NeXus: a common data format for neutron, x-ray, and muon science

Release 3.1

http://nexusformat.org
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http://www.nexusformat.org/
1.1 NeXus Introduction

NeXus is an effort by an international group of scientists to define a common data exchange format for neutron, X-ray, and muon experiments. NeXus is built on top of the scientific data format HDF5 and adds domain-specific rules for organizing data within HDF5 files in addition to a dictionary of well-defined domain-specific field names. The NeXus data format has two purposes. First, NeXus defines a format that can serve as a container for all relevant data associated with a scientific instrument or beamline. This is a very important use case. Second, NeXus defines standards in the form of application definitions for the exchange of data between applications. NeXus provides structures for raw experimental data as well as for processed data.

In recent years, a community of scientists and computer programmers working in neutron and synchrotron facilities around the world came to the conclusion that a common data format would fulfill a valuable function in the scattering community. As instrumentation becomes more complex and data visualization becomes more challenging, individual scientists, or even institutions, find it difficult to keep up with new developments. A common data format makes it easier, both to exchange experimental results and to exchange ideas about how to analyze them. It promotes greater cooperation in software development and stimulates the design of more sophisticated visualization tools. Additional background information is given in the chapter titled "Brief history of NeXus."

This section is designed to give a brief introduction to NeXus, the data format and tools that have been developed in response to these needs. It explains what a modern data format such as NeXus is and how to write simple programs to read and write NeXus files.

The programmers who produce intermediate files for storing analyzed data should agree on simple interchange rules.

1.1.1 What is NeXus?

The NeXus data format has four components:

A set of design principles to help people understand what is in the data files.

A set of data storage objects (Base Class Definitions and Application Definitions) to allow the development of portable analysis software.
A set of subroutines, utilities, and examples to make it easy to read and write NeXus data files.

A scientific community to provide the scientific data, advice, and continued involvement with the NeXus standard. NeXus provides a forum for the scientific community to exchange ideas in data storage.

In addition, NeXus relies on a set of low-level file formats to actually store NeXus files on physical media. Each of these components are described in more detail in the Physical File format section.

The NeXus Application-Programmer Interface (NAPI), which provides the set of subroutines for reading and writing NeXus data files, is described briefly in NAPI: The NeXus Application Programming Interface. (Further details are provided in the NAPI chapter.)

The principles guiding the design and implementation of the NeXus standard are described in the NeXus Design chapter.

Base classes, which comprise the data storage objects used in NeXus data files, are detailed in the Base Class Definitions chapter.

Additionally, a brief list describing the set of NeXus Utilities available to browse, validate, translate, and visualise NeXus data files is provided in the NeXus Utilities chapter.

A Set of Design Principles

NeXus data files contain four types of entity: data groups, data fields, attributes, and links.

Data Groups Data groups are like folders that can contain a number of fields and/or other groups.

Data Fields Data fields can be scalar values or multidimensional arrays of a variety of sizes (1-byte, 2-byte, 4-byte, 8-byte) and types (characters, integers, floats). In HDF, fields are represented as HDF Scientific Data Sets (also known as SDS).

Data Attributes Extra information required to describe a particular group or field, such as the data units, can be stored as a data attribute.

Links Links are used to reference the plottable data from NXdata when the data is provided in other groups such as NXmonitor or NXdetector.

In fact, a NeXus file can be viewed as a computer file system. Just as files are stored in folders (or subdirectories) to make them easy to locate, so NeXus fields are stored in groups. The group hierarchy is designed to make it easy to navigate a NeXus file.

Example of a NeXus File

The following diagram shows an example of a NeXus data file represented as a tree structure.
Example of a NeXus Data File

Note that each field is identified by a name, such as `counts`, but each group is identified both by a name and, after a colon as a delimiter, the class type, e.g., `monitor:NXmonitor`). The class types, which all begin with `NX`, define the sort of fields that the group should contain, in this case, counts from a beamline monitor. The hierarchical design, with data items nested in groups, makes it easy to identify information if you are browsing through a file.

Important Classes

Here are some of the important classes found in nearly all NeXus files. A complete list can be found in the `NeXus Design` chapter.

**Note:** `NXentry` and `NXdata` are the only two classes necessary to store the minimum amount of information in a valid NeXus data file.

**NXentry Required:** The top level of any NeXus file contains one or more groups with the class `NXentry`. These contain all the data that is required to describe an experimental run or scan. Each `NXentry` typically contains a number of groups describing sample information (class `NXsample`), instrument details (class `NXinstrument`), and monitor counts (class `NXmonitor`).

**NXdata Required:** Each `NXentry` group contains one or more groups with class `NXdata`. These groups contain the experimental results in a self-contained way, i.e., it should be possible to generate a sensible plot of the data from the information contained in each `NXdata` group. That means it should contain the axis labels and titles as well as the data.

**NXsample** A `NXentry` group will often contain a group with class `NXsample`. This group contains information pertaining to the sample, such as its chemical composition, mass, and environment variables (temperature, pressure, magnetic field, etc.).

**NXinstrument** There might also be a group with class `NXinstrument`. This is designed to encapsulate all the instrumental information that might be relevant to a measurement, such as flight paths, collimation, chopper frequencies, etc.
Since an instrument can include several beamline components each defined by several parameters, the components are each specified by a separate group. This hides the complexity from generic file browsers, but makes the information available in an intuitively obvious way if it is required.

Simple Example

NeXus data files do not need to be complicated. In fact, the following diagram shows an extremely simple NeXus file (in fact, the simple example shows the minimum information necessary for a NeXus data file) that could be used to transfer data between programs. (Later in this section, we show how to write and read this simple example.)

Example structure of a simple data file

This illustrates the fact that the structure of NeXus files is extremely flexible. It can accommodate very complex instrumental information, if required, but it can also be used to store very simple data sets. Here is the structure of a very simple NeXus data file (examples/verysimple.nx5):

Structure of a very simple NeXus Data file

```plaintext
verysimple.nx5 : NeXus data file
file_name = verysimple.nx5
file_time = 2014-09-08T09:07:11.939912
NXus_version = 4.3.0
```
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NeXus files are easy to create. This example NeXus file was created using a short Python program and NeXpy:

```python
#!/usr/bin/env python

# This example uses NeXpy to build the verysimple.nx5 data file.

from nexpy.api import nexus

angle = [18.9094, 18.9096, 18.9098, 18.91, 18.9102,
         18.9104, 18.9106, 18.9108, 18.911, 18.9112,
         18.9114, 18.9116, 18.9118, 18.912, 18.9122]

diode = [1193, 4474, 53220, 274310, 515430, 827880,
         1227100, 1434640, 1330280, 1037070, 598720,
         316460, 56677, 1000, 1000]

two_theta = nexus.SDS(angle, name="two_theta",
                       units="degrees",
                       long_name="two_theta (degrees)")

counts = nexus.SDS(diode, name="counts",
                   long_name="photodiode counts")

data = nexus.NXdata(counts, [two_theta])
data.save("verysimple.nx5")
```

### A Set of Data Storage Objects

If the design principles are followed, it will be easy for anyone browsing a NeXus file to understand what it contains, without any prior information. However, if you are writing specialized visualization or analysis software, you will need to know precisely what specific information is contained in advance. For that reason, NeXus provides a way of defining the format for particular instrument types, such as time-of-flight small angle neutron scattering. This requires some agreement by the relevant communities, but enables the development of much more portable software.

The set of data storage objects is divided into three parts: base classes, application definitions, and contributed definitions. The base classes represent a set of components that define the dictionary of all possible terms to be used with that component. The application definitions specify the minimum required information to satisfy a particular scientific or data analysis software interest. The contributed definitions have been submitted by the scientific community for incubation before they are adopted by the NIAC or for availability to the community.
These instrument definitions are formalized as XML files, using NXDL, (as described in the NXDL chapter) to specify the names of data fields, and other NeXus data objects. The following is an example of such a file for the simple NeXus file shown above.

A very simple NeXus Definition Language (NXDL) file

```xml
<?xml version="1.0" ?>
<definition
   xmlns="http://definition.nexusformat.org/nxdl/3.1"
   xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance"
   xsi:schemaLocation="http://definition.nexusformat.org/nxdl/3.1 ../nxdl.xsd"
   category="base"
   name="verysimple"
   version="1.0"
   type="group" extends="NXobject">
   <doc>
      A very simple NeXus NXDL file
   </doc>
   <group type="NXentry">
      <group type="NXdata">
         <field name="counts" type="NX_INT" units="NX_UNITLESS">
            <doc>counts recorded by detector</doc>
         </field>
         <field name="two_theta" type="NX_FLOAT" units="NX_ANGLE">
            <doc>rotation angle of detector arm</doc>
         </field>
      </group>
   </group>
</definition>
```

Complete examples of reading and writing NeXus data files are provided later. This chapter has several examples of writing and reading NeXus data files. If you want to define the format of a particular type of NeXus file for your own use, e.g. as the standard output from a program, you are encouraged to publish the format using this XML format. An example of how to do this is shown in the Creating a NXDL Specification section.

A Set of Subroutines

NeXus data files are high-level so the user only needs to know how the data are referenced in the file but does not need to be concerned where the data are stored in the file. Thus, the data are most easily accessed using a subroutine library tuned to the specifics of the data format.

In the past, a data format was defined by a document describing the precise location of every item in the data file, either as row and column numbers in an ASCII file, or as record and byte numbers in a binary file. It is the job of the subroutine library to retrieve the data. This subroutine library is commonly called an application-programmer interface or API.

For example, in NeXus, a program to read in the wavelength of an experiment would contain lines similar to the following:
Simple example of reading data using the NeXus API

```c
NXopendata (fileID, "wavelength");
NXgetdata (fileID, lambda);
NXclosedata (fileID);
```

In this example, the program requests the value of the data that has the label `wavelength`, storing the result in the variable `lambda`. `fileID` is a file identifier that is provided by NeXus when the file is opened.

We shall provide a more complete example when we have discussed the contents of the NeXus files.

Scientific Community

NeXus began as a group of scientists with the goal of defining a common data storage format to exchange experimental results and to exchange ideas about how to analyze them.

The NeXus Community provides the scientific data, advice, and continued involvement with the NeXus standard. NeXus provides a forum for the scientific community to exchange ideas in data storage through the NeXus wiki.

The NeXus International Advisory Committee (NIAC) supervises the development and maintenance of the NeXus common data format for neutron, X-ray, and muon science. The NIAC supervises a technical committee to oversee the NAPI: NeXus Application Programmer Interface (frozen) and the Introduction to NeXus definitions.

