



# ENSDF Modernization

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Not export controlled



# Outline

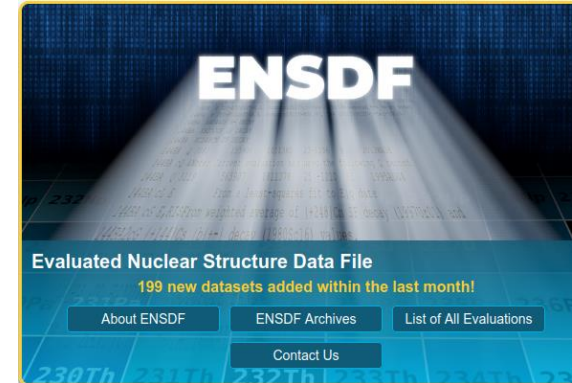
- Review of the modernization project
- Updates on the JSON format
- Other thoughts on modernization

# How it started

The design of ENSDF effectively envisions two kinds of users:

1. Evaluators
2. Journal readers

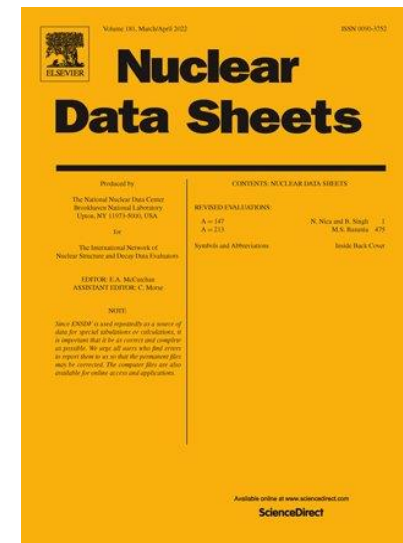
# Here be evaluators



```

143BA L 0.0 5/2- 14.5 3 A
143BAX L XREF=ABC
143BA2 L 46-100
143BA CL JShfs (1988we07,1983mu12,1981ne06), [p from analysis of [m and
143BA2CL Large negative Q. They suggest decoupled configuration with
143BA2CL main components [n(1/2-)(530)] and [n(1/2-)(532)]
143BA CL T$from vt av (same as LMI based on T[1/2-]=14.33 s [18] (1986ok03),
143BA2CL 15.2 s [12] (1979en02), 15.17 s [138] (1976Am2w), 14.5 s [15]
143BA2CL (1978Pm1), 13.2 s [13] (1969Ru14), 12 s [13] (1962Ma36). See also
143BA2CL 1973Ta13
143BA2 L MOMM1=+0.443 11 (1968we07,2011S2Z)
143BA CL MOMM1 Other: +0.454 [108] hfs (1983Mu12)
143BA3 L MOMM2=-0.88 2 (1988we07,2011S2Z)
143BA CL MOMM2 Other: -0.81 [17] hfs (1983Mu12)
143BA L 33.29 3(1/2)-
143BAX L XREF=AC
143BA CL J [g to 5/2- is E2, log[ (If)-6.0 via 3/2- parent
143BA G 33.46 100 E2 125.7
143BA CG MSFrom [a(L)(exp)=107 [121] and ce(L)/ce(M)[75 ln (+143)Cs [b(-)]
143BA2CG Decay, after normalizing electron and [g-ray intensities using
143BA2CG [a(K)(117)]g, E2, theory=ub,747.
143BAS G LC=98.9 142MC=21.8 3SNC+=+5.05 7
143BAS G NC=4.48 750C=0.569 85PC=0.000631 9
143BA L 117.358 14 9/2- 3.5 ns B A
143BAX L XREF=ABC
143BA2 L MOMM1=+0.5 3 (1995sm05,2011S2Z)
143BA CL T$from LMI based on T[1/2]=2.6 ns [18] (1995sm05), 3.8 ns [112] from
143BA2CL [b(-)] decay (1979sc11), 6 ns [12] from (+252)cf SF (1974cl2X).
143BA2CL Other: 2095fo17.
143BA CL [m from g-factor (1995sm05)
143BA G 117.32 5 100 E2 1.094
143BAB G BE2W=1, BE2 4
143BAS G KC=0.741 135LC=0.278 49MC=0.0605 95NC=+0.01432 21
143BAS G NC=0.01260 1890C=0.001686 245PC=3.52E-5 5
143BA CG M [a(K)(exp)=0.846 [125], [a(L)(exp)=0.378 [1113], [a(M)(exp)=0.102
143BA2CG [135]
  
