



Fission Yield Interpolation and Energy Dependent Fission Product Yield Covariance Format

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Nuclear Data Week 2025 – January 6 – 9, 2026

LA-UR-26-20127



Fission Product Yield Interpolation

The current format and manual recommend linear interpolation between incident energies

- From the ENDF-6 manual:

```
[MAT, 8, MT/ ZA, AWR, LE+1, 0, 0]HEAD
[MAT, 8, MT/ E1, 0.0, LE, 0, NN, NFP/ Cn(E1) ]LIST
[MAT, 8, MT/ E2, 0.0, I, 0, NN, NFP/ Cn(E2) ]LIST
[MAT, 8, MT/ E3, 0.0, I, 0, NN, NFP/ Cn(E3) ]LIST
-----
[MAT, 8, 0 /0.0, 0.0, 0, 0, 0]SEND
```

I_i Interpolation scheme (see paragraph on Two-dimensional Interpolation Schemes in Section 0.5.2) to be used between the E_{i-1} and E_i energy points.

Yields for the same fission product nuclides should be given at each energy point. This will facilitate interpolation of yield data between incident energy points. Also, a linear-linear interpolation scheme should be used.



But most evaluators recommend *not interpolating at all* for previous evaluations

- The incident energies are “effective” energies representing a spectrum
 - Are there any cases for which this is not true?
- The new LANL FPY’s are energy-dependent, and linear interpolation can be performed between the incident energies
- How can this be communicated?



ENDF-6 Format Change

- A new flag in the HEAD record to indicate if interpolation is recommended:
 - 0 : not recommended (“effective” energies)
 - 1 : recommended by the law given in the LIST records

```
[MAT, 8, MT/ ZA, AWR, LE+1, 0, 0, 0]HEAD
[MAT, 8, MT/ E1, 0.0, LE, 0, NN, NFP/ Cn(E1) ]LIST
[MAT, 8, MT/ E2, 0.0, I, 0, NN, NFP/ Cn(E2) ]LIST
[MAT, 8, MT/ E3, 0.0, I, 0, NN, NFP/ Cn(E3) ]LIST
-----
[MAT, 8, 0 /0.0, 0.0, 0, 0, 0, 0]SEND
```



GNDS Format Change

- The format currently does not hold any interpolation information (not in an XYs2d)
- Could an interpolation attribute be added to the `<incidentEnergies>` node?
- Or should `<incidentEnergies>` be changed to an XYs2d?
- This will be brought up at the EG-GNDS meeting in May 2026

```
<fissionFragmentData>
  <delayedNeutrons>...</delayedNeutrons>
  <fissionEnergyReleased>...</fissionEnergyReleased>
  <productYields>
    <productYield label="eval">
      <nuclides> V66 V67 Cr66 ... </nuclides>
      <elapsedTimes>
        <elapsedTime label="initial">
          <time>
            <double label="initial" value="0.0" unit="s"/></time>
        <incidentEnergies>
          <incidentEnergy label="0">
            <energy>
              <double label="0" value="500000.0" unit="eV"/></energy>
            <yields>
              <nuclides href="../../../../nuclides"/>
              <values>1.87064e-17 2.02328e-18 6.25013e-13 ... </values>
              <uncertainty>
                <covariance>
                  <array shape="1267,1267" compression="diagonal">
                    <values>...</values></array></covariance>
                </uncertainty>
              </yields>
            </incidentEnergy>
            <incidentEnergy label="1">...</incidentEnergy>
          </incidentEnergy>
        </incidentEnergies>
      </elapsedTime>
    </productYield>
  </productYields>
</fissionFragmentData>
```





Energy Dependent Fission Product Yield Covariances

A format is needed for FPY Covariances

- File 8 holds fission product yield uncertainties, but not correlations
- Considerations:
 - Backward compatible with uncertainties in MF8
 - Not a compact format
 - Allow cross-correlations between energies and MT's
 - Hold incident-energy grouped covariances (like File 33)



ENDF-6 Format Proposal

- Leave uncertainties in File 8 in the point-wise structure
 - Allow linear interpolation on the absolute uncertainty values
- Put correlations only in File 38 in a grouped structure
- Require the same list of nuclides in both for all energies and MT's

The structure of a section is:

```
[MAT, 38, MT/ ZA, AWR, 0, 0, NB, 0] HEAD
[MAT, 38 ,MT/ 0.0, 0.0, 0, 0, NG+1, 0/ El]LIST
[MAT, 38 ,MT/ 0.0, 0.0, 0, 0, 2*NFP, NFP/ ZAFPl, FPSl]LIST
    <NB subsections>
[MAT, 38, MT/ 0.0, 0.0, 0, 0, 0] SEND
```

Subsections

Each subsection represents a single submatrix of the full correlation matrix, associated with a given row and column energy group. Each subsection has the following structure:

```
[MAT,38, MT/ IROW, ICOL, LS, 0, NT, 0/ Cl]LIST
```



GNDS Format Proposal

- The FPY objects already hold covariances within a single incident energy
- For cross-correlations (different energies):
 - Make use of `rowData` and `columnData` attributes
- This will be brought up at the EG-GNDS meeting in May 2026

```
<fissionFragmentData>
  <delayedNeutrons>...</delayedNeutrons>
  <fissionEnergyReleased>...</fissionEnergyReleased>
  <productYields>
    <productYield label="eval">
      <nuclides> V66 V67 Cr66 ... </nuclides>
      <elapsedTimes>
        <elapsedTime label="initial">
          <time>
            <double label="initial" value="0.0" unit="s"/></time>
        <incidentEnergies>
          <incidentEnergy label="0">
            <energy>
              <double label="0" value="500000.0" unit="eV"/></energy>
            <yields>
              <nuclides href="../../../../nuclides"/>
              <values>1.87064e-17 2.02328e-18 6.25013e-13 ... </values>
            <uncertainty>
              <covariance>
                <array shape="1267,1267" compression="diagonal">
                  <values>...</values></array></covariance>
              </uncertainty>
            </yields>
          </incidentEnergy>
        <incidentEnergy label="1">...</incidentEnergy>
      </incidentEnergies>
    </elapsedTime>
  </productYield>
</productYields>
```



Example File with 3 Energy Groups

- ^{235}U energy-dependent FPY's
- 21 energies in File 8
- 859 nuclides, consistent across all energies and MT's
- 3 energy bins in File 38
- Cross-correlations between all nuclides and energy bins
- Correlations for Independent and Cumulative, no cross-correlation between them
- File size: 144MB



Saving Space

- Store a limited set of eigenvalues and eigenvectors of the matrices
 - The evaluators would choose how many are needed
- Relax the requirement of having the same list of nuclides in File 8 and File 38
 - The evaluators would provide a default correlation value to use