Representation of data examples

Most of the examples of data files have been written in a format intended to show the structure of the file rather than the data content. In some cases, where it is useful, some of the data is shown. Consider this prototype example:

```plaintext
example of NeXus data file structure
```

```c
every
entry:NXentry
  instrument:NXinstrument
    detector:NXdetector
      data:[ ]
        axes = "bins"
        long_name = "strip detector 1-D array"
        signal = 1
        bins:[0, 1, 2, ... 1023]
        long_name = "bin index numbers"
      sample:NXsample
        name = "zeolite"
    data:NXdata
      data --> /entry/instrument/detector/data
      bins --> /entry/instrument/detector/bins
```

Some words on the notation:
- Hierarchy is represented by indentation. Objects on the same indentation level are in the same group.
- The combination `name:NXclass` denotes a NeXus group with name `name` and class `NXclass`.
- A simple name (no following class) denotes a data field. An equal sign is used to show the value, where this is important to the example.
• Sometimes, a data type is specified and possibly a set of dimensions. For example, `energy:NX_NUMBER[NE]` says `energy` is a 1-D array of numbers (either integer or floating point) of length `NE`.

• Attributes are noted as `@name="value"` pairs. The `@` symbol only indicates this is an attribute and is not part of the attribute name.

• Links are shown with a text arrow `-->` indicating the source of the link (using HDF5 notation listing the sequence of names).

Line 1 shows that there is one group at the root level of the file named `entry`. This group is of type `NXentry` which means it conforms to the specification of the `NXentry NeXus base class. Using the HDF5 nomenclature, we would refer to this as the `/entry` group.

Lines 2, 10, and 12: The `/entry` group contains three subgroups: `instrument`, `sample`, and `data`. These groups are of type `NXinstrument`, `NXsample`, and `NXdata`, respectively.

Line 4: The data of this example is stored in the `/entry/instrument/detector` group in the dataset called `data` (HDF5 path is `/entry/instrument/detector/data`). The indication of `data:\[\]` says that `data` is an array of unspecified dimension(s).

Lines 5-7: There are three attributes of `/entry/instrument/detector/data`: `axes`, `long_name`, and `signal`.

Line 8 (reading `bins:\[0, 1, 2, ... 1023\]`) shows that `bins` is a 1-D array of length presumably 1024. A small, representative selection of values are shown.

Line 9: an attribute that shows a descriptive name of `/entry/instrument/detector/bins`. This attribute might be used by a NeXus client while plotting the data.

Line 11 (reading `name = "zeolite"`) shows how a string value is represented.

Lines 13-14: The `/entry/data` group has two datasets that are actually linked as shown. (As you will see later, the `NXdata` group is required and enables NeXus clients to easily determine what to offer for display on a default plot.)

**Class path specification**

In some places in this documentation, a path may be shown using the class types rather than names. For example:

`/NXentry/NXinstrument/NXcrystal/wavelength`

identifies a dataset called `wavelength` that is inside a group of type `NXcrystal` ...

As it turns out, this syntax is the syntax used in NXDL `link` specifications. This syntax is also used when the exact name of each group is either unimportant or not specified.

If default names are taken for each class, then the above class path is expressed as this equivalent HDF5 path:

`/entry/instrument/crystal/wavelength`

In some places in this documentation, where clarity is needed to specify both the path and class name, you may find this equivalent path:

`/entry:NXentry/instrument:NXinstrument/crystal:NXcrystal/wavelength`

**Motivations for the NeXus standard in the Scientific Community**

By the early 1990s, several groups of scientists in the fields of neutron and X-ray science had recognized a common and troublesome pattern in the data acquired at various scientific instruments and user facilities. Each of these instruments
and facilities had a locally defined format for recording experimental data. With lots of different formats, much of the scientists’ time was being wasted in the task of writing import readers for processing and analysis programs. As is common, the exact information to be documented from each instrument in a data file evolves, such as the implementation of new high-throughput detectors. Many of these formats lacked the generality to extend to the new data to be stored, thus another new format was devised. In such environments, the documentation of each generation of data format is often lacking.

Three parallel developments have led to NeXus:

1. **June 1994**: Mark Könnecke (Paul Scherer Institute, Switzerland) made a proposal using netCDF for the European neutron scattering community while working at the ISIS pulsed neutron facility.

2. **August 1994**: Jon Tischler and Mitch Nelson (Oak Ridge National Laboratory, USA) proposed an HDF-based format as a standard for data storage at the Advanced Photon Source (Argonne National Laboratory, USA).

3. **October 1996**: Przemek Klosowski (National Institute of Standards and Technology, USA) produced a first draft of the NeXus proposal drawing on ideas from both sources.

These scientists proposed methods to store data using a self-describing, extensible format that was already in broad use in other scientific disciplines. Their proposals formed the basis for the current design of the NeXus standard which was developed across three workshops organized by Ray Osborn (ANL), SoftNeSS’94 (Argonne Oct. 1994), SoftNeSS’95 (NIST Sept. 1995), and SoftNeSS’96 (Argonne Oct. 1996), attended by representatives of a range of neutron and X-ray facilities. The NeXus API was released in late 1997. Basic motivations for this standard were:

1. **Simple plotting**

2. **Unified format for reduction and analysis**

3. **Defined dictionary of terms**

**Simple plotting** An important motivation for the design of NeXus was to simplify the creation of a default plot view. While the best representation of a set of observations will vary, depending on various conditions, a good suggestion is often known *a priori*. This suggestion is described in the `NXdata` element so that any program that is used to browse NeXus data files can provide a *best representation* without request for user input.

**Unified format for reduction and analysis** Another important motivation for NeXus, indeed the *raison d’etre*, was the community need to analyze data from different user facilities. A single data format that is in use at a variety of facilities would provide a major benefit to the scientific community. This should be capable of describing any type of data from the scientific experiments, at any step of the process from data acquisition to data reduction and analysis. This unified format also needs to allow data to be written to storage as efficiently as possible to enable use with high-speed data acquisition.

*Self-description*, combined with a reliance on a *multi-platform* (and thereby *portable*) data storage format, are valued components of a data storage format where the longevity of the data is expected to be longer than the lifetime of the facility at which it is acquired. As the name implies, self-description within data files is the practice where the structure of the information contained within the file is evident from the file itself. A multi-platform data storage format must faithfully represent the data identically on a variety of computer systems, regardless of the bit order or byte order or word size native to the computer.

The scientific community continues to grow the various types of data to be expressed in data files. This practice is expected to continue as part of the investigative process. To gain broad acceptance in the scientific user community, any data storage format proposed as a standard would need to be *extendable* and continue to provide a means to express the latest notions of scientific data.

The maintenance cost of common data structures meeting the motivations above (self-describing, portable, and extendable) is not insurmountable but is often well-beyond the research funding of individual members of the muon, neutron, and X-ray science communities. Since it is these members that drive the selection of a data storage format, it is necessary for the user cost to be as minimal as possible. In this case, experience has shown that the format must be

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in the public-domain for it to be commonly accepted as a standard. A benefit of the public-domain aspect is that the source code for the API is open and accessible, a point which has received notable comment in the scientific literature.

More recently, NeXus has recognized that part of the scientific community with a desire to write and record scientific data, has small data volumes and a large aversion to the requirement of a complicated API necessary to access data in binary files such as HDF. For such information, the NeXus API (NAPI) has been extended by the addition of the eXtensible Markup Language (XML)\(^1\) as an alternative to HDF. XML is a text-based format that supports compression and structured data and has broad usage in business and e-commerce. While possibly complicated, XML files are human readable, and tools for translation and extraction are plentiful. The API has routines to read and write XML data and to convert between HDF and XML.

**NeXus as a Common Data Exchange Format**  By the late 1980s, it had become common practice for a scientific instrument or facility to define its own data format, often at the convenience of the local computer system. Data from these facilities were not easily interchanged due to various differences in computer systems and the compression schemes of binary data. It was necessary to contact the facility to obtain a description so that one could write an import routine in software. Experience with facilities closing (and subsequent lack of access to information describing the facility data format) revealed a significant limitation with this common practice. Further, there existed a \(N \times N\) number of conversion routines necessary to convert data between various formats. In \(N\) separate file formats, circles represent different data file formats while arrows represent conversion routines. Note that the red circle only maps to one other format.

![Diagram of N separate file formats](image)

One early idea has been for NeXus to become the common data exchange format, and thereby reduce the number of data conversion routines from \(N \times N\) down to \(2N\), as shown in \(N\) separate file formats joined by a common NeXus converter.

**Defined dictionary of terms**  A necessary feature of a standard for the interchange of scientific data is a ‘defined dictionary (or lexicography) of terms. This dictionary declares the expected spelling and meaning of terms when they are present so that it is not necessary to search for all the variant forms of energy when it is used to describe data (e.g., \(E, e, keV, eV, nrg, \ldots\)).

NeXus recognized that each scientific specialty has developed a unique dictionary and needs to categorize data using those terms. The NeXus Application Definitions provide the means to document the lexicography for use in data files of that scientific specialty.

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\(^1\) XML: [http://www.w3.org/XML/](http://www.w3.org/XML/). There are many other descriptions of XML, for example: [http://en.wikipedia.org/wiki/XML](http://en.wikipedia.org/wiki/XML)
NAPI: The NeXus Application Programming Interface

The NeXus API consists of routines to read and write NeXus data files. It was written to provide a simple to use and consistent common interface for all supported backends (XML, HDF4 and HDF5) to scientific programmers and other users of the NeXus Data Standard.

Note: It is not necessary to use the NAPI to write or read NeXus data files. The intent of the NAPI is to simplify the programming effort to use the HDF programming interface. There are Examples of writing and reading NeXus data files to help you understand.

This section will provide a brief overview of the available functionality. Further documentation of the NeXus Application Programming Interface (NAPI) for bindings to specific programming language can be found in the NAPI chapter and may be downloaded from the NeXus development site. 2

For an even more detailed description of the internal workings of NAPI see NeXusIntern.pdf, copied from the NeXus code repository. That document is written for programmers who want to work on the NAPI itself. If you are new to NeXus and just want to implement basic file reading or writing you should not start by reading that.

How do I write a NeXus file? The NeXus Application Program Interface (NAPI) provides a set of subroutines that make it easy to read and write NeXus files. These subroutines are available in C, Fortran 77, Fortran 90, Java, Python, C++, and IDL.

The API uses a very simple state model to navigate through a NeXus file. (Compare this example with NAPI Simple 2-D Write Example (C, F77, F90), in the NAPI chapter, using the native HDF5 commands.) When you open a file, the API provides a file handle, which stores the current location, i.e. which group and/or field is currently open. Read and write operations then act on the currently open entity. Following the simple example titled Example structure of a simple data file, we walk through a schematic of NeXus program written in C (without any error checking or real data).

