

# TXT<sub>2</sub>NTAB

## Polaris cross-section generation for Core analysis

Antonella Labarile, Rafael Miró Herrero

University Research Institute for Industrial, Radiophysical and Environmental Safety (ISIRYM)

Universitat Politècnica València (UPV)

---

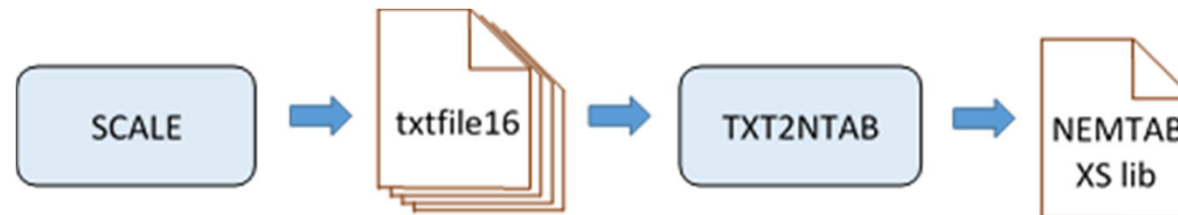
**SCALE USERS GROUP WORKSHOP**

**ORNL – July 2020**



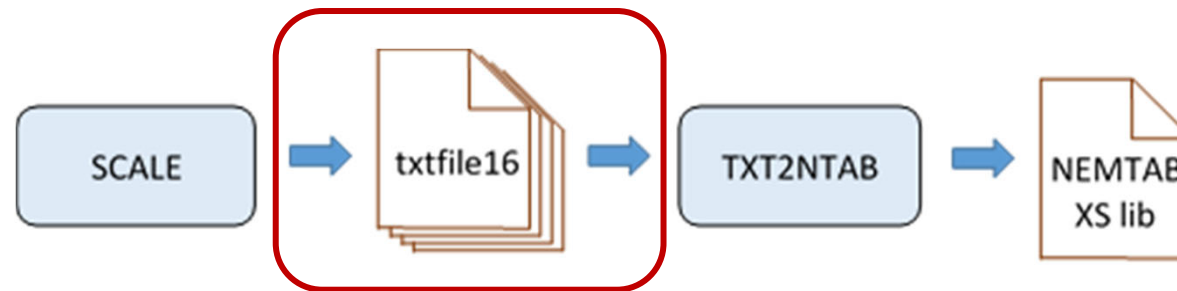
## 1. INTRODUCTION

**TXT2NTAB** is a set of **MATLAB** scripts developed to **convert** cross sections predicted by **SCALE** (**txtfile16** format) into **NEMTAB** formatted cross section libraries (to use with PARCS for example).



## 1. INTRODUCTION

TXT2NTAB is a set of MATLAB scripts developed to convert cross sections predicted by SCALE (txtfile16 format) into NEMTAB formatted cross section libraries (to use with PARCS for example).



**Each** reactor **segment** must be run in **SCALE** to provide all **txtfile16** needed for a complete core representation.

## 1. INTRODUCTION

- ❑ TXT2NTAB was used by the **UAM-LWR benchmark** organizers. To generate NEMTAB to **distribute** among benchmark participants.
- ❑ It handles **TRITON-NEWT** and **Polaris** modules, but also CASMO and SERPENT results.
- ❑ **Burn-up** and **History** dependence feature.
- ❑ It is provided with a friendly **user interface** in ASCII format.
- ❑ Optionally, it also generates **GEOM** auxiliary file for PARCS.



### Feedback parameters...

- ❑ It is the user **responsibility** to introduce proper feedback parameters.
- ❑ These are set using **branches** in SCALE.
- ❑ There are three important things to consider.
  - Feedback **parameters** to be included: Tfuel, Dmod, Crod (this is fixed in the NEMTAB format).
  - Number of **points** for each parameter. More accurate library, but more computational time.
  - The **range** for the feedback parameters should cover the range of all possible operational conditions.

## 2. GUIDE

### TRITON-NEWT

There are:

$i$  moderator density points

$j$  fuel temperature points

one *boron* concentration point

two control rod states (*in/out*)

The reactor history is constant for all branches (nominal state).