```

# Everyone else





# Problems for non-experts

The 80-column ENSDF format is hard to use

- For the standard one-card records:
  - No delimiters - must remember field widths
  - No labels - must remember field locations
  - Inconsistent units - must remember what/where
  - Asymmetric errors almost never supported
- For the continuation items:
  - Labels can be confusing, often used inconsistently
  - Units are not allowed
  - Multiple ways to indicate limits or approximations

```
143BA L 0.0          5/2-          14.5 S   3          A
143BAX L XREF=ABC
143BA2 L %B=-100
143BA cL J$hfs (1988We07,1983Mu12,1981Ne06), |p from analysis of |m and
143BA2cL large negative Q. They suggest decoupled configuration with
143BA3cL main components |n(1/2-[530]) and |n(3/2-[532])
143BA cL T$from wt av (same as LWM) based on T{-1/2}=14.33 s {I8} (19860k03),
143BA2cL 15.2 s {I2} (1979En02), 15.17 s {I38} (1976AmZW), 14.5 s {I5}
143BAxcL (1978Pa01), 13.2 s {I3} (1969Ru14), 12 s {I3} (1962Wa36). See also
143BAxcL 1973Ta13
143BA2 L MOMM1=+0.443 11 (1988We07,2011StZZ)
143BA cL MOMM1 Other: +0.454 {I20} hfs (1983Mu12)
143BA3 L MOME2=-0.88 2 (1988We07,2011StZZ)
143BA cL MOME2 Other: -0.81 {I7} hfs (1983Mu12)
143BA L 33.29      3(1/2)-
143BAX L XREF=AC
143BA cL J          |g to 5/2- is E2, log| {Ift}=6.0 via 3/2+ parent
143BA G 33.46      100      E2          125.7
143BA cG M$From |a(L){exp}=107 {I21} and ce(L)/ce(M)|?5 in {+143}Cs |b{+-}
143BA2cG Decay, after normalizing electron and |g-ray intensities using
143BA3cG |a(K)(117|g, E2, theory)=0.747.
143BAS G LC=98.9 14$MC=21.8 3$NC+=5.05 7
143BAS G NC=4.48 7$OC=0.569 8$PC=0.000631 9
143BA L 117.368   24 9/2-          3.5 NS   8          A
143BAX L XREF=ABC
143BA2 L MOMM1=+0.5 3 (1999Sm05,2011StZZ)
143BA cL T$from LWM based on T{-1/2}=2.6 ns {I8} (1999Sm05), 3.8 ns {I12} from
143BAxcL |b{+-} decay (1979Sc11), 6 ns {I2} from {+252}Cf SF (1974ClZX).
143BA2cL Other: 2005Fo17.
143BA cL          |m from g-factor (1999Sm05)
143BA G 117.32   5 100      E2          1.094
143BAB G BE2W=1.0E+2 4
143BAS G KC=0.741 11$LC=0.278 4$MC=0.0605 9$NC+=0.01432 21
143BAS G NC=0.01260 18$OC=0.001686 24$PC=3.52E-5 5
143BA cG M          |a(K)exp=0.846 {I25}, |a(L)exp=0.378 {I113}, |a(M)exp=0.102
143BAxcG {I35}
```

# New paradigm

1. Evaluators interact with ENSDF via an editor (c.f. upcoming talk by D. Mason)
2. Human readers interact with ENSDF via PDFs
3. Computational users interact with ENSDF via new JSON format

## Benefits:

- The representation of the data is decoupled from the data itself
  - E.g. evaluators do not have to worry about format changes, the editor handles those details
- JSON enjoys widespread adoption in computing
  - Much of the tool-development work is done for us