Writing a simple NeXus file using NAPI

```c
#include "napi.h"
```

int main()
{
    NXhandle fileID;
    NXopen ("NXfile.nxs", NXACC_CREATE, &fileID);
    NXmakegroup (fileID, "Scan", "NXEntry");
    NXopengroup (fileID, "Scan", "NXEntry");
    NXmakegroup (fileID, "data", "NXdata");
    NXopengroup (fileID, "data", "NXdata");
    /* somehow, we already have arrays tth and counts, each length n*/
    NXmakedata (fileID, "two_theta", NX_FLOAT32, 1, &n);
    NXopendata (fileID, "two_theta");
    NXputdata (fileID, tth);
    NXputattr (fileID, "units", "degrees", 7, NX_CHAR);
    NXclosedata (fileID); /* two_theta */
    NXmakedata (fileID, "counts", NX_FLOAT32, 1, &n);
    NXopendata (fileID, "counts");
    NXputdata (fileID, counts);
    NXclosedata (fileID); /* counts */
    NXclosegroup (fileID); /* data */
    NXclosegroup (fileID); /* Scan */
    NXclose (&fileID);
    return;
}

program analysis

1. line 6: Open the file NXfile.nxs with create access (implying write access). NAPI
   returns a file identifier of type NXhandle.

2. line 7: Next, we create the NXEntry group to contain the scan using NXmakegroup() and then open it for
   access using NXopengroup().

3. line 9: The plottable data is contained within an NXdata group, which must also be created and opened.

4. line 12: To create a field, call NXmakedata(), specifying the data name, type (NX_FLOAT32), rank (in this
   case, 1), and length of the array (n). Then, it can be opened for writing.

5. line 14: Write the data using NXputdata().

6. line 15: With the field still open, we can also add some data attributes, such as the data units, which
   are specified as a character string (type="NX_CHAR") that is 7 bytes long.

7. line 16: Then we close the field before opening another. In fact, the API will do this automatically if you
   attempt to open another field, but it is better style to close it yourself.

8. line 17: The remaining fields in this group are added in a similar fashion. Note that the indentation whenever a
   new field or group are opened is just intended to make the structure of the NeXus file more transparent.

9. line 20: Finally, close the groups (NXdata and NXentry) before closing the file itself.

How do I read a NeXus file? Reading a NeXus file works in the same way by traversing the tree with the handle.

3 NAPI: NeXus Application Programmer Interface (frozen)
4 See the chapter Base Class Definitions for more information.
5 The NeXus Data Types section describes the available data types, such as NX_FLOAT32 and NX_CHAR.
6 NeXus Data Units
7 The NeXus rule about data units is described in the NeXus Data Units section.
8 see Data Types allowed in NXDL specifications
This schematic C code will read the two-theta array created in the example above. (Again, compare this example with Reading a simple NeXus file using native HDF5 commands in C.)

Reading a simple NeXus file using NAPI

```
NXopen ("NXfile.nxs", NXACC_READ, &fileID);
NXopengroup (fileID, "Scan", "NXentry");
NXopengroup (fileID, "data", "NXdata");
NXopendata (fileID, "two_theta");
NXgetinfo (fileID, &rank, dims, &datatype);
NXmalloc ((void **) &tth, rank, dims, datatype);
NXgetdata (fileID, tth);
NXclosedata (fileID);
NXclosegroup (fileID);
NXclosegroup (fileID);
NXclose (fileID);
```

How do I browse a NeXus file? NeXus files can also be viewed by a command-line browser, nxbrowse, which is included as a helper tool in the NeXus API distribution. The following is an example session of nxbrowse to view a data file.

Using nxbrowse

```
%> nxbrowse lrcs3701.nxs
NXBrowse 3.0.0. Copyright (C) 2000 R. Osborn, M. Koennecke, P. Klosowski
NXus_version = 1.3.3
file_name = lrcs3701.nxs
file_time = 2001-02-11 00:02:35-0600
user = EAG/RO
NX> dir
NX Group : Histogram1 (NXentry)
NX Group : Histogram2 (NXentry)
NX> open Histogram1
NX/Histogram1> dir
NX Data : title[44] (NX_CHAR)
NX Data : analysis[7] (NX_CHAR)
NX Data : start_time[24] (NX_CHAR)
NX Data : end_time[24] (NX_CHAR)
NX Data : run_number (NX_INT32)
NX Group : sample (NXsample)
NX Group : LRMECS (NXinstrument)
NX Group : monitor1 (NXmonitor)
NX Group : monitor2 (NXmonitor)
NX Group : data (NXdata)
NX/Histogram1> open data
NX/Histogram1/data> dir
NX Data : title[44] (NX_CHAR)
NX Data : data[148,750] (NX_INT32)
```
program analysis

1. line 1: Start nxbrowse from the UNIX command line and open file lrcs3701.nxs from IPNS/LRMECS.
2. line 8: List the contents of the current group.
3. line 11: Open the NeXus group Histogram1.
4. line 23: Print the contents of the NeXus data labeled title.
5. line 41: Close the current group.
6. line 43: Quits nxbrowse.

The source code of nxbrowse provides an example of how to write a NeXus reader. The test programs included in the NeXus API may also be useful to study.

1.2 NeXus Design

This chapter actually defines the rules to use for writing valid NeXus files. An explanation of NeXus objects is followed by the definition of NeXus coordinate systems, the rules for structuring files and the rules for storing single items of data.

The structure of NeXus files is extremely flexible, allowing the storage both of simple data sets, such as a single data array and its axes, and also of highly complex data, such as the simulation results or an entire multi-component instrument. This flexibility is a necessity as NeXus strives to capture data from a wild variety of applications in X-ray, muSR and neutron scattering. The flexibility is achieved through a hierarchical structure, with related fields collected together into groups, making NeXus files easy to navigate, even without any documentation. NeXus files are self-describing, and should be easy to understand, at least by those familiar with the experimental technique.

Note: In this manual, we use the terms field, data field, and data item synonymously to be consistent with their meaning between NeXus data file instances and NXDL specification files.

9 https://svn.nexusformat.org/code/trunk/applications/NXbrowse/NXbrowse.c
1.2.1 NeXus Objects and Terms

Before discussing the design of NeXus in greater detail it is necessary to define the objects and terms used by NeXus. These are:

- **Data Groups**: Group data fields and other groups together. Groups represent levels in the NeXus hierarchy.
- **Data Fields**: Multidimensional arrays and scalars representing the actual data to be stored.
- **Data Attributes**: Additional metadata which can be assigned to groups or data fields.
- **Links**: Elements which point to data stored in another place in the file hierarchy.
- **NeXus Base Classes**: Dictionaries of names possible in the various types of NeXus groups.
- **NeXus Application Definitions**: Describe the minimum content of a NeXus file for a particular usage case.

In the following sections these elements of NeXus files will be defined in more detail.

**Data Groups**

NeXus files consist of data groups, which contain fields and/or other groups to form a hierarchical structure. This hierarchy is designed to make it easy to navigate a NeXus file by storing related fields together. Data groups are identified both by a name, which must be unique within a particular group, and a class. There can be multiple groups with the same class but they must have different names (based on the HDF rules).

For the class names used with NeXus data groups the prefix NX is reserved. Thus all NeXus class names start with NX.

**Data Fields**

Data fields contain the essential information stored in a NeXus file. They can be scalar values or multidimensional arrays of a variety of sizes (1-byte, 2-byte, 4-byte, 8-byte) and types (integers, floats, characters). The fields may store both experimental results (counts, detector angles, etc), and other information associated with the experiment (start and end times, user names, etc). Data fields are identified by their names, which must be unique within the group in which they are stored. Some fields have engineering units to be specified. In some cases, such in `NXdetector/data`, a field is expected to have be an array of several dimensions.

**Examples of data fields**

- `variable (NX_NUMBER)` Dimension scale defining an axis of the data.
- `variable_errors (NX_NUMBER)` Errors (uncertainties) associated with axis `variable`.
- `wavelength (NX_FLOAT)` Wavelength of radiation, `units="NX_FLOAT"`.
- `chemical_formula (NX_CHAR)` The chemical formula specified using CIF conventions.
- `name (NX_CHAR)` Name of user responsible for this entry.
- `data (NX_NUMBER)` Data values from the detector, `units="NX_ANY"`.

**Data Attributes**

Attributes are extra (meta-)information that are associated with particular fields. They are used to annotate the data, e.g. with physical units or calibration offsets, and may be scalar numbers or character strings. In addition, NeXus uses attributes to identify plottable data and their axes, etc. A description of some of the many possible attributes can be found in the next table:
Examples of data attributes

**units (NX_CHAR)**  Data units given as character strings, must conform to the NeXus units standard. See the NeXus Data Units section for details.

**signal (NX_POSINT)**  Defines which data set contains the signal to be plotted, use signal=1 for main signal, signal=2 for a second item to plot, and so on.

**axes (NX_CHAR)**  axes defines the names of the dimension scales for this data set as a colon-delimited list. Note that some legacy data files may use a comma as delimiter.

For example, suppose data is an array with elements data[j][i] (C) or data(i,j) (Fortran), with dimension scales time_of_flight[i] and polar_angle[j], then data would have an attribute axes="polar_angle:time_of_flight" in addition to an attribute signal=1.

**axis (NX_POSINT)**  The original way of designating data for plotting, now superceded by the axes attribute. This defines the rank of the signal data for which this data set is a dimension scale in order of the fastest varying index (see a longer discussion in the section on NXdata structure), i.e. if the array being stored is data, with elements data[j][i] in C and data(i,j) in Fortran, axis would have the following values: ith dimension (axis=1), jth dimension (axis=2), etc.

**primary (NX_POSINT)**  Defines the order of preference for dimension scales which apply to the same rank of signal data. Use primary=1 to indicate preferred dimension scale

**long_name (NX_CHAR)**  Defines title of signal data or axis label of dimension scale

**calibration_status (NX_CHAR)**  Defines status of data value - set to Nominal or Measured

**offset (NX_INT)**  Rank values of offsets to use for each dimension if the data is not in C storage order

**stride (NX_INT)**  Rank values of steps to use when incrementing the dimension

**vector (NX_FLOAT)**  3 values describing the axis of rotation or the direction of translation

**interpretation (NX_CHAR)**  Describes how to display the data. Allowed values include:

- scaler (0-D data)
- spectrum (1-D data)
- image (2-D data)
- vertex (3-D data)

Finally, NeXus files themselves have global attributes which are listed in the next table. These attributes identify the NeXus version, file creation time, etc. All attributes are identified by their names, which must be unique within each field.

Examples of global attributes

**file_name (NX_CHAR)**  File name of original NeXus file to assist in identification if the external name has been changed

**file_time (ISO 8601)**  Date and time of file creation

**file_update_time (ISO 8601)**  Date and time of last file change at close

**NeXus_version (NX_CHAR)**  Version of NeXus API used in writing the file

**creator (NX_CHAR)**  Facility or program where the file originated
Links

Links are pointers to existing data somewhere else. The concept is very much like symbolic links in a unix filesystem. The NeXus definition sometimes requires to have access to the same data in different groups in the same file. For example: detector data is stored in the `NXinstrument/NXdetector` group but may be needed in `NXdata` for automatic plotting. Rather then replicating the data, NeXus uses links in such situations. See the figure for a more descriptive representation of the concept of linking.

