enrimo

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NAME

enrimo - Investigate the influence of an enriched Mo isotope

SYNOPSIS

DESCRIPTION

```
This Perl program generates datasets for investigating the influence of an enriched Mo isotope on its associated Mo material, Mo element, and companion isotopes.

The following quantities, as functions of the enrichment level of the Mo isotope to be enriched, are calculated for a Mo material:

- Amount fractions and mass fractions of Mo and O isotopes

- Mass fractions of Mo and O elements

- Mass and number densities of the Mo material, Mo and O elements, and their isotopes

- Density change coefficients (DCCs) of Mo and O isotopes
```

OPTIONS

```
Multiple values are separated by the comma (,).

--materials=mo_mat ... (short: --mats, default: momet)
all
```

```
All of the following mo_mat's.
   momet
       Mo metal
   moo2
       Mo(IV) oxide (aka Mo dioxide)
   moo3
       Mo(VI) oxide (aka Mo trioxide)
--isotope=mo_isot (short: --isot, default: mo100)
   Mo isotope to be enriched.
   mo92
   mo94
   mo95
   mo96
   mo97
   mo98 <= Mo-98(n,g)Mo-99
   mo100 \leftarrow Mo-100(g,n)Mo-99, Mo-100(n,2n)Mo-99, Mo-100(p,2n)Tc-99m
--enri_lev_type=frac_type (short: --type, default: amt_frac)
   The fraction type to refer to the enrichment level.
   amt frac
   mass_frac
--enri_lev_range=frac_range (short: --range, default: 0,0.0001,1)
   The range of enrichment levels to be examined.
   e.g. 0.1,0.5
                    (beg,end; incre is automatically determined)
   e.g. 0,0.001,1 (beg,incre,end)
   e.g. 0,0.00001,1 (beg,incre,end)
--min_depl_lev_global=enri_lev (short: --global, default: 0.0000)
    The minimum depletion level that applies to all the nuclides
   associated \mbox{with} the designated Mo materials. Overridden, \mbox{if} given,
   by nuclide-specific minimum depletion levels.
   e.g. 0.0007
--depl order=option (short: --order, default: ascend)
   The order in which the Mo isotopes other than the to-be-enriched one
   will be depleted.
   ascend (short: asc)
       Ascending order of mass number
   descend (short: desc)
       Descending order of mass number
   random (short: rand, alt: shuffle)
       Random order
--inp=fname (short: -i)
   An input file specifying the nuclide-specific minimum depletion levels
   and the calculation precision. See the sample input file for the syntax.
   e.g. 0p9739.enr
--out_path=path (short: --path, default: the value of -isotope)
   Path for the output files.
--out_fmts=ext ... (short: -o, default: dat,xlsx)
   Output file formats.
   all
       All of the following ext's.
   dat
       Plain text
       LaTeX tabular environment
       comma-separated value
   xlsx
       Microsoft Excel 2007
    json
       JavaScript Object Notation
   yaml
--projectiles=particle ... (short: --projs, default: none)
   Reaction projectiles {f for} associating the product nuclides {f with} DCCs.
   If designated, the relevant reporting files are generated
   in addition to the default output files.
```

```
all
All of the following particles.

g
Photon <= Mo-100(g,n)Mo-99

n
Neutron <= Mo-98(n,g)Mo-99, Mo-100(n,2n)Mo-99

p
Proton <= Mo-100(p,2n)Tc-99m

--verbose (short: --verb)
Display the calculation process in real time. This will pause the shell each time a core calculation routine is called; use it only when debugging or checking part of the calculation process.

--nofm
The front matter will not be displayed at the beginning of program.

--nopause
The shell will not be paused at the end of program.
Use it for a batch run.
```

EXAMPLES

```
perl enrimo.pl --type=mass_frac --range=0,0.00001,1
perl enrimo.pl --mats=moo3 --global=0.0005 --verb
perl enrimo.pl --mats=momet,moo3 --range=0.0974,0.0001,0.9739 --inp=0p9739.enr
```

REQUIREMENTS

```
Perl 5
Text::CSV, Excel::Writer::XLSX, JSON, YAML
```

SEE ALSO

enrimo on GitHub

enrimo on Zenodo

enrimo in a paper: J. Phys. Commun. 3 (2019) 055015

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