# What is JSON?

- A highly structured data interchange format
- Governed by a simple set of rules:
  - Data entries are key-value pairs
  - Keys are (unique) strings
  - Values can have three types:
    - Basic: string, integer, number, boolean, NULL
    - Object: A collection of key-value pairs enclosed in { }
    - Array: An ordered list of values enclosed in [ ]
- Trivially easy to deserialize

affiliations.json

```
[
  {
    "institution": "University of Nowhere",
    "address": {
      "street": "University Ave",
      "number": 1,
      "zip": 12345
    },
    "presentAddress": true
  }
]
```

# JSON with Python

deserialize.py

```
import json

with open("affiliations.json") as jsonfile:
    jsondata = json.load(jsonfile)
    for item in jsondata:
        print(item["institution"])
```

```
ensdf@nndc:~$ python deserialize.py
University of Nowhere
```

affiliations.json

```
[
  {
    "institution": "University of Nowhere",
    "address": {
      "street": "University Ave",
      "number": 1,
      "zip": 12345
    },
    "presentAddress": true
  }
]
```



# Updates on the new format

The new files are available at <https://www.nndc.bnl.gov/ensdf-json>

NB: These are still considered a beta release

# Organization

- Adopted dataset
  - All other datasets (decays, reactions, comments)
  - Header (Z, A, ...)
  - Comments
  - Various info (e.g. Q-values)
  - Levels table
    - Level properties (energy, spin-parity, ...)
    - Cross-link to radiation tables (alpha, beta, gamma...)
  - Radiation (alpha, beta, gamma...) tables
    - Radiation properties (multipolarity, hindrance factor, ...)
    - Cross link to levels table

# Datasets

There are currently 14 defined types of datasets in the JSON format  
The number of datasets in each category is given in parentheses

adopted (3411)	general reaction (7001)
alpha decay (831)	isomer decay (589)
beta decay (2369)	neutron capture (608)
charge exchange (140)	prompt-particle decay (49)
coulomb excitation (391)	general decay (266)
delayed-particle decay (280)	transfer (2570)
fluorescence (200)	comments (276)

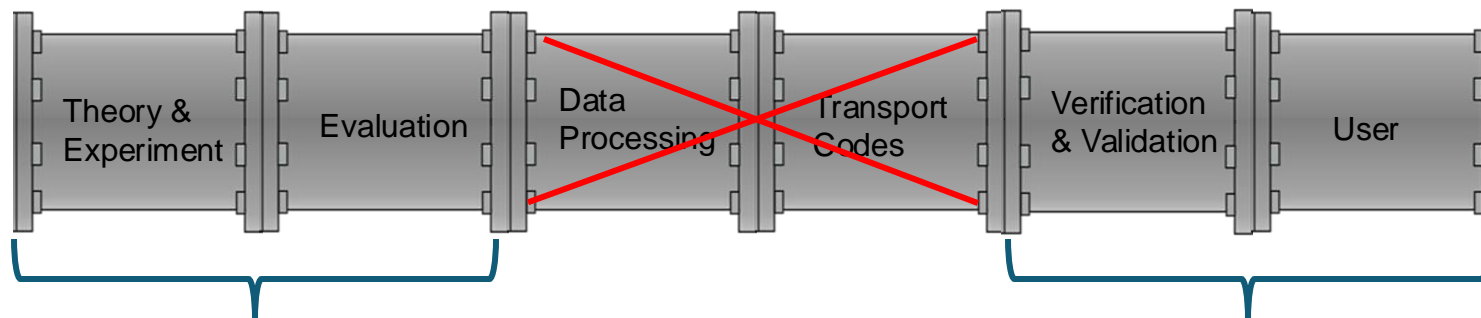
# Other updates

- Addressing feedback from NSDD meeting
  - For each nuclide, the Adopted dataset holds all other datasets in a single file
  - For error checking, individual quantities hold a string version of their numerical values
- XUNDL has been converted (11,066 datasets)
  - Big thanks to Ben Shu for this effort
- Unit tests are being developed for the JSON schema validator