![Figure 1.3: Linking in a NeXus file](image)

NeXus also allows for links to external files. Consider the case where an instrument uses a detector with a closed-system software support provided by a commercial vendor. This system writes its images into a NeXus HDF5 file. The instrument's data acquisition system writes instrument metadata into another NeXus HDF5 file. In this case, the instrument metadata file might link to the data in the detector image file. Here is an example (from Diamond Light Source) showing an external file link in HDF5:

```hdf5
EXTERNAL_LINK "data" {
    TARGETFILE "/dls/i22/data/2012/sm7594-1/i22-69201-Pilatus2M.h5"
    TARGETPATH "entry/instrument/detector/data"
}
```

NeXus Base Classes

Data groups often describe objects in the experiment (monitors, detectors, monochromators, etc.), so that the contents (both data fields and/or other data groups) comprise the properties of that object. NeXus has defined a set of standard objects, or base classes, out of which a NeXus file can be constructed. This is each data group is identified by a name and a class. The group class, defines the type of object and the properties that it can contain, whereas the group name defines a unique instance of that class. These classes are defined in XML using the NeXus Definition Language (NXDL) format. All NeXus class types adopted by the NIAC must begin with NX. Classes not adopted by the NIAC must not start with NX.

Note: NeXus base classes are the components used to build the NeXus data structure.
Not all classes define physical objects. Some refer to logical groupings of experimental information, such as plottable data, sample environment logs, beam profiles, etc. There can be multiple instances of each class. On the other hand, a typical NeXus file will only contain a small subset of the possible classes.

**Note:** The groups, fields, links, and attributes of a base class definition are all optional, with a few particular exceptions in NXentry and NXdata. They are named in the specification to describe the exact spelling and usage of the term when it appears.

NeXus base classes are not proper classes in the same sense as used in object oriented programming languages. In fact the use of the term classes is actually misleading but has established itself during the development of NeXus. NeXus base classes are rather dictionaries of field names and their meanings which are permitted in a particular NeXus group implementing the NeXus class. This sounds complicated but becomes easy if you consider that most NeXus groups describe instrument components. Then for example, a NXmonochromator base class describes all the possible field names which NeXus allows to be used to describe a monochromator.

Most NeXus base classes represent instrument components. Some are used as containers to structure information in a file (NXentry, NXcollection, NXinstrument, NXprocess, NXparameter). But there are some base classes which have special uses which need to be mentioned here:

- **NXdata**  
  NXdata is used to identify the default plottable data. The notion of a default plot of data is a basic motivation of NeXus.

- **NXlog**  
  NXlog is used to store time stamped data like the log of a temperature controller. Basically you give a start time, and arrays with a difference in seconds to the start time and the values read.

- **NXnote**  
  This group provides a place to store general notes, images, video or whatever. A mime type is stored together with a binary blob of data. Please use this only for auxiliary information, for example an image of your sample, or a photo of your boss.

- **NXgeometry**  
  NXgeometry and its subgroups NXtranslation, NXorientation, NXshape are used to store absolute positions in the laboratory coordinate system or to define shapes.

These groups can appear anywhere in the NeXus hierarchy, where needed. Preferably close to the component they annotate or in a NXcollection. All of the base classes are documented in the reference manual.

**NXdata Facilitates Automatic Plotting**

The most notable special base class (or group in NeXus) is NXdata. NXdata is the answer to a basic motivation of NeXus to facilitate automatic plotting of data. NXdata is designed to contain the main dataset and its associated dimension scales (axes) of a NeXus data file. The usage scenario is that an automatic data plotting program just opens a NXentry and then continues to search for any NXdata groups. These NXdata groups represent the plottable data. An algorithm for identifying the default plottable data is presented in the chapter titled Rules for Storing Data Items in NeXus Files.

**Where to Store Metadata**

There are many ways to store metadata about your experiments. Already there are many fields in the various base classes to store the more common or general metadata, such as wavelength. (For wavelength, see the Strategies: The wavelength section.)

One common scheme is to store the metadata all in one group. If the group is to be validated for content, then there are several possibilities, as shown in the next table:
NeXus Application Definitions

The objects described so far provide us with the means to store data from a wide variety of instruments, simulations, or processed data as resulting from data analysis. But NeXus strives to express strict standards for certain applications of NeXus, too. The tool which NeXus uses for the expression of such strict standards is the NeXus Application Definition. A NeXus Application Definition describes which groups and data items have to be present in a file in order to properly describe an application of NeXus. For example for describing a powder diffraction experiment. Typically an application definition will contain only a small subset of the many groups and fields defined in NeXus. NeXus application definitions are also expressed in the NeXus Definition Language (NXDL). A tool exists which allows one to validate a NeXus file against a given application definition.

Note: NeXus application definitions define the minimum required information necessary to satisfy data analysis or other data processing.

Another way to look at a NeXus application definition is as a contract between a file producer (writer) and a file consumer (reader).

The contract reads: If you write your files following a particular NeXus application definition, I can process these files with my software.

Yet another way to look at a NeXus application definition is to understand it as an interface definition between data files and the software which uses this file. Much like an interface in the Java or other modern object oriented programming languages.

In contrast to NeXus base classes, NeXus supports inheritance in application definitions.

Please note that a NeXus Application Definition will only define the bare minimum of data necessary to perform common analysis with data. Practical files will nearly always contain more data. One of the beauties of NeXus is that it is always possible to add more data to a file without breaking its compliance with its application definition.

1.2.2 NeXus Coordinate Systems

The NeXus coordinate system is shown below. Note that it is the same as that used by McStas (http://mcstas.risoe.dk).

Note: The NeXus definition of +z is opposite to that in the IUCr International Tables for Crystallography, volume G, and consequently, +x is also reversed.

Coordinate systems in NeXus have undergone significant development. Initially, only motor positions of the relevant motors were stored without further standardization. This soon proved to be too little and the NeXus polar coordinate system was developed. This system still is very close to angles that are meaningful to an instrument scientist but allows to define general positions of components easily. Then users from the simulation community approached the NeXus team and asked for a means to store absolute coordinates. This was implemented through the use of the NXgeometry class on top of the McStas system. We soon learned that all the things we do can be expressed through the McStas
coordinate system. So it became the reference coordinate system for NeXus. NXgeometry was expanded to allow the description of shapes when the demand came up. Later, members of the CIF team convinced the NeXus team of the beauty of transformation matrices and NeXus was enhanced to store the necessary information to fully map CIF concepts. Not much had to be changed though as we choose to document the existing angles in CIF terms. The CIF system allows to store arbitrary operations and nevertheless calculate absolute coordinates in the laboratory coordinate system. It also allows to convert from local, for example detector coordinate systems, to absolute coordinates in the laboratory system.

Please note that NXgeometry and the polar coordinate system are suggested to be deprecated. For new projects, rather use the CIF approach.

**McStas and NXgeometry System**

As stated above, NeXus uses the McStas coordinate system (http://mcstas.risoe.dk) as its laboratory coordinate system. The instrument is given a global, absolute coordinate system where the z axis points in the direction of the incident beam, the x axis is perpendicular to the beam in the horizontal plane pointing left as seen from the source, and the y axis points upwards. See below for a drawing of the McStas coordinate system. The origin of this coordinate system is the sample position or, if this is ambiguous, the center of the sample holder with all angles and translations set to zero. The McStas coordinate system is illustrated in the next figure:

The NeXus NXgeometry class directly uses the McStas coordinate system. NXgeometry classes can appear in any component in order to specify its position. The suggested name to use is geometry. In NXgeometry the NXtranslation/values field defines the absolute position of the component in the McStas coordinate system. The NXorientation/value field describes the orientation of the component as a vector of in the McStas coordinate system.

Please note that it is planned to deprecate NXgeometry in favour of the transformation based system described below.
Simple (Spherical Polar) Coordinate System

In this system, the instrument is considered as a set of components through which the incident beam passes. The variable distance is assigned to each component and represents the effective beam flight path length between this component and the sample. A sign convention is used where negative numbers represent components pre-sample and positive numbers components post-sample. At each component there is local spherical coordinate system with the angles polar_angle and azimuthal_angle. The size of the sphere is the distance to the previous component.

In order to understand this spherical polar coordinate system it is helpful to look initially at the common condition that azimuthal_angle is zero. This corresponds to working directly in the horizontal scattering plane of the instrument. In this case polar_angle maps directly to the setting commonly known as two theta. Now, there are instruments where components live outside of the scattering plane. Most notably detectors. In order to describe such components we first apply the tilt out of the horizontal scattering plane as the azimuthal_angle. Then, in this tilted plane, we rotate to the component. The beauty of this is that polar_angle is always two theta. Which, in the case of a component out of the horizontal scattering plane, is not identical to the value read from the motor responsible for rotating the component. This situation is shown in Polar Coordinate System.

Please note that it is planned to deprecate this polar system in favour of the transformation based system described below.

Coordinate Transformations

Another way to look at coordinates is through the use of transformation matrices. In this world view, the absolute position of a component or a detector pixel with respect to the laboratory coordinate system is calculated by applying
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

a series of translations and rotations. These operations are commonly expressed as transformation matrices and their combination as matrix multiplication. A very important aspect is that the order of application of the individual operations does matter. Another important aspect is that any operation transforms the whole coordinate system and gives rise to a new local coordinate system. The mathematics behind this is well known and used in such applications such as industrial robot control, space flight and computer games. The beauty in this comes from the fact that the operations to apply map easily to instrument settings and constants. It is also easy to analyze the contribution of each individual operation: this can be studied under the condition that all other operations are at a zero setting.

In order to use coordinate transformations, several morsels of information need to be known:

- **Type** The type of operation: rotation or translation
- **Direction** The direction of the translation or the direction of the rotation axis
- **Value** The angle of rotation or the length of the translation
- **Order** The order of operations to apply to move a component into its place.

NeXus chooses to encode this information in the following way:

- **Type** Through a data set attribute `transformation_type`. This can take the value of either `translation` or `rotation`.
- **Direction** Through a data set attribute `vector`. This is a set of three values describing either the components of the rotation axis or the direction along which the translation happens.
- **Value** This is represented in the actual data of the data set. In addition, there is the `offset` attribute which has three components describing a translation to apply before applying the operation of the real axis. Without the offset attribute additional virtual translations would need to be introduced in order to encode mechanical offsets in the axis.
- **Order** The order is encoded through the `depends_on` attribute on a data set. The value of the `depends_on` attribute is the axis upon which the current axis sits. If the axis sits in the same group it is just a name, if it is in another group it is a path to the dependent axis. In addition, for each beamline component, there is a `depends_on` field which points to the data set at the head of the axis dependency chain. Take as an example an eulerian cradle as used on a four-circle diffractometer. Such a cradle has a dependency chain of `phi:chi:rotation_angle`. Then the `depends_on` field in `NXsample` would have the value `phi`.