Branch index	Feedback parameters	Comment
0	hdm, htm, htf, hbo, hcr=0	Nominal branch.
1	dm1, tm1, tf1, cb1, cr0	
2	dm2, tm2, tf1, cb1, cr0	
...		
$i$	dmi, tmi, tf1, cb1, cr0	$i$ branches with dm, tm points.
...		
$i \cdot j - i + 1$	dm1, tm1, tfj, cb1, cr0	Repeat same branches $j$ times with different tf points.
$i \cdot j - i + 2$	dm2, tm2, tfj, cb1, cr0	
...		
$i \cdot j$	dmi, tmi, tfj, cb1, cr0	
$i \cdot j + 1$	dm1, tm1, tf1, cb1, cr1	Repeat $i \cdot j$ branches with control rod inserted.
$i \cdot j + 2$	dm2, tm2, tf1, cb1, cr1	
...		
$i \cdot j + i$	dmi, tmi, tf1, cb1, cr1	
...		
$2 \cdot i \cdot j - i + 1$	dm1, tm1, tfj, cb1, cr1	
$2 \cdot i \cdot j - i + 2$	dm2, tm2, tfj, cb1, cr1	
...		
$2 \cdot i \cdot j$	dmi, tmi, tfj, cb1, cr1	Total number of branches $2 \cdot i \cdot j + 1$



### POLARIS

In order to maintain the analogy with txtfile16 generated by TRITON-NEWT, the first branch specified in Polaris must define the nominal state (or reactor history).

```
state FUEL : temp=htf
state COOL : temp=htm
state COOL : dens=hdm
state COOL : boron=hbo
state BANK : in=false

'branch - instantaneous change
read branch
'   BRANCH:      NOM| Branch 1| Branch i| Branch ij | Branch ij+i CR| ...
  add FUEL : temp=htf   tf1 ... tf1 ... tfj ... ... tf1 ... ...
                        tm1 ... tmi ... tmi ... ... tmi ... ...
                        dm1 ... dmi ... dmi ... ... dmi ... ...
                        BANK : in=false ... .. true ... ..
end branch
```

### POLARIS

- File txtfile16 generated by Polaris has some **discrepancies** compared to the **txtfile16** generated by TRITON.
  
- User must **complete** some information for a couple of data blocks:
  - **Block 3:** Polaris does not write this block only for fresh fuel conditions.
  - **Block 4:** user must complete the feedback parameter information.
  
- This has been **automatized** in TXT2NTAB.



### Input files and data.

- fueltype.dat** contains the **fuel** type **distribution** for the particular reactor (radial mapping).
- burnup.dat** only provided for burn-up conditions (no fresh fuel), contains the **burn-up distribution** as a 3D map.
- input.dat** is the input data provided by the **user** to define the reactor dimensions, segment information, parameters, feedback parameters...
  - Define as many fuel assemblies and segments as needed.
  - Option for cross section extrapolation for low moderator densities.
  - Three neutronic composition options.
  - NEMTAB generation for a specific burn-up point.



