1 ORSO standard for reflectometry data files

1.1 human-readable data file

The header should be formatted using YAML.

1.1.1 Structure of the header

```
The first line should state what the file is. E.g. #reflectivity data file orso file format 0.0
```

The header is structured into information on the

- creator ownership of the data file
- data source ownership and provenience of the raw data
- reduction software and reduction steps
- misc non-orso content
- data column description and units

using key words and structure as listed in the dictionary.

And finally a one-line column description referring to the data section of the type # 1 Qz 2 RQz 3 sRQz 4 sQz 5 ... or just # 1 2 3 4 5 ...

1.1.2 Dictionary of the key words used in the header

```
creator (required)
```

This section referes to the creation of this file, not the data.

name (required) NX_CHAR

Name of the person who created this file

affiliation (optional) NX_CHAR

Affiliation of the person who created this file

time (optional) NX_DATE_TIME

Date and time of the creation of this file

system (optional) NX_CHAR

Computer and user who created this file

data source (required)

This section deals with the source of the data used for generating this file.

```
origin (required)
```

This referes to the legal ownership of the raw data.

```
owner NX CHAR
```

Name of the owner of the raw data facility

Name of the facility where the measurement has been berformed.

experiment ID (required, if applicable) NX CHAR

The proposal number or experiment ID under which the data were cllected.

```
experiment date (optional) NX_DATE_TIME
```

Dates when the experiement was performed (the whole period rather than the individual measurement).

title Title of the experiment / the measurement campain.

```
experiment (required)
```

instrument Name and if applicable type of the instrument used.

```
probe (required)
```

Radiation used during the experiemnt. Either neutrons or x-rays.

polarisation (optional)

For neutrons the polarisation might be given as +1 for fully spin up polarised, -1 for fully spin down polarised and θ for unpolarised.

Partial polarisation can be expressed as

measurement (required)

How and parameters

```
scheme (optional)
```

Measurement scheme / geometry. This might be angle- of erergy dispersive, or both.

wavelength range (optional) NX FLOAT

Value and unit for angle dispersive scheme

Format <value> # <unit>

Value range and unit for wavelength dispersive scheme.

Format [<lower limit>, <upper limit>] # <unit>

angular range (optional) NX_FLOAT

Value and unit for wavelength dispersive scheme

Format <value> # <unit>

Value range and unit for angle dispersive scheme.

```
Format [<lower limit>, <upper limit>] # <unit>
sample (required)
   Description of the measured sample.
   name (required)
     A name uniquely identifying the sample
   description (optional)
     Nominal composition of the sample if known.
     Format suggestion following GenX nomenclature
     - amb: air
     - layer: {material: Ni, thickness: 100 nm}
      - subs: Si
links (optional)
   List of links to related data, publications, instruments and so
   on. Free format, e.g.
   related extensive file : fulldatafile.hdf
                             : orso2020.123456.789
                           : doi:10.1016/j.nima.2016.03.007
   instrument reference
```

reduction (required)

Information on the reduction steps performed to obtain the data below from the raw data set(s) listed here.

software (required)

Name and version of the software.

call (required)

Echo of the call of the software or soemthing similar which allows to reproduce the data content of this file.

comments (optional)

Plain text with comments about the data reduction. This allows to explain details of the reduction algorithm or what assumptions have been made.

corrections (optional)

List of reduction steps that have been performed. Probably with reference to a standadised procedure (orso repository) or to a publication.

binning (optional)

Description of the binning applied to the data.

several ranges require a repetition of the block.

```
Qz range [:0.01] # Aa^-1
              type linear
              delta Qz 0.001 \# Aa^-1
     input files (required)
          Data files used for creating the data below.
          references (required if applicable)
              List of files used for normalisation of the data.
              file File name
              created Date of creation (measurement?) of the raw file
                  Format YYYY/MM/DD:hh:mm:ss
          datafiles (required)
              List of files containing the raw data.
              file File name
              created Date of creation (measurement?) of the raw file
                  Format YYYY/MM/DD:hh:mm:ss
     data state (optional)
          key word like summary of the reduction steps
          Format ':
misc (optional)
```

Optional section to be used with non-orso-standard key words.

data (required)

Column description and data array containing the reduced data and related quantities.

The content of columns 1 to 4 is defined. Further columns may contain whatever the creator wants - as long as it is clearly stated what it is and what the units are.

```
column 1 (required)
```

Must be one of Qz, alpha_i or lambda.

Together with the unit, i.e. nm^-1, Aa^-1, deg, rad, nm or Aa.

column 2 (required)

Must be the reflectivity or intensity as a function column 1.

If applicable with unit.

column 3 (required)

Must be the uncertainty of the quantity in column 2.

This might be the standard deviation (sigma), FWHM, or the like.

Including appropriate units.

column 4 (optional, but defined if present)

If available the uncertainty of the quantity in column 1. This might be the standard deviation (sigma), FWHM, or the like. Including appropriate units.

 ${f column} \ {f 5} \ ({f optional})$

. . .