**NeXus Transformation encoding**

Transformation encoding for an eulerian cradle on a four-circle diffractometer

```plaintext
sample:NXsample
    rotation_angle
        @transformation_type=rotation
        @vector=0,1,0
        @offset=0,0,0,
    chi
        @transformation_type=rotation
        @vector=0,0,1
        @offset=0,0,0,
        @depends_on=rotation_angle
    phi
        @transformation_type=rotation
        @vector=0,1,0
        @offset=0,0,0,
        @depends_on=chi
```
The type and direction of the NeXus standard operations is documented below in the table: *Actions of standard NeXus fields*. The rule is to always give the attributes to make perfectly clear how the axes work. The CIF scheme also allows to store and use arbitrarily named axes in a NeXus file.

**Actions of standard NeXus fields**

**Transformation Actions**

<table>
<thead>
<tr>
<th>Field Name</th>
<th>transformation_type</th>
<th>vector</th>
</tr>
</thead>
<tbody>
<tr>
<td>polar_angle</td>
<td>rotation</td>
<td>0 1 0</td>
</tr>
<tr>
<td>azimuthal_angle</td>
<td>rotation</td>
<td>0 0 1</td>
</tr>
<tr>
<td>meridional_angle</td>
<td>rotation</td>
<td>1 0 0</td>
</tr>
<tr>
<td>distance</td>
<td>translation</td>
<td>0 0 1</td>
</tr>
<tr>
<td>height</td>
<td>translation</td>
<td>0 1 0</td>
</tr>
<tr>
<td>x_translation</td>
<td>translation</td>
<td>1 0 0</td>
</tr>
<tr>
<td>chi</td>
<td>rotation</td>
<td>0 0 1</td>
</tr>
<tr>
<td>phi</td>
<td>rotation</td>
<td>0 1 0</td>
</tr>
</tbody>
</table>

For the NeXus spherical coordinate system, the order is implicit and is given in the next example.

**implicit order of NeXus spherical coordinate system**

azimuthal_angle:polar_angle:distance

This is also a nice example of the application of transformation matrices:

1. You first apply azimuthal_angle as a rotation around $z$. This rotates the whole coordinate out of the plane.
2. Then you apply polar_angle as a rotation around $y$ in the tilted coordinate system.
3. This also moves the direction of the $z$ vector. Along which you translate the component to place by distance.

### 1.2.3 Rules and Underlying File Formats

#### Rules for Structuring Information in NeXus Files

All NeXus files contain one or many groups of type NXentry at root level. Many files contain only one NXentry group, then the name is entry. The NXentry level of hierarchy is there to support the storage of multiple related experiments in one file. Or to allow the NeXus file to serve as a container for storing a whole scientific workflow from data acquisition to publication ready data. Also, NXentry class groups can contain raw data or processed data. For files with more than one NXentry group, since HDF requires that no two items at the same level in an HDF file may have the same name, the NeXus fashion is to assign names with an incrementing index appended, such as entry1, entry2, entry3, etc.

In order to illustrate what is written in the text, example hierarchies like the one in figure *Raw Data* are provided.

#### Content of a Raw Data NXentry Group

An example raw data hierarchy is shown in figure *Raw Data* (only showing the relevant parts of the data hierarchy). In the example shown, the data field in the NXdata group is linked to the 2-D detector data (a 512x512 array of 32-bit
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

... integers) which has the attribute signal = 1. Note that [ , ] represents a 2D array.

**NeXus Raw Data Hierarchy**

```plaintext
entry: NXentry
   instrument: NXinstrument
   source: NXsource
   ...,
   detector: NXdetector
   data: NX_INT32[512,512]
      signal = 1
   sample: NXsample
   control: NXmonitor
   data: NXdata
      data --> /entry/instrument/detector/data
```

An NXentry describing raw data contains at least a NXsample, one NXmonitor, one NXdata and a NXinstrument group. It is good practice to use the names sample for the NXsample group, control for the NXmonitor group holding the experiment controlling monitor and instrument for the NXinstrument group. The NXinstrument group contains further groups describing the individual components of the instrument as appropriate.

The NXdata group contains links to all those data items in the NXentry hierarchy which are required to put up a default plot of the data. As an example consider a SAXS instrument with a 2D detector. The NXdata will then hold a link to the detector image. If there is only one NXdata group, it is good practice to name it data. Otherwise, the name of the detector bank represented is a good selection.

**Content of a processed data NXentry group**

Processed data, see figure *Processed Data*, in this context means the results of a data reduction or data analysis program. Note that [ ] represents a 1D array.

**NeXus Processed Data Hierarchy**

```plaintext
entry: NXentry
   reduction: NXprocess
      program_name = "pyDataProc2010"
      version = "1.0a"
   input: NXparameter
      filename = "sn2013287.nxs"
   sample: NXsample
   data: NXdata
      data
         signal = 1
```

NeXus stores such data in a simplified NXentry structure. A processed data NXentry has at minimum a NXsample, a NXdata and a NXprocess group. Again the preferred name for the NXsample group is sample. In the case of processed data, the NXdata group holds the result of the processing together with the associated axis data. The NXprocess group holds the name and version of the program used for this processing step and further NXparameter groups. These groups ought to contain the parameters used for this data processing step in suitable detail so that the processing step can be reproduced.

Optionally a processed data NXentry can hold a NXinstrument group with further groups holding relevant information about the instrument. The preferred name is again instrument. Whereas for a raw data file, NeXus strives...
to capture as much data as possible, a \texttt{NXinstrument} group for processed data may contain a much-reduced subset.

\textbf{NXsubentry or Multi-Method Data}

Especially at synchrotron facilities, there are experiments which perform several different methods on the sample at the same time. For example, combine a powder diffraction experiment with XAS. This may happen in the same scan, so the data needs to be grouped together. A suitable \texttt{NXentry} would need to adhere to two different application definitions. This leads to name clashes which cannot be easily resolved. In order to solve this issue, the following scheme was implemented in NeXus:

- The complete beamline (all data) is stored in an appropriate hierarchy in an \texttt{NXentry}.
- The \texttt{NXentry} group contains further \texttt{NXsubentry} groups, one for each method. Each \texttt{NXsubentry} group is constructed like a \texttt{NXentry} group. It contains links to all those data items required to fulfill the application definition for the particular method it represents.

See figure \textit{NeXus Multi Method Hierarchy} for an example hierarchy. Note that \[,] represents a 2D array.

\textbf{Rules for Special Cases}

\textbf{Scans}  
Scans are difficult to capture because they have great variety. Basically, any variable can be scanned. Such behaviour cannot be captured in application definitions. Therefore NeXus solves this difficulty with a set of rules. In this section, \texttt{NP} is used as a symbol for the number of scan points.
• The scan dimension $NP$ is always the first dimension of any multi-dimensional dataset. The reason for this is that HDF allows the first dimension of a dataset to be unlimited. Which means, that data can be appended to the dataset during the scan.

• All data is stored as arrays of dimensions $NP$, original dimensions of the data at the appropriate position in the NXentry hierarchy.

• The NXdata group has to contain links to all variables varied during the scan and the detector data. Thus the NXdata group mimics the usual tabular representation of a scan.

• Datasets in an NXdata group must contain the proper attributes to enable the default plotting, as described in the section titled NXdata Facilitates Automatic Plotting.

Simple scan Examples may be in order here. Let us start with a simple case, the sample is rotated around its rotation axis and data is collected in a single point detector. See figure Simple Scan for an overview. Then we have:

• A dataset at NXentry/NXinstrument/NXdetector/data of length $NP$ containing the count data.

• A dataset at NXentry/NXsample/rotation_angle of length $NP$ containing the positions of rotation_angle at the various steps of the scan.

• NXdata contains links to:
  - NXentry/NXinstrument/NXdetector/data
  - NXentry/NXsample/rotation_angle

• All other data fields have their normal dimensions.

NeXus Simple Scan Example

```python
entry:NXentry
    instrument:NXinstrument
        detector:NXdetector
data[VP
signal = 1
sample:NXsample
    rotation_angle[VP
        axis=1
control:NXmonitor
data[N
    data --> /entry/instrument/detector/data
    rotation_angle --> /entry/sample/rotation_angle
```

Simple scan with area detector The next example is the same scan but with an area detector with $xsize$ times $ysize$ pixels. The only thing which changes is that /NXentry/NXinstrument/NXdetector/data will have the dimensions $NP$, $xsize$, $ysize$. See figure Simple Scan with Area Detector for an overview.

NeXus Simple Scan Example with Area Detector

```python
entry:NXentry
    instrument:NXinstrument
        detector:NXdetector
data[VP
        data --> /entry/instrument/detector/data
        rotation_angle --> /entry/sample/rotation_angle
```
Complex \textit{hkl} scan The next example involves a complex movement along an axis in reciprocal space which requires multiple motors of a four-circle diffractometer to be varied during the scan. We then have:

- A dataset at \texttt{NXentry/NXinstrument/NXdetector/data} of length \texttt{NP} containing the count data.
- A dataset at \texttt{NXentry/NXinstrument/NXdetector/polar\_angle} of length \texttt{NP} containing the positions of the detector’s polar\_angle at the various steps of the scan.
- A dataset at \texttt{NXentry/NXsample/rotation\_angle} of length \texttt{NP} containing the positions of rotation\_angle at the various steps of the scan.
- A dataset at \texttt{NXentry/NXsample/chi} of length \texttt{NP} containing the positions of chi at the various steps of the scan.
- A dataset at \texttt{NXentry/NXsample/phi} of length \texttt{NP} containing the positions of phi at the various steps of the scan.
- A dataset at \texttt{NXentry/NXsample/h} of length \texttt{NP} containing the positions of the reciprocal coordinate h at the various steps of the scan.
- A dataset at \texttt{NXentry/NXsample/k} of length \texttt{NP} containing the positions of the reciprocal coordinate k at the various steps of the scan.
- A dataset at \texttt{NXentry/NXsample/l} of length \texttt{NP} containing the positions of the reciprocal coordinate l at the various steps of the scan.

- \texttt{NXdata} contains links to:
  - \texttt{NXentry/NXinstrument/NXdetector/data}
  - \texttt{NXentry/NXinstrument/NXdetector/polar\_angle}
  - \texttt{NXentry/NXsample/rotation\_angle}
  - \texttt{NXentry/NXsample/chi}
  - \texttt{NXentry/NXsample/phi}
  - \texttt{NXentry/NXsample/h}
  - \texttt{NXentry/NXsample/k}
  - \texttt{NXentry/NXsample/l}

The datasets in \texttt{NXdata} must have the appropriate attributes as described in the axis location section.