# Other thoughts on modernization

Disclaimer: These aren't proposals, just brainstorming

# Evaluating the ENSDF evaluation pipeline



Graphic shamelessly stolen from Gustavo Nobre

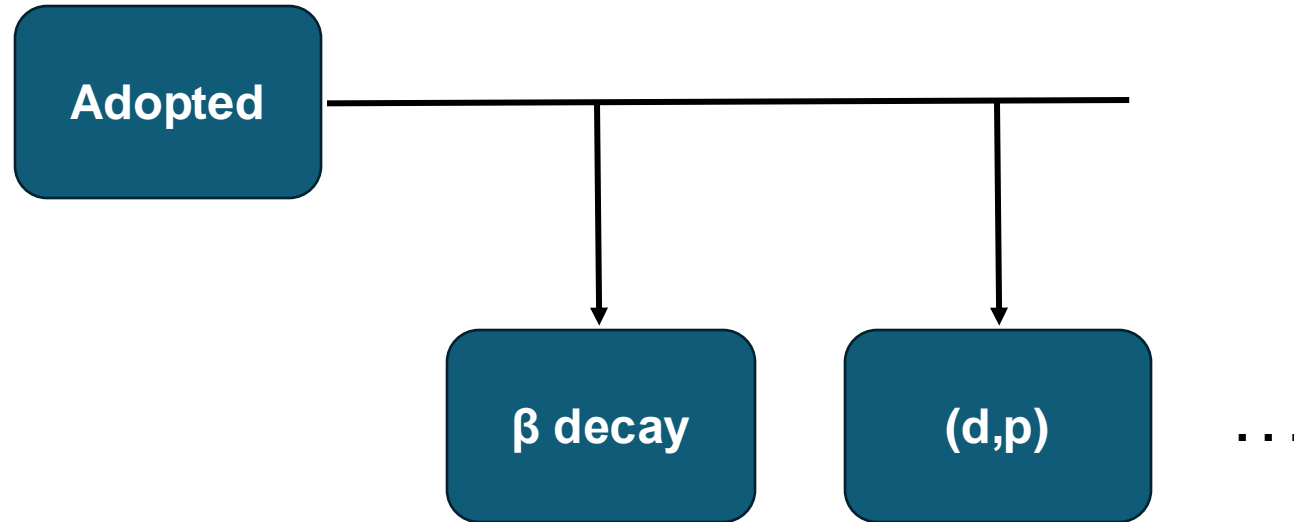
## Other opportunities?

- More thorough documentation of individual data sources?
- Backporting/preserving "bibliographic" info relevant to ENSDF?

