IV strength functions

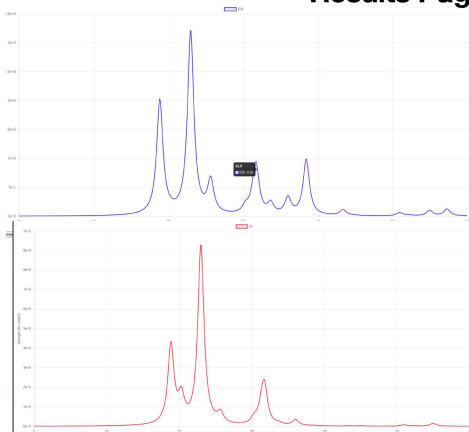
The main output file

Send results via email

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

## Results Page



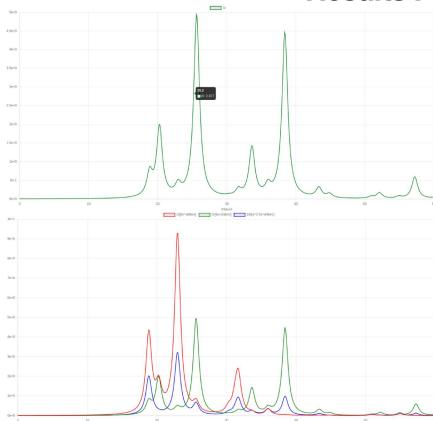
EM strength function

IS strength function

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## Results Page



IV strength function

Combined chart (EM, IS, IV)



- ① Binding energies, density distributions and mean square radii ([skyrme\\_rpa](#), [hfbcqs\\_qrpa](#))
  - ② Energies and wave functions/transition densities of the excited states, as well as electromagnetic transition probabilities to the ground state ([skyrme\\_rpa](#), [hfbcqs\\_qrpa](#))
  - ③ Calculations of charge-changing transitions
  - ④ Beta-decay half-lives
- [skyrme\\_rpa](#)  
*G. Colò et al., Comp. Phys. Comm. 184, 142 (2013)*
- [hfbcqs\\_qrpa](#)  
*G. Colò and X. Roca-Maza, arXiv:2102.06562 [nucl-th]*

# MileStones (MS)

WP2.4: Research  
infrastructures  
offering  
theoretical  
support for  
experiments

Manuela  
Rodríguez  
Gallardo

- **MS10 (Month 18)**: Contracted personnel for Theo4Exp VA in place and first codes available for users in the virtual facility
- **MS11 (Month 42)**: All software released and validated by IRP

So we are on time!

# Final remarks

WP2.4: Research infrastructures offering theoretical support for experiments

Manuela Rodríguez Gallardo

- By the **end of the year**, the calculation platforms will be **available** with several programs.
- There will be a common portal (institucional.us.es/theo4exp), including the link to the different installations.
- We will start then the users counting (Units of Access).
- The user registration is through (**Please, try!**) <https://iam-eurolabs.ijclab.in2p3.fr>.
- Try to favour the use of the installations when available and, in particular, for **summer schools**, **doctoral training**.
- Provide feedbacks!

# Time for discussion in Kraków

WP2.4: Research  
infrastructures  
offering  
theoretical  
support for  
experiments

Manuela  
Rodríguez  
Gallardo