### Input.dat example

- **Structured** with variables and cards.
- **Comments** are available with % symbol.
- In-code **checks** to reduce bugs.
- Whole description in user **guide**.
- And more ...

```

%radial reflector
ftyp 1          %fuel type for this assembly
segm 1 1 1     %segment number for this assembly
fuel 0 0 0     %1 for fuel segments, 0 for reflector segments, nsegm
absz -15.00 0.00 375.00 390.00 %absolute segment height, nsegm+1

%CALCULATION PARAMETERS
pola 1          %optional, 1 if txtfile16 were obtained with Polaris module, 0 otherwise, default is 0 ('
nobu 1          %optional, 1 for fresh fuel, variable burn is then omitted, default is 1
intp linear    %optional, interpolation method accepted by MATLAB: linear or spline, default and recomm
extp 0          %optional, 1 for cross section extrapolation for low moderator densities, default is 0 (:
dlim 0.17075;  %limit for extrapolation, omitted if variable extp is 0
geop 0          %optional, 1 to create GEOM file to be used with PARCS, default is 0
nkop 0          %optional, options are 0, 1 or 2. Option to compute number of neutronic composition, the:

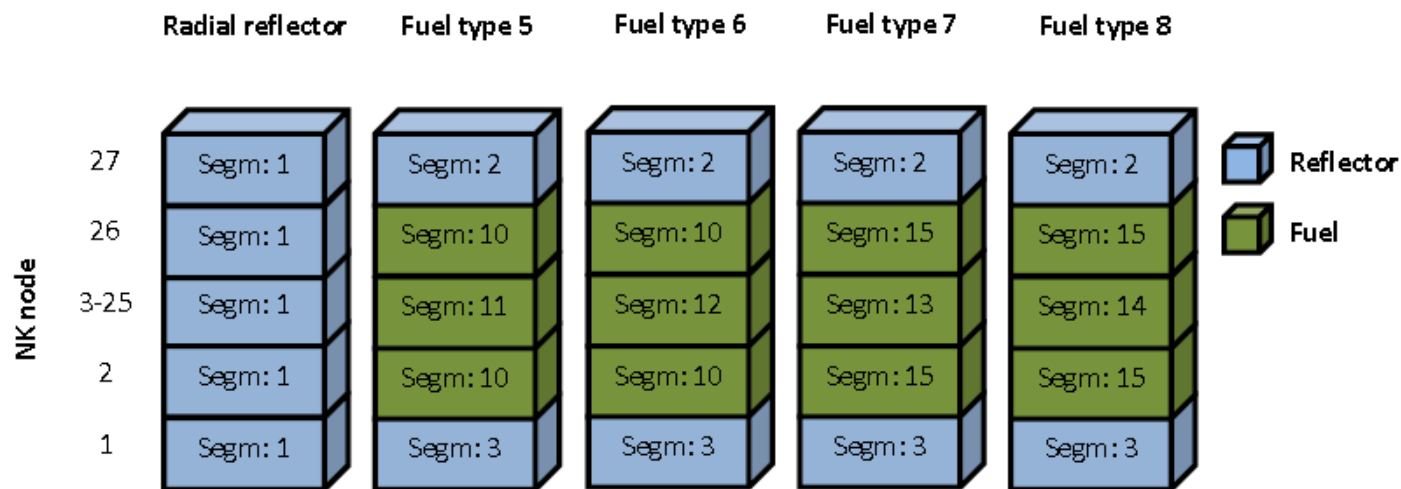
%FEEDBACK PARAMETERS
%to look up in txtfile16
dm16 0.03814 0.17753 0.45632 0.73511 0.84034 0.94281 0.99829 %moderator density, increasing order
tm16 561.4 561.4 561.4 561.4 493 393 293 %moderator temperature for each moderator den:
tf16 293 660.8 879.5 1028.6 1396.5 1764.3 2132.2 %fuel temperature, increasing order
cb16 1935 %scalar only (no variation with boron concentration yet)

```



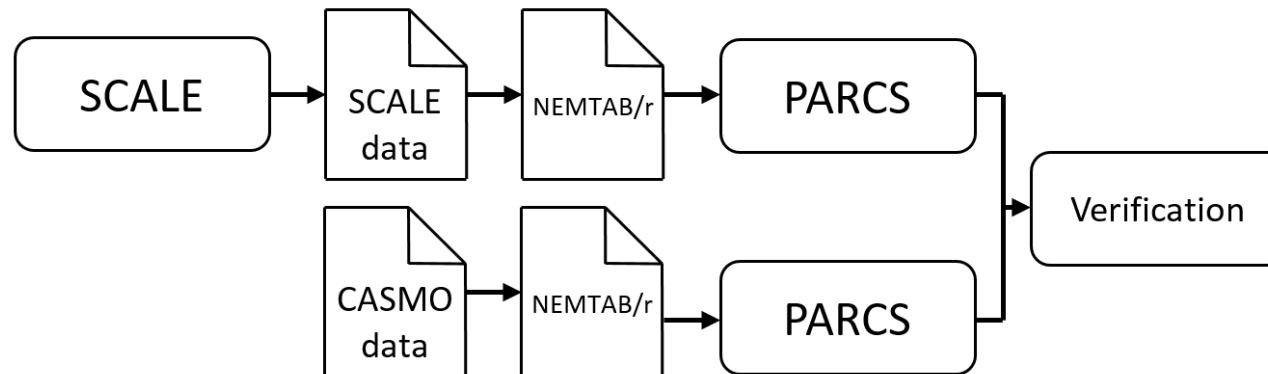
### 3. EXAMPLE CASE

- ❑ Segment information
- ❑ **Simplified** case to **test** the methodology.
- ❑ 6 fuel segments and 3 reflector segments.
- ❑ **Burn-up** conditions **implemented** (with interpolation).
- ❑ **History** conditions **implemented** (with interpolation).



### 3. EXAMPLE CASE

- ❑ **Verification** with a code-to-code comparison.
- ❑ Complementary **MATLAB** tools were developed.
- ❑ **NEMTABs** can be **obtained** out of CASMO/SCALE/SERPENT results.
- ❑ The same **PARCS model** can be run with those NEMTAB sets.