- All other data fields have their normal dimensions.
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

NeXus Complex hkl Scan

```plaintext
entry:NXentry
  instrument:NXinstrument
    detector:NXdetector
      data[NE][NT]
      signal = 1
      polar_angle[NE][NT]
      axis = 1
    name
  sample:NXsample
    name
    rotation_angle[NE][NT]
      axis=1
    chi[NE][NT]
      axis=1
    phi[NE][NT]
      axis=1
    h[NE][NT]
      axis=1
      primary=1
    k[NE][NT]
      axis=1
    l[NE][NT]
      axis=1
  control:NXmonitor
    data[NE][NT]
      data --> /entry/instrument/detector/data
      rotation_angle --> /entry/sample/rotation_angle
      chi --> /entry/sample/chi
      phi --> /entry/sample/phi
      polar_angle --> /entry/instrument/detector/polar_angle
      h --> /entry/sample/h
      k --> /entry/sample/k
      l --> /entry/sample/l
```

Multi-parameter scan: XAS  
Data can be stored almost anywhere in the NeXus tree. While the previous examples showed data arrays in either NXdetector or NXsample, this example demonstrates that data can be stored in other places. Links are used to reference the data.

The example is for X-ray Absorption Spectroscopy (XAS) data where the monochromator energy is step-scanned and counts are read back from detectors before ($I_0$) and after ($I$) the sample. These energy scans are repeated at a sequence of sample temperatures to map out, for example, a phase transition. While it is customary in XAS to plot $\log(I_0/I)$, we show them separately here in two different NXdata groups to demonstrate that such things are possible. Note that the length of the 1-D energy array is $NE$ while the length of the 1-D temperature array is $NT$.

NeXus Multi-parameter scan: XAS

```plaintext
entry:NXentry
  instrument:NXinstrument
    detector:NXdetector
      data: NX_NUMBER[NE,NT]
      signal = 1
      axes = "energy:temperature"
```
Rastering  Rastering is the process of making experiments at various locations in the sample volume. Again, rasterisation experiments can be variable. Some people even raster on spirals! Rasterisation experiments are treated the same way as described above for scans. Just replace \texttt{NP} with \texttt{P}, the number of raster points.

Special rules apply if a rasterisation happens on a regular grid of size \(xraster, yraster\). Then the variables varied in the rasterisation will be of dimensions \(xraster, yraster\) and the detector data of dimensions \(xraster, yraster, \text{(original dimensions)}\) of the detector. For example, an area detector of size \(xsize, ysize\) then it is stored with dimensions \(xraster, yraster, xsize, ysize\).

\textbf{Warning:} Be warned: if you use the 2D rasterisation method with \(xraster, yraster\) you may end up with invalid data if the scan is aborted prematurely. This cannot happen if the first method is used.

\textbf{NXcollection}  On demand from the community, NeXus introduced a more informal method of storing information in a NeXus file. This is the \texttt{NXcollection} class which can appear anywhere underneath \texttt{NXentry}. \texttt{NXcollection} is a container for holding other data. The foreseen use is to document collections of similar data which do not otherwise fit easily into the \texttt{NXinstrument} or \texttt{NXsample} hierarchy, such as the intent to record all motor positions on a synchrotron beamline. Thus, \texttt{NXcollection} serves as a quick point of access to data for an instrument scientist or another expert. \texttt{NXcollection} is also a feature for those who are too lazy to build up the complete NeXus hierarchy. An example usage case is documented in figure \textit{NXcollection example}.

\textbf{NXcollection Example}

```python
entry:NXentry
  positioners:NXcollection
    mxx:NXpositioner
    mzz:NXpositioner
    sgu:NXpositioner
    ttv:NXpositioner
    hugo:NXpositioner
    ....
  scalars:NXcollection
title (dataset)
```

1.2. NeXus Design
Rules for Storing Data Items in NeXus Files

This section describes the rules which apply for storing single data fields in data files.

Naming Conventions

Group and field Names used within NeXus follow a naming convention described by the following rules:

- The names of NeXus group and field items must only contain a restricted set of characters. This set may be described by a regular expression syntax regular expression, as described below.
- For the class names of NeXus group items, the prefix NX is reserved. Thus all NeXus class names start with NX. The chapter titled NeXus: Reference Documentation lists the available NeXus class names as either base classes, application definitions, or contributed definitions.

Regular expression pattern for NXDL group and field names

It is recommended that all group and field names contain only these characters:

- lower case letters
- digits
- “_” (underscore character)

and that they begin with a lower case letter. This is the regular expression used to check this recommendation.

\[^[a-z_][a-z_\d_]+\]

The length should be limited to no more than 63 characters (imposed by the HDF5 rules for names).

It is recognized that some facilities will construct group and field names with upper case letters. NeXus data files with upper case characters in the group or field names might not be accepted by all software that reads NeXus data files. Hence, group and field names that do not pass the regular expression above but pass this expression:

\[^[A-Za-z_][\w_]+\]

will be flagged as a warning during data file validation.

Use of underscore in descriptive names

Sometimes it is necessary to combine words in order to build a descriptive name for a data field or a group. In such cases lowercase words are connected by underscores.

---

10 The class name is the value assigned to the NX_class attribute of an HDF5 group in the NeXus data file. This class name is different than the name of the HDF5 group. This is important when not using the NAPI to either read or write the HDF5 data file.
For all data fields, only names from the NeXus base class dictionaries should be used. If a data field name or even a complete component is missing, please suggest the addition to the *NIAC: The NeXus International Advisory Committee.* The addition will usually be accepted provided it is not a duplication of an existing field and adequately documented.

**Note:** The NeXus base classes provide a comprehensive dictionary of terms that can be used for each class. The expected spelling and definition of each term is specified in the base classes. It is not required to provide all the terms specified in a base class. Terms with other names are permitted but might not be recognized by standard software. Rather than persist in using names not specified in the standard, please suggest additions to the *NIAC: The NeXus International Advisory Committee.*

### NeXus Array Storage Order

NeXus stores multi-dimensional arrays of physical values in C language storage order, where the last dimension is the fastest varying. This is the rule. *Good reasons are required to deviate from this rule.*

It is possible to store data in storage orders other than C language order.

As well it is possible to specify that the data needs to be converted first before being useful. Consider one situation, when data must be streamed to disk as fast as possible and conversion to C language storage order causes unnecessary latency. This case presents a good reason to make an exception to the standard rule.

**Non C Storage Order** In order to indicate that the storage order is different from C storage order two additional data set attributes, offset and stride, have to be stored which together define the storage layout of the data. Offset and stride contain rank numbers according to the rank of the multidimensional data set. Offset describes the step to make when the dimension is multiplied by 1. Stride defines the step to make when incrementing the dimension. This is best explained by some examples.

**Offset and Stride for 1 D data:**

```plaintext
* raw data = 0 1 2 3 4 5 6 7 8 9
size[1] = { 10 }  // assume uniform overall array dimensions

* default stride:
  stride[1] = { 1 }
  offset[1] = { 0 }
  for i:
    result[i]:
      0 1 2 3 4 5 6 7 8 9

* reverse stride:
  stride[1] = { -1 }
  offset[1] = { 9 }
  for i:
    result[i]:
      9 8 7 6 5 4 3 2 1 0
```
Offset and Stride for 2D Data

```plaintext
* raw data = 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19
size[2] = { 4, 5 } // assume uniform overall array dimensions

* row major (C) stride:
  stride[2] = { 5, 1 }
  offset[2] = { 0, 0 }
  for i:
    for j:
      result[i][j]:
        0 1 2 3 4
        5 6 7 8 9
        10 11 12 13 14
        15 16 17 18 19

* column major (Fortran) stride:
  stride[2] = { 1, 4 }
  offset[2] = { 0, 0 }
  for i:
    for j:
      result[i][j]:
        0 4 8 12 16
        1 5 9 13 17
        2 6 10 14 18
        3 7 11 15 19

* "crazy reverse" row major (C) stride:
  stride[2] = { -5, -1 }
  offset[2] = { 4, 5 }
  for i:
    for j:
      result[i][j]:
        19 18 17 16 15
        14 13 12 11 10
        9 8 7 6 5
        4 3 2 1 0
```

Offset and Stride for 3D Data

```plaintext
* raw data = 0 1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19
  20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 39
  40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57 58 59
size[3] = { 3, 4, 5 } // assume uniform overall array dimensions

* row major (C) stride:
  stride[3] = { 20, 5, 1 }
  offset[3] = { 0, 0, 0 }
  for i:
    for j:
      for k:
        result[i][j][k]:
          0 1 2 3 4
          5 6 7 8 9
          10 11 12 13 14
          15 16 17 18 19
```

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* column major (Fortran) stride:
  
  \[
  \text{stride}[3] = (1, 3, 12) \\
  \text{offset}[3] = (0, 0, 0)
  \]

For\[i:]

\[
\text{for } j:\n\text{for } k:\n\text{result}[i][j][k] =
\]

\[
0 12 24 36 48 \\
3 15 27 39 51 \\
6 18 30 42 54 \\
9 21 33 45 57 \\
1 13 25 37 49 \\
4 16 28 40 52 \\
7 19 31 43 55 \\
10 22 34 46 58 \\
2 14 26 38 50 \\
5 17 29 41 53 \\
8 20 32 44 56 \\
11 23 35 47 59
\]

**NeXus Data Types**

<table>
<thead>
<tr>
<th>description</th>
<th>matching regular expression</th>
</tr>
</thead>
<tbody>
<tr>
<td>integer</td>
<td>NX_INT(</td>
</tr>
<tr>
<td>floating-point</td>
<td>NX_FLOAT(32</td>
</tr>
<tr>
<td>array</td>
<td>([{0-9}}]?</td>
</tr>
<tr>
<td>valid item name</td>
<td>^[A-Za-z-z_]{A-Za-z-a-z0-9_}+$</td>
</tr>
<tr>
<td>valid class name</td>
<td>^NX[A-Za-z-a-z0-9_]+$</td>
</tr>
</tbody>
</table>

NeXus supports numeric data as either integer or floating-point numbers. A number follows that indicates the number of bits in the word. The table above shows the regular expressions that matches the data type specifier.

**integers**  NX_INT8, NX_INT16, NX_INT32, or NX_INT64

**floating-point numbers**  NX_FLOAT32 or NX_FLOAT64

**date / time stamps**  NX_DATE_TIME or ISO8601: Dates and times are specified using ISO-8601 standard definitions. Refer to *NeXus dates and times*.

