

Nuclear Structure and Decay Data (NSDD)
Network

Brookhaven National Laboratory
November 6-7, 2003

Decay Data Evaluation Project (DDEP)

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1. Objectives

- Applied Research and Detector Calibration
 - Reactor Materials
 - Nuclear Medicine
 - Astrophysics
- Data Content and Presentation
 - Atomic and Nuclear Properties
 - Measured Values and Evaluation Procedures
- Publication
 - Laboratory reports (Saclay, PTB)
 - WWW (<http://www.bnm.fr/bnm-Inhb/DDEP.htm>)
 - Bureau International des Poids et Mesures (BIPM)

2. Members of the DDEP Collaboration

- CEA/LNHB (**France**): Dr. Marie-Martine Bé, Dr. Vanessa Chiste, Dr. B. Duchemin
- PTB (**Germany**): Dr. Herbert Janssen, Dr. E. Schönfeld
- LBNL (**USA**): Dr. Edgardo Browne, Dr. Coral Baglin, Dr. Shiu-Chin Wu
- INEEL (**USA**): Dr. Richard G. Helmer, Dr. Charles Reich
- BNL (USA): Dr. Jagdish K. Tuli
- ANL (**USA**): Dr. Filip Kondev
- KRI (**Russia**): Dr. Valery P. Chechev
- NPL (**UK**): Dr. T. Desmond MacMahon
- IAEA (**Vienna**): Dr. Alan L. Nichols
- USP (**Brazil**): Dr. Vito Vanin, Dr. Rui M. Castro

Other collaborators

- CIEMAT (Spain): Dr. Jose M. Los Arcos
- UNED (Spain): Dr. Amalia Williard Torres)
- PTB (Germany): Dr. Rainer Dersch

3. Formation of the DDEP Collaboration

- 1991** Dick Helmer initiated discussions with members of PTB and Saclay
- 1991 – 1993** PTB and Saclay had reached a collaborative agreement
- 1994** Evaluators from the US received approval for participating in this effort. Soon Russian and English evaluators joined them.
- 1995** First meeting of the DDEP collaboration in Paris

4. Nuclear and Atomic Properties in DDEP Evaluations

Most properties evaluated by the DDEP collaboration are relevant to applied research, such as

- Radionuclide half-life
- Nuclear radiations (α , β^\pm , γ) energies, emission and transition probabilities, and equilibrium emission probabilities.
- Energies and emission probabilities of atomic radiations such as X rays, conversion and Auger electrons, and electron-positron pairs emitted in nuclear disintegrations.

Properties reported, but usually not evaluated, are:

- Q-values (from Audi et al. [3])
- Theoretical internal conversion coefficients (Rösel et al. [4]) and pair creation (Schlüter et al. [14]) coefficients.
- γ -ray multipolarities and mixing ratios (from ENSDF)
- Level schemes: levels spin, parity, and half-life (from ENSDF)
- Auger electron energies (from Schönfeld and Rodloff [5]) and X-ray energies and relative emission probabilities (from Schönfeld and Rodloff [6])
- Atomic data such as K-fluorescence yields (ω_K), mean L-shell fluorescence yields (ω_L), and total number of L vacancies created by a $K \rightarrow L$ transfer (η_{KL}) (from Schönfeld and Janssen [7]).
- Log ft's for β^\pm decay and α -decay hindrance factors (from ENSDF, or calculated with computer programs used for ENSDF evaluations.)

5. Evaluation Procedures and Computer Programs

- *Limitation of Relative Statistical Weight* method (W.L. Zipf, M.U. Rajput, and T.D. Macmahon)
- Program *LWEIGHT* (Developed at NPL and LBNL)
- Program *EMISSION* (Developed at PTB)
- Program *EC-CAPTURE* (Developed at PTB and LBNL)

6. Current Status of Evaluations

- **259** selected radionuclides
- As of October 1, 2003:
 - 99** evaluations have been produced
 - 29** evaluations have been published in the French *Table of Radionuclides*
 - 74** evaluations are available on the WWW (<http://www.bnm.fr/bnm-Inhb/DDEP.htm>)
 - 5** evaluations are in progress

7. DDEP, IAEA-CRP, and ENSDF

- The DDEP Collaboration has evaluated most of the radionuclides included in the IAEA-CRP list.
- A large fraction of the data presented in DDEP evaluations is from ENSDF
- Very few DDEP evaluations have been included in ENSDF. Main difficulties:

Different data formats: Microsoft Access for DDEP, 80-character ASCII records for ENSDF

Different theoretical conversion coefficients: Rösel values in DDEP, Hager and Seltzer values in ENSDF

Integrating DDEP and ENSDF

- Using the same theoretical internal conversion coefficients (“Dirac-Fock Internal Conversion Coefficients”, I.M. Band et al., *Atom. Data and Nucl. Data Tables* **81**, 1 (2002)?)
- Using new software (from Saclay) to produce ENSDF data sets from DDEP evaluations
- Establishing a new database (similar to XUNDL) in Brookhaven for DDEP evaluations in ENSDF format

Nucleide

The screenshot shows the Nucleide software interface for Gallium-66. At the top, the element symbol 'Ga' is displayed in a box, with the atomic number 31 and mass number 66. To the right, the name 'Gallium' is written twice. Below this is a 'General Information' section with a 'Daughter Choice' set to 'Zn66'. The interface is divided into two main panels: 'INPUT' (red background) and 'CONSULT' (blue background). The 'INPUT' panel has sub-sections for 'Constants' (Half-lives, Q, Atomic) and 'Emissions' (Electrons, Photons, α , Elec. Capt.). The 'CONSULT' panel has buttons for 'Dosimetry', 'Auger Spectra', 'Activity <> Mass', and 'External'. At the bottom, the version '2002- mars' and file path 'C:\saisinuc02\bdvide_2000.mdb' are shown, along with the logos for CEA and BNM (Laboratoire National Henri Becquerel).

66
Ga
31 35
Gallium
Gallium

General Information

Daughter Choice **Zn66**

INPUT

Constants

Half-lives, Q Atomic

Emissions

Electrons Photons

α Elec. Capt.

CONSULT

Dosimetry

Auger Spectra

Activity <> Mass

External

Version 2002- mars C:\saisinuc02\bdvide_2000.mdb

cea Laboratoire National Henri Becquerel BNM

ENSDF Data set

```
66ZN      66GA EC DECAY                      2002BA38,1994EN02
66ZN C   Others: 2002Ga20, 1971Ca14, 1970Ph01.
66ZN  N 1.0                      1.0          1.0          1.0
66ZN  L0                          0              STABLE
66ZN  E                          50             4 0.47       4 7.88
66ZN 2 E CK=0.8850 15$CL=0.0978 12$CM+=0.0161 4
66ZN  L1039.2268 212                      1.65 PS      6
66ZN  G 1039.220  3 37                     3 E2
66ZN  L1872.7633 242                      0.19 PS      7
66ZN  G 833.5324  215.9                    5 M1+E2      -1.6        2
66ZN  G 1872.740  6 0.0229 24 [E2]
66ZN  L2372.352  4 0
66ZN  E                          0.30         3 0.038      3 8.46
66ZN 2 E CK=0.8847 15$CL=0.0980 12$CM+=0.0162 4
66ZN  G 499.590   6 0.0048 12E2+M3
66ZN  G 1333.112  5 1.17                   9 E2
```