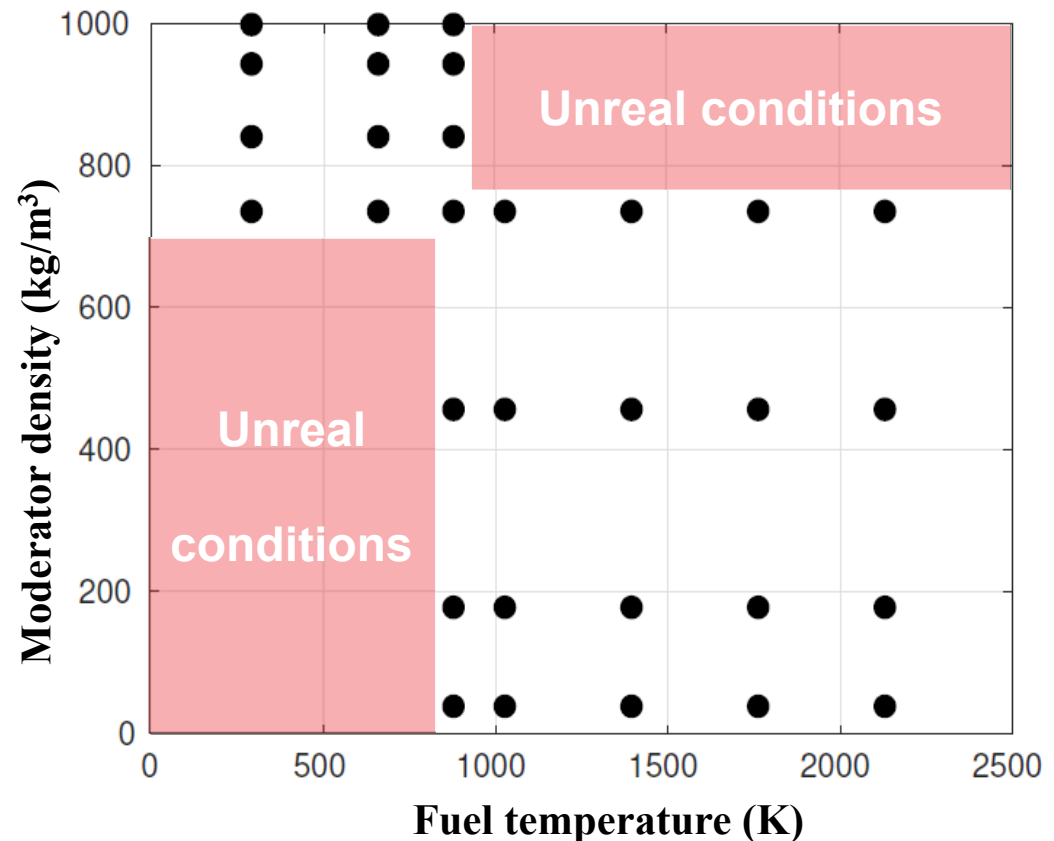


### 3. EXAMPLE CASE

☐ User must decide what **feedback parameters** (branches) to include for each segment.

☐ Some **branches** can be **discarded** to save computational time.

☐ However, there is no problem if the user wants to simulate **all branches**.



### 3. EXAMPLE CASE

Data shown hereafter are full core results obtained with **PARCSv3.2**, either with NEMTABs out of SCALE (with TXT2NTAB) or with NEMTABs out of CASMO.