## ENSDF modernization goals

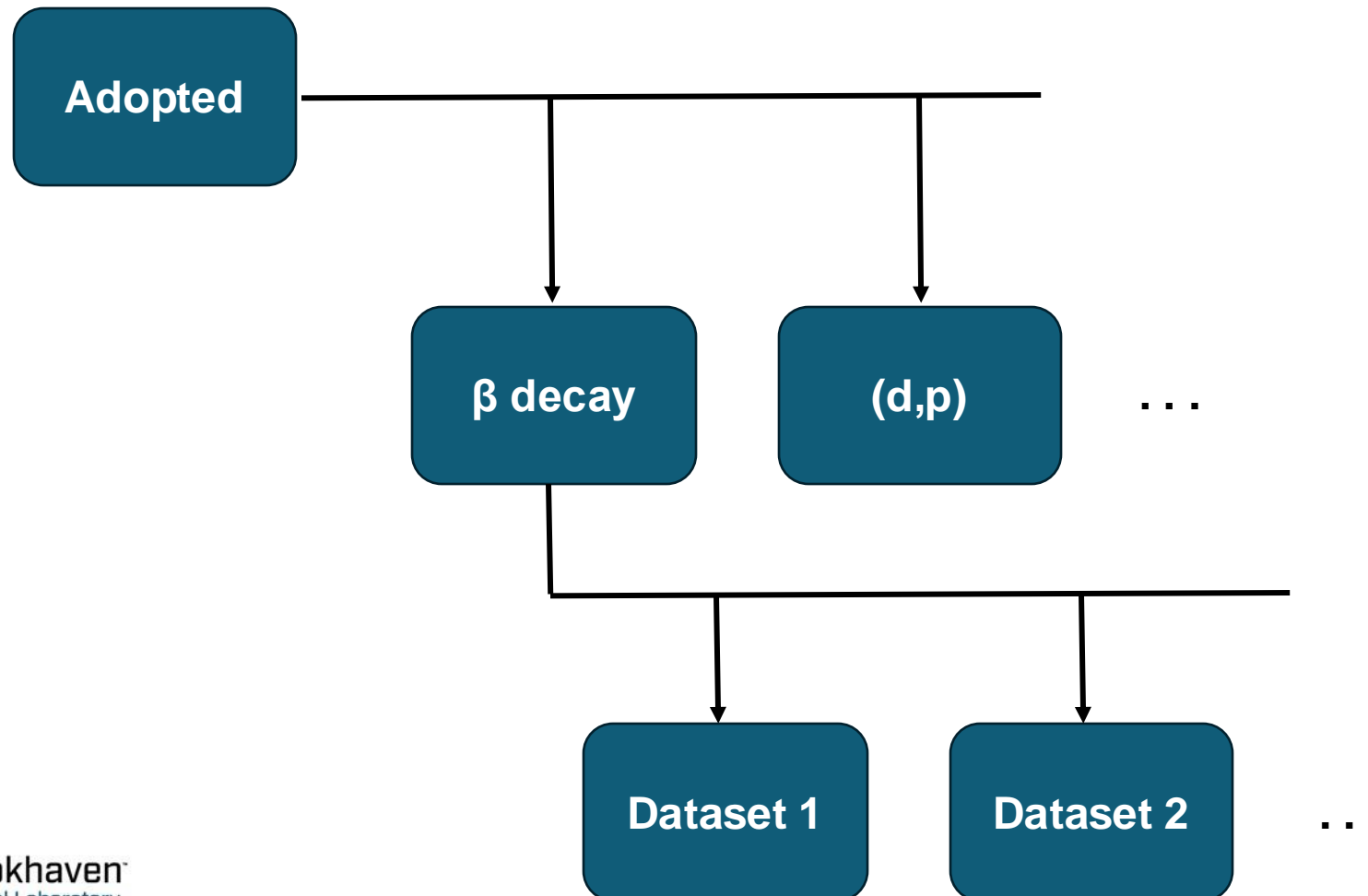
- Adopt a modern format for easy tooling
- Convenient for users
- Enable expansion of the format for emerging needs

# Documenting data sources



- Currently, ENSDF is a two-level system
  - Adopted
  - All other datasets

# Documenting data sources



- Currently, ENSDF is a two-level system
  - Adopted
  - All other datasets
- We could add another layer for individual publications
  - Similar to XUNDL
- Potential benefits
  - Data encapsulation
  - Simple to import new XUNDL datasets



# Better bibliographies

Quick Search | Text Search | Indexed Search | Keynumber Search | DOI Search

Author  
'Brown', 'B.A.Brown', 'Brown,B.A.'

Nuclide  
31Na, 144Xe, etc. 240Pu

Reaction  
n,g or (n,g) or (16O,16O)

Filters

Publication Year: [ ] to [ ]

Entries added since: [ mm / dd / yyyy ]

Subject / Reference: Measured / Experimental

Primary Only  EXFOR Data Available

Search Reset

Download New Tab Copy Key Numbers Only

Printer-Friendly

- search-type: 'quick'
- subject-type: 'measured'
- nuclide: '240Pu'
- page: '1'
- size: '100'

441 results found!

Some nuclei generate *many* publications  
How to review all this literature?

- Ideal case, you have records from last evaluation. Not true for newer evaluators.
- Assume previous evaluation caught everything relevant? Dangerous.
- Read everything; very inefficient!

Many references in an NSR search are not relevant to ENSDF for a given nuclide. It would be helpful to capture this information as part of a mass-chain evaluation.

- One possibility: add new field to NSR entries to mark a key number as (ir)relevant for a given ENSDF evaluation

# The end!

Questions? Comments?