**strings**  All strings are to be encoded in UTF-8. Since most strings in a NeXus file are restricted to a small set of characters and the first 128 characters are standard across encodings, the encoding of most of the strings in a NeXus file will be a moot point. Where encoding in UTF-8 will be important is when recording peoples names in NXuser and text notes in NXnotes. Because the few places where encoding is important also have unpredictable content, as well as the way in which current operating systems handle character encoding,
it is practically impossible to test the encoding used. Hence, `nxvalidate` provides no messages relating to character encoding.

**Binary data** Binary data is to be written as `UINT8`.

**Images** Binary image data is to be written using `UINT8`, the same as binary data, but with an accompanying image mime-type. If the data is text, the line terminator is `[CR][LF]`.

**NeXus dates and times** NeXus dates and times should be stored using the ISO 8601 \(^\text{11}\) format, e.g. `1996-07-31T21:15:22+0600`. The standard also allows for time intervals in fractional seconds with \(1\) or more digits of precision. This avoids confusion, e.g. between U.S. and European conventions, and is appropriate for machine sorting.

**`strftime()` format specifiers for ISO-8601 time**

```
%Y-%m-%dT%H:%M:%S%z
```

**Note:** Note that the `T` appears literally in the string, to indicate the beginning of the time element, as specified in ISO 8601. It is common to use a space in place of the `T`, such as `1996-07-31 21:15:22+0600`. While human-readable (and later allowed in a relaxed revision of the standard), compatibility with libraries supporting the ISO 8601 standard is not assured with this substitution. The `strftime()` format specifier for this is “%Y-%m-%d %H:%M:%S%z”.

**NeXus Data Units**

Given the plethora of possible applications of NeXus, it is difficult to define units to use. Therefore, the general rule is that you are free to store data in any unit you find fit. However, any data field must have a units attribute which describes the units. Wherever possible, SI units are preferred. NeXus units are written as a string attribute (`NX_CHAR`) and describe the engineering units. The string should be appropriate for the value. Values for the NeXus units must be specified in a format compatible with Unidata UDunits \(^\text{12}\). Application definitions may specify units to be used for fields using an enumeration.

**Linking Multi Dimensional Data with Axis Data**

NeXus allows to store multi dimensional arrays of data. In most cases it is not sufficient to just have the indices into the array as a label for the dimensions of the data. Usually the information which physical value corresponds to an index into a dimension of the multi dimensional data set. To this purpose a means is needed to locate appropriate data arrays which describe what each dimension of a multi dimensional data set actually corresponds too. There is a standard HDF facility to do this: it is called dimension scales. Unfortunately, at a time, there was only one global namespace for dimension scales. Thus NeXus had to come up with its own scheme for locating axis data which is described here. A side effect of the NeXus scheme is that it is possible to have multiple mappings of a given dimension to physical data. For example a TOF data set can have the TOF dimension as raw TOF or as energy.

There are two methods of linking each data dimension to its respective dimension scale. The preferred method uses the `axes` attribute to specify the names of each dimension scale. The original method uses the `axis` attribute to identify with an integer the axis whose value is the number of the dimension. After describing each of these methods, the two methods will be compared. A prerequisite for both methods is that the data fields describing the axis are stored

\(^{11}\) ISO 8601: http://www.w3.org/TR/NOTE-datetime

\(^{12}\) The UDunits specification also includes instructions for derived units. At present, the contents of NeXus `units` attributes are not validated in data files.
together with the multi dimensional data set whose axes need to be defined in the same NeXus group. If this leads to data duplication, use links.

**Linking by name using the axes attribute** The preferred method is to define an attribute of the data itself called *axes*. The axes attribute contains the names of each dimension scale as a colon (or comma) separated list in the order they appear in C. For example:

```
data: NXdata
  time_of_flight = 1500.0 1502.0 1504.0 ...
  polar_angle = 15.0 15.6 16.2 ...
  some_other_angle = 0.0 0.0 2.0 ...
  data = 5 7 14 ...
  @axes = polar_angle:time_of_flight
  @signal = 1
```

**Linking by dimension number using the axis attribute** The original method is to define an attribute of each dimension scale called *axis*. It is an integer whose value is the number of the dimension, in order of fastest varying dimension. That is, if the array being stored is data with elements *data[j][i]* in C and *data(i,j)* in Fortran, where *i* is the time-of-flight index and *j* is the polar angle index, the NXdata group would contain:

```
data: NXdata
  time_of_flight = 1500.0 1502.0 1504.0 ...
  @axis = 1
  @primary = 1
  polar_angle = 15.0 15.6 16.2 ...
  @axis = 2
  @primary = 1
  some_other_angle = 0.0 0.0 2.0 ...
  @axis = 1
  data = 5 7 14 ...
  @signal = 1
```

The axis attribute must be defined for each dimension scale. The primary attribute is unique to this method of linking.

There are limited circumstances in which more than one dimension scale for the same data dimension can be included in the same NXdata group. The most common is when the dimension scales are the three components of an (hkl) scan. In order to handle this case, we have defined another attribute of type integer called primary whose value determines the order in which the scale is expected to be chosen for plotting, i.e.

- 1st choice: primary=1
- 2nd choice: primary=2
- etc.

If there is more than one scale with the same value of the axis attribute, one of them must have set primary=1. Defining the primary attribute for the other scales is optional.
Note: The primary attribute can only be used with the first method of defining dimension scales discussed above. In addition to the signal data, this group could contain a data set of the same rank and dimensions called errors containing the standard deviations of the data.

Discussion of the two linking methods In general the method using the axes attribute on the multi dimensional data set should be preferred. This leaves the actual axis describing data sets unannotated and allows them to be used as an axis for other multi dimensional data. This is especially a concern as an axis describing a data set may be linked into another group where it may describe a completely different dimension of another data set.

Only when alternative axes definitions are needed, the axis method should be used to specify an axis of a data set. This is shown in the example above for the some_other_angle field where axis=1 denotes another possible primary axis for plotting. The default axis for plotting carries the primary=1 attribute.

Both methods of linking data axes will be supported in NeXus utilities that identify dimension scales, such as NXUfindaxis().

Storing Detectors

There are very different types of detectors out there. Storing their data can be a challenge. As a general guide line: if the detector has some well defined form, this should be reflected in the data file. A linear detector becomes a linear array, a rectangular detector becomes an array of size xsize times ysize. Some detectors are so irregular that this does not work. Then the detector data is stored as a linear array, with the index being detector number till ndet. Such detectors must be accompanied by further arrays of length ndet which give azimuthal_angle, polar_angle and distance for each detector.

If data from a time of flight (TOF) instrument must be described, then the TOF dimension becomes the last dimension, for example an area detector of xsize vs. ysize is stored with TOF as an array with dimensions xsize, ysize, ntof.

Monitors are Special

Monitors, detectors that measure the properties of the experimental probe rather than the sample, have a special place in NeXus files. Monitors are crucial to normalize data. To emphasize their role, monitors are not stored in the NXinstrument hierarchy but on NXentry level in their own groups as there might be multiple monitors. Of special importance is the monitor in a group called control. This is the main monitor against which the data has to be normalized. This group also contains the counting control information, i.e. counting mode, times, etc.

Monitor data may be multidimensional. Good examples are scan monitors where a monitor value per scan point is expected or time-of-flight monitors.

Find the plottable data

Any program whose aim is to identify the default plottable data should use the following procedure:

1. Start at the top level of the NeXus data file.
2. Loop through the groups with class NXentry until the next step succeeds.
3. Open the NXentry group and loop through the subgroups with class NXdata until the next step succeeds.
4. Open the NXdata group and loop through the fields for the one field with attribute signal="1". Note: There should be only one field that matches.
This is the default plottable data.

(a) If this field has an attribute `axes`:
   
i. The `axes` attribute value contains a colon (or comma) delimited list (in the C-order of the data array) with the names of the dimension scales associated with the plottable data. Such as:
   ```
   axes="polar_angle:time_of_flight"
   ```
   ii. Parse `axes` and open the datasets to describe your dimension scales

(b) If this field has no attribute `axes`:
   
i. Search for datasets with attributes `axis=1`, `axis=2`, etc.
   
ii. These are the fields describing your axis. There may be several fields for any axis, i.e. there may be multiple fields with the attribute `axis=1`. Among them the field with the attribute `primary=1` is the preferred one. All others are alternative dimension scales.

5. Having found the default plottable data and its dimension scales: make the plot.

Consult the `NeXus API` section, which describes the routines available to program these operations. In the course of time, generic NeXus browsers will provide this functionality automatically.

**Physical File format**

This section describes how NeXus structures are mapped to features of the underlying physical file format. This is a guide for people who wish to create NeXus files without using the NeXus-API.

**Choice of HDF as Underlying File Format**

At its beginnings, the founders of NeXus identified the Hierarchical Data Format (HDF) as a capable and efficient multi-platform data storage format. HDF was designed for large data sets and already had a substantial user community. HDF was developed and maintained initially by the National Center for Supercomputing Applications (NCSA) at the University of Illinois at Urbana-Champaign (UIUC) and later spun off into its own group called The HDF Group (THG: http://www.hdfgroup.org/). Rather then developing its own unique physical file format, the NeXus group choose to build NeXus on top of HDF.

HDF (now HDF5) is provided with software to read and write data (this is the application-programmer interface, or API) using a large number of computing systems in common use for neutron and X-ray science. HDF is a binary data file format that supports compression and structured data.

**Mapping NeXus into HDF**

NeXus data structures map directly to HDF structures. NeXus `groups` are HDF4 `vgroups` or HDF5 `groups`, NeXus data sets (or `fields`) are HDF4 `SDS` (scientific data sets) or HDF5 `datasets`. Attributes map directly to HDF group or dataset attributes.

The only special case is the NeXus class name. HDF4 supports a group class which is set with the `Vsetclass()` call and read with `VGetclass()`. HDF-5 has no group class. Thus the NeXus class is stored as an attribute to the HDF-5 group with the name `NX_class` and value of the NeXus class name.

A NeXus `link` directly maps to the HDF linking mechanisms.

**Note:** Examples are provided in the `Examples of writing and reading NeXus data files` chapter. These examples include software to write and read NeXus data files using the NAPI, as well as other software examples that use native (non-NAPI) libraries. In some cases the examples show the content of the NeXus data files that are produced. Here are links to some of the examples:
Perhaps the easiest way to view the implementation of NeXus in HDF5 is to view how the data structures look. For this, we use the h5dump command-line utility provided with the HDF5 support libraries. Short examples are provided for the basic NeXus data components:

- **group**: created in C NAPI by:
  
  ```c
  NXmakegroup (fileID, "entry", "NXentry");
  ```

- **field**: created in C NAPI by:
  
  ```c
  NXmakedata (fileID, "two_theta", NX_FLOAT32, 1, &n);
  NXopendata (fileID, "two_theta");
  NXputdata (fileID, tth);
  ```

- **attribute**: created in C NAPI by:
  
  ```c
  NXputattr (fileID, "units", "degrees", 7, NX_CHAR);
  ```

- **link**: created in C NAPI by:
  
  ```c
  # --tba--
  # TODO: write some text about HDF5 hard links
  # until then, see the h5dump example below
  ```

See the sections *NAPI Simple 2-D Write Example (C, F77, F90)* and *NAPI Python Simple 3-D Write Example* in the *Examples of writing and reading NeXus data files* chapter for examples that use the native HDF5 calls to write NeXus data files.