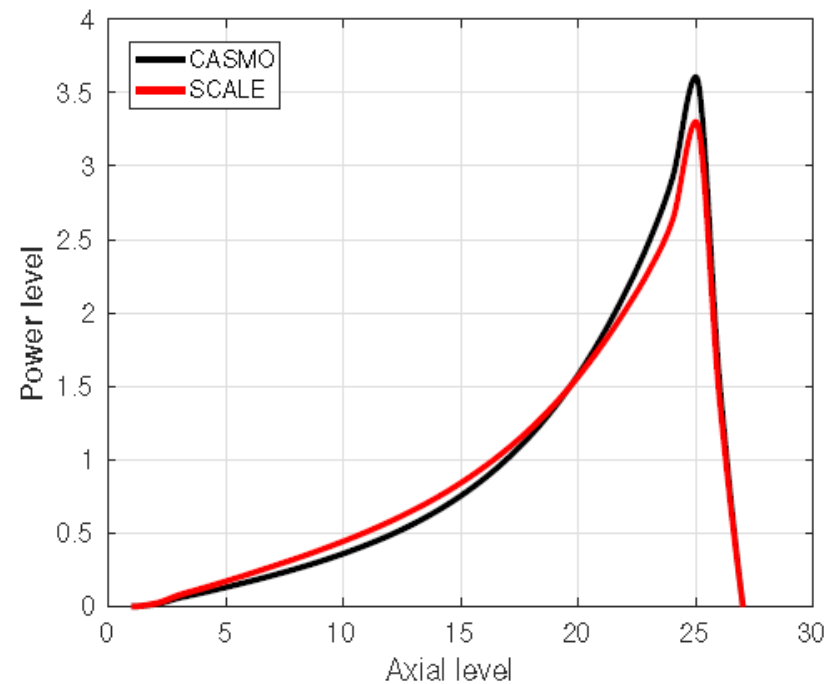
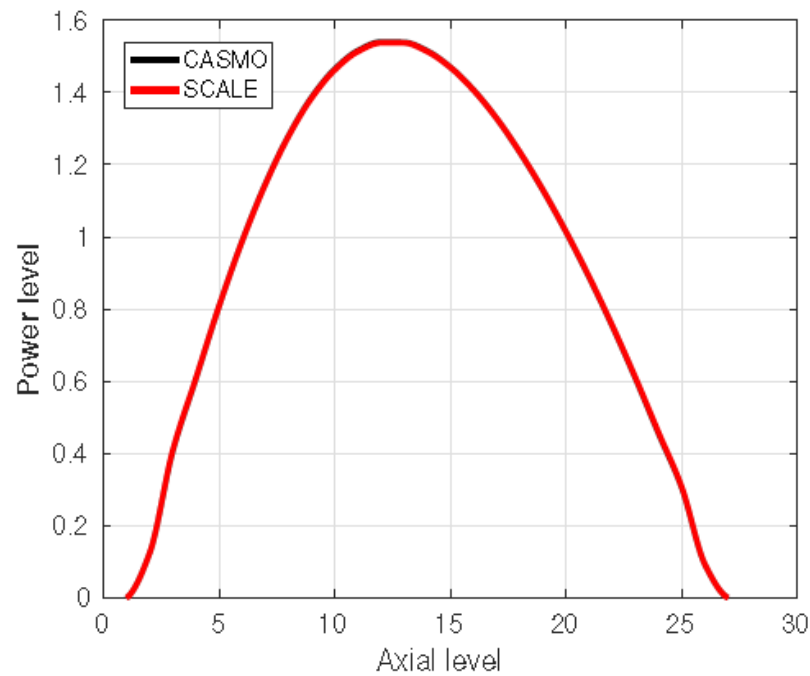
ARO	CASMO	SCALE v7-56		SCALE v7-252	
		Default	MCDANCOFF	Default	MCDANCOFF
$k_{\text{eff}}$	1.03171	1.03601	1.03417	1.03533	1.03416
<b>Error (pcm)</b>	-	429.5	246.0	362.1	<b>244.9</b>

ARI	CASMO	SCALE v7-56		SCALE v7-252	
		Default	MCDANCOFF	Default	MCDANCOFF
$k_{\text{eff}}$	0.79702	0.80023	0.79893	0.80011	0.79925
<b>Error (pcm)</b>	-	321.1	<b>191.5</b>	308.9	223.3

### 3. EXAMPLE CASE

ARO	RMS <sub>axial</sub> (%)	
	V7-56	V7-252
Default	0.24	<b>0.21</b>
MCDANCOFF	0.24	<b>0.21</b>

ARI	RMS <sub>axial</sub> (%)	
	V7-56	V7-252
Default	11.98	<b>11.53</b>
MCDANCOFF	12.09	11.56

