

# Development of multi-group neutron activation cross-section library from JENDL/AD-2017

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## Introduction

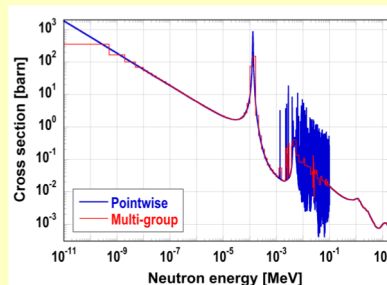
- ❑ JENDL Activation Cross Section File for Nuclear Decommissioning 2017 (**JENDL/AD-2017**) [1] was released in 2018.
- ❑ This file includes neutron-induced nuclear reaction cross-sections of 311 nuclides from  $10^{-5}$  eV to 20 MeV.
- ❑ Dr. Okumura et al. developed a multi-group neutron activation cross-section library (**MAXS2015**) based on the nuclear data libraries JENDL-4.0 and JEFF-3.0/A for activation calculations in nuclear facility decommissioning [2].
- ❑ Thus a multi-group neutron activation cross-section library (**MAXS/AD-2017**) with the same format as MAXS2015 has been developed from JENDL/AD-2017 for validation of JENDL/AD-2017 and activation calculations.

## How to make MAXS/AD-2017

- ❑ JENDL/AD-2017 includes total production cross sections (MF3) of radioactive and stable nuclides, branching ratios (MF9) and partial production cross sections (MF10) for the ground and isomer states of nuclides.
- ❑ JENDL/AD-2017 has four versions.
  - MF3, MF9 and MF10 at 0 K
  - MF3, MF9 and MF10 at 293.6 K
  - MF3 and MF10 at 0 K (for NJOY processing)
  - MF3 and MF10 at 293.6 K (for NJOY processing)
- ❑ MAXS-2015 was produced with the NJOY2012 [3] code.
- ❑ However it was found that **GENDF files produced with the group module in NJOY2012 did not include production cross sections to isomer states.**
- ❑ Then the **PREPRO 2018** [4] code was adopted for producing a group-wise file of JENDL/AD-2017 (**MF3, MF9 and MF10 at 0 K**).
- ❑ The following modules in PREPRO 2018 were used; **ENDF2C, LINEAR, RECENT, SIGMA1, ACTIVATE, FIXUP, DICTIN, GROUPIE.**
- ❑ The calculation conditions are as follows,
  - Temperature : **300 K**
  - Group structure : **199 groups (VITAMIN-B6)**
  - Weighting spectrum : **Maxwell + 1/E + Fission (-2)**
  - **Infinite dilution** cross section
- ❑ The produced group-wise file of JENDL/AD-2017 was converted to **MAXS/AD-2017** of the MAXS format [2] with a small program.
- ❑ The following issues were pointed out in this processing
  - No information of decay data (MF8) in the capture reaction of  $^{187}\text{W}$  and  $^{193}\text{Os} \rightarrow$  Add
  - The MT number of the (n,t) reaction of  $^6\text{Li}$  is changed from 105 to 107 for ORIGEN-S.
  - The MAXS format includes **no data for the (n,n') reaction**  $\rightarrow$  MAXS files with data for the (n,n') reaction are also produced (**Extended MAXS/AD-2017**), though ORIGEN-S cannot treat the (n,n') reaction.
- ❑ A similar procedure for a **DCHAIN-SP library** was also established, and was provided to the PHITS group. Users can use the DCHAIN-SP library of JENDL/AD-2017 in the latest PHITS (**PHITS3.16**).

← Capture cross section of  $^{59}\text{Co}$  in JENDL/AD-2017 (Red line : MAXS/AD-2017)

```
#MAXS-xs Library
# Nuclide ID & Name
270590
Co059
# background XS (sigz<0:effective XS)
1.000000E+10
# Temperature (K)
3.000000E+02
#Number of Energy groups (NGN)
199
#Number of Reaction Types (NMT)
11
#
# No. Energy MT numbers
```



```
16 22 28 32 41 102 103 104 105 107 112
# 2n Co058 na Mn055 np Fe058 nd Fe057 2np Fe057 g Co060 p Fe059 d Fe058 t Fe057 a Mn056 pa Cr055
1 1.964000E+07 7.882728E-01 2.628246E-02 1.512986E-01 1.256140E-04 3.566710E-06 5.124860E-04 3.529885E-02 2.690402E-02 1.156921E-03 1.648094E-02 3.545590E-06
2 1.733200E+07 7.809816E-01 1.699300E-02 1.299114E-01 2.515740E-07 0.000000E+00 6.795890E-04 3.812621E-02 2.143026E-02 5.896660E-04 2.089150E-02 1.429740E-07
3 1.690500E+07 7.732319E-01 1.426557E-02 1.234424E-01 1.600430E-08 0.000000E+00 7.415650E-04 3.930724E-02 1.954275E-02 4.555710E-04 2.226716E-02 5.172710E-08
4 1.648700E+07 7.582623E-01 1.012589E-02 1.132496E-01 8.194510E-09 0.000000E+00 8.278450E-04 4.112929E-02 1.694981E-02 3.063550E-04 2.451590E-02 1.102070E-08
}
197 5.000000E-03 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 1.002084E+02 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00
198 2.000000E-03 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 1.660360E+02 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00
199 5.000000E-04 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 3.544969E+02 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00
200 1.000000E-05 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00 0.000000E+00
```

Example of MAXS/AD-2017 ( $^{59}\text{Co}$ )

## Concluding Remarks

- ❑ A multi-group neutron activation cross-section library (MAXS/AD-2017) with the MAXS format was developed from JENDL/AD-2017 for activation calculations in nuclear facility decommissioning.
- ❑ Next MAXS/AD-2017 will be converted to ORIGEN libraries and be tested with the JPDR decommissioning data. Then MAXS/AD-2017 will be released.

## References

- [1] <https://www.ndc.jaea.go.jp/ftpnd/jendl/jendl-ad-2017.html>
- [2] K. Okumura, K. Kojima, K. Tanaka, "Development of multi-group neutron activation cross-section library for decommissioning of nuclear facilities," Proc. of 2014 Symposium on Nuclear Data, p. 43, JAEA-Conf 2015-003(2016).
- [3] R. E. MacFarlane, D. W. Muir, R. M. Boicourt, A. C. Kahler, "The NJOY Nuclear Data Processing System, Version 2012," LA-UR-12-27079, Los Alamos National Laboratory (2012).
- [4] <https://www.nds.iaea.org/public/endf/prepro2018/>