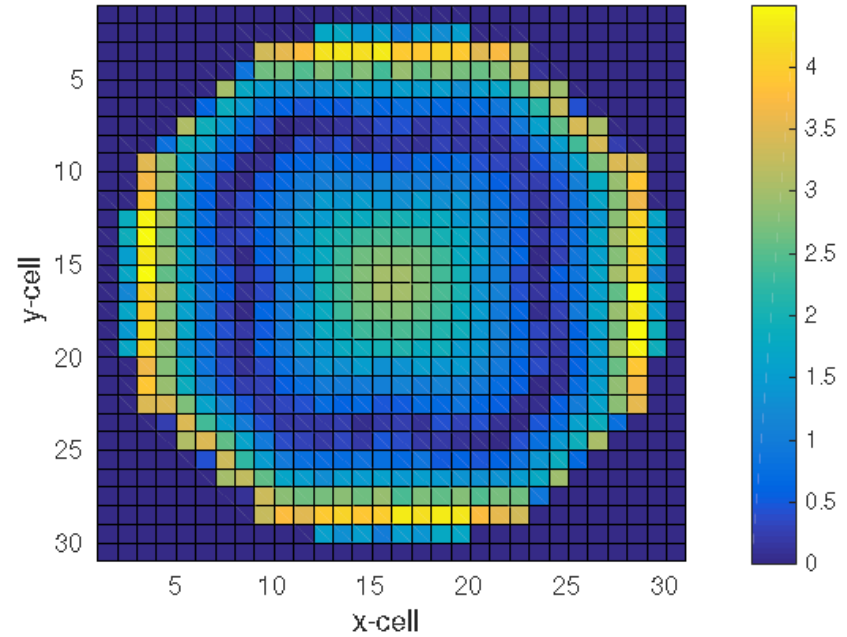
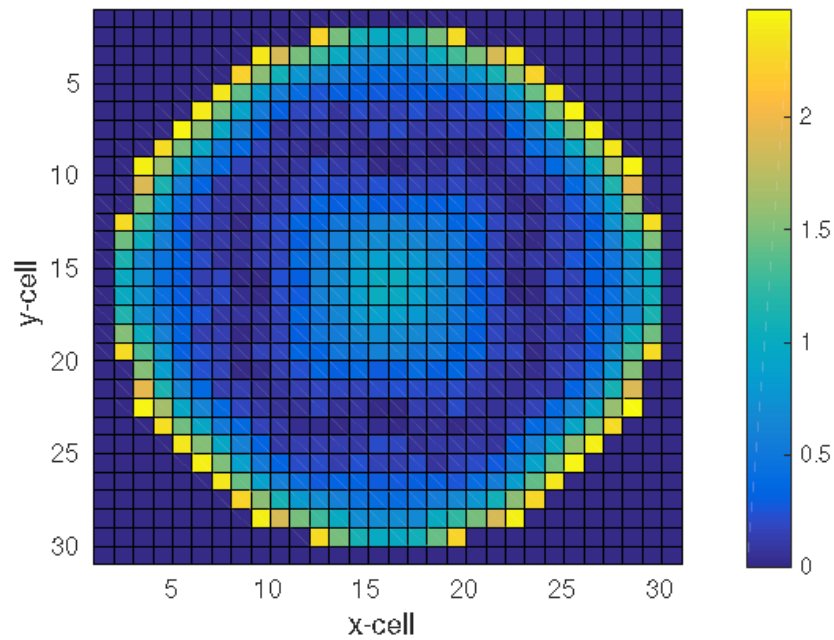


### 3. EXAMPLE CASE

ARO	RMS <sub>radial</sub> (%)	
	V7-56	V7-252
Default	0.53	0.53
MCDANCOFF	0.53	0.53

ARI	RMS <sub>radial</sub> (%)	
	V7-56	V7-252
Default	1.75	1.73
MCDANCOFF	1.74	<b>1.72</b>

#### Absolute radial error (%)





## 4. CONCLUSIONS

- ✓ TXT2NTAB program is developed to fill the **gap** between SCALE and a core simulator.
- ✓ Problem-dependent cross section libraries in NEMTAB format can be obtained from **TRITON** or **Polaris** in the SCALE code system, but also from SERPENT and CASMO code.
- ✓ The choice of **feedback parameters** is crucial in the lattice simulation.
- ✓ **Burn-up** and **history** dependencies are handle in TXT2NTAB.
- ✓ A **friendly** ASCII user **interface** is used to introduce the data with ease.
- ✓ A user **guide** is provided.



TXT2NTAB is actively being updated...

- ✓ For **U&S** analysis. Process **SAMPLER** perturbed txtfile16 to generate...
  - One set of NEMTAB/r with the **average** response.
  - One set of NEMTAB/r with their **standard deviation**.
- ✓ Improve the use of TXT2NTAB while using **Polaris** without the user intervention.
  - Block3 and Block 4 are not present in the txtfile16
  - The axial and radial discretization in Polaris slows down the simulation a lot

# TXT2NTAB

## Polaris cross-section generation for Core analysis

[alabarile@iqn.upv.es](mailto:alabarile@iqn.upv.es)

University Research Institute for Industrial, Radiophysical and Environmental Safety (ISIRYM)

Universitat Politècnica València (UPV)

---

**SCALE USERS GROUP WORKSHOP**

**ORNL – July 2020**