**h5dump of a NeXus NXentry group**

```bash
GROUP "entry" {
  ATTRIBUTE "NX_class" {
    DATATYPE H5T_STRING {
      STRSIZE 7;
      STRPAD H5T_STR_NULLPAD;
      CSET H5T_CSET_ASCII;
      CTYPE H5T_C_S1;
    }
    DATASPACE SCALAR
    DATA {
      (0): "NXentry"
    }
  }
  ... group contents
}
```
**h5dump of a NeXus field (HDF5 dataset)**

```plaintext
DATASET "two_theta" {
  DATATYPE H5T_IEEE_F64LE
  DATASPACE SIMPLE { ( 31 ) / ( 31 ) }
  DATA {
    (0): 17.9261, 17.9259, 17.9258, 17.9256, 17.9254, 17.9252,
    (6): 17.9251, 17.9249, 17.9247, 17.9246, 17.9244, 17.9243,
    (12): 17.9241, 17.9239, 17.9237, 17.9236, 17.9234, 17.9232,
    (18): 17.9231, 17.9229, 17.9228, 17.9226, 17.9224, 17.9222,
    (24): 17.9221, 17.9219, 17.9217, 17.9216, 17.9214, 17.9213,
    (30): 17.9211
  }
  ATTRIBUTE "units" {
    DATATYPE H5T_STRING {
      STRSIZE 7;
      STRPAD H5T_STR_NULLPAD;
      CSET H5T_CSET_ASCII;
      CTYPE H5T_C_S1;
    }
    DATASPACE SCALAR
    DATA {
      (0): "degrees"
    }
  }
} # ... other attributes
```

**h5dump of a NeXus attribute**

```plaintext
ATTRIBUTE "axes" {
  DATATYPE H5T_STRING {
    STRSIZE 9;
    STRPAD H5T_STR_NULLPAD;
    CSET H5T_CSET_ASCII;
    CTYPE H5T_C_S1;
  }
  DATASPACE SCALAR
  DATA {
    (0): "two_theta"
  }
}
```

**h5dump of a NeXus link**

```plaintext
# NeXus links have two parts in HDF5 files.
# The dataset is created in some group.
# A "target" attribute is added to indicate the HDF5 path to this dataset.
ATTRIBUTE "target" {
  DATATYPE H5T_STRING {
    STRSIZE 21;
    STRPAD H5T_STR_NULLPAD;
  }
  DATASPACE SCALAR
  DATA {
  }
}
```

1.2. NeXus Design
# NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

```c
CSET H5T_CSET_ASCII;
CTYPE H5T_C_S1;

) 
DATASPACE SCALAR 
DATA {
(0): "entry/data/two_theta"
}
}
#
# then, the hard link is created that refers to the original dataset
# (Since the name is "two_theta" in this example, it is understood that
# this link is created in a different HDF5 group than "/entry/data").

DATASET "two_theta" {
HARDLINK "/entry/data/two_theta"

}
```

## Mapping NeXus into XML

This takes a bit more work than HDF. At the root of NeXus XML file is a XML element with the name `NXroot`. Further XML attributes to `NXroot` define the NeXus file level attributes. An example NeXus XML data file is provided in the *NeXus Introduction* chapter as Example *A very simple NeXus Data file (in XML)*.

NeXus groups are encoded into XML as elements with the name of the NeXus class and an XML attribute `name` which defines the NeXus name of the group. Further group attributes become XML attributes. An example:

### NeXus group element in XML

```xml
<NXentry name="entry">
</NXentry>
```

NeXus data sets are encoded as XML elements with the name of the data. An attribute `NAPItype` defines the type and dimensions of the data. The actual data is stored as `PCDATA` in the element. Another example:

### NeXus data elements

```xml
<mode NAPItype="NX_CHAR[7]"> monitor </mode>
<counts NAPItype="NX_INT32[4]"> 21 456 127876 319 </counts>
```

Data are printed in appropriate formats and in C storage order. The codes understood for `NAPItype` are all the NeXus data type names. The dimensions are given in square brackets as a comma separated list. No dimensions need to be given if the data is just a single value. Data attributes are represented as XML attributes. If the attribute is not a text string, then the attribute is given in the form: `type:value`, for example: `signal="NX_POSINT:1"`.

NeXus links are stored in XML as XML elements with the name `NAPIlink` and a XML attribute `target` which stores the path to the linked entity in the file. If the item is linked under a different name, then this name is specified as a XML attribute name to the element `NAPIlink`.

---

13 `PCDATA` is the XML term for *parsed character data* (see: [http://www.w3schools.com/xml/xml_cdata.asp](http://www.w3schools.com/xml/xml_cdata.asp)).
The authors of the NeXus API worked with the author of the miniXML XML library to create a reasonably efficient way of handling numeric data with XML. Using the NeXus API handling something like 400 detectors versus 2000 time channels in XML is not a problem. But you may hit limits with XML as the file format when data becomes too large or you try to process NeXus XML files with general XML tools. General XML tools are normally ill prepared to process large amounts of numbers.

**Special Attributes**

NeXus makes use of some special attributes for its internal purposes. These attributes are stored as normal group or data set attributes in the respective file format. These are:

- **target** This attribute is automatically created when items get linked. The target attribute contains a text string with the path to the source of the item linked.

- **napimount** The napimount attribute is used to implement external linking in NeXus. The string is a URL to the file and group in the external file to link too. The system is meant to be extended. But as of now, the only format supported is:

  `nxfile://path-to-file#path-infile`

  This is a NeXus file in the file system at `path-to-file` and the group `path-infile` in that NeXus file.

- **NAPIlink** NeXus supports linking items in another group under another name. This is only supported natively in HDF-5. For HDF-4 and XML a crutch is needed. This crutch is a special class name or attribute `NAPIlink` combined with the target attribute. For groups, `NAPIlink` is the group class, for data items a special attribute with the name `NAPIlink`.

### 1.3 Constructing NeXus Files and Application Definitions

In *NeXus Design*, we discussed the design of the NeXus format in general terms. In this section a more tutorial style introduction in how to construct a NeXus file is given. As an example a hypothetical instrument named WONI will be used.

**Note:** If you are looking for a tutorial on reading or writing NeXus data files using the NeXus API, consult the *NAPI: NeXus Application Programmer Interface (frozen)* chapter. For code examples, refer to *Code Examples that use the NAPI* chapter. Alternatively, there are examples in the *Example NeXus C programs using native HDF5 commands* chapter of writing and reading NeXus data files using the native HDF5 interfaces in C. Further, there are also some Python examples using the `h5py` package in the *Python Examples using h5py* section.

#### 1.3.1 The WONderful New Instrument (WONI)

Consider yourself to be responsible for some hypothetical WONderful New Instrument (WONI). You are tasked to ensure that WONI will record data according to the NeXus standard. For the sake of simplicity, WONI bears a strong resemblance to a simple powder diffractometer, but let’s pretend that WONI cannot use any of the existing NXDL application definitions.

WONI uses collimators and a monochromator to illuminate the sample with neutrons of a selected wavelength as described in *The (fictional) WONI example powder diffractometer*. The diffracted beam is collected in a large, banana-shaped, position sensitive detector. Typical data looks like *Example Powder Diffraction Plot from (fictional) WONI at HYNES*. There is a generous background to the data plus quite a number of diffraction peaks.
Figure 1.7: The (fictional) WONI example powder diffractometer
Figure 1.8: Example Powder Diffraction Plot from (fictional) WONI at HYNES
1.3.2 Constructing a NeXus file for WONI

The starting point for a NeXus file for WONI will be an empty basic NeXus file hierarchy as documented in the next figure. In order to arrive at a full NeXus file, the following steps are required:

1. For each instrument component, decide which parameters need to be stored
2. Map the component parameters to NeXus groups and parameters and add the components to the NXinstrument hierarchy
3. Decide what needs to go into NXdata
4. Fill the NXsample and NXmonitor groups

Basic structure of a NeXus file

```plaintext
entry: NXentry
   NXdata
   NXinstrument
   NXmonitor
   NXsample
```

Decide which parameters need to be stored

Now the various groups of this empty NeXus file shell need to be filled. The next step is to look at a design drawing of WONI. Identify all the instrument components like collimators, detectors, monochromators etc. For each component decide which values need to be stored. As NeXus aims to describe the experiment as good as possible, strive to capture as much information as practical.

Mapping parameters to NeXus

With the list of parameters to store for each component, consult the reference manual section on the NeXus base classes. You will find that for each of your instruments components there will be a suitable NeXus base class. Add this base class together with a name as a group under NXinstrument in your NeXus file hierarchy. Then consult the possible parameter names in the NeXus base class and match them with the parameters you wish to store for your instruments components.

As an example, consider the monochromator. You may wish to store: the wavelength, the d-value of the reflection used, the type of the monochromator and its angle towards the incoming beam. The reference manual tells you that NXcrystal is the right base class to use. Suitable fields for your parameters can be found in there to. After adding them to the basic NeXus file, the file looks like in the next figure:

Basic structure of a NeXus file with a monochromator added

```plaintext
entry: NXentry
   NXdata
   NXinstrument
   monochromator: NXcrystal
      wavelength
      d_spacing
      rotation_angle
      reflection
      type
```
If a parameter or even a whole group is missing in order to describe your experiment, do not despair! Contact the NIAC and suggest to add the group or parameter. Give a little documentation what it is for. The NIAC will check that your suggestion is no duplicate and sufficiently documented and will then proceed to enhance the base classes with your suggestion.

A more elaborate example of the mapping process is given in the section *Creating a NXDL Specification*.

**Decide on NXdata**

The NXdata/ group is supposed to contain the data required to put up a quick plot. For WONI this is a plot of counts versus two theta (polar_angle in NeXus) as can be seen in *Example Powder Diffraction Plot from (fictional) WONI at HYNES*. Now, in NXdata, create links to the appropriate data items in the NXinstrument hierarchy. In the case of WONI, both parameters live in the detector:NXdetector group.

**Fill in auxiliary Information**

Look at the section on NXsample in the NeXus reference manual. Choose appropriate parameters to store for your samples. Probably at least the name will be needed.

In order to normalize various experimental runs against each other it is necessary to know about the counting conditions and especially the monitor counts of the monitor used for normalization. The NeXus convention is to store such information in a control:NXmonitor group at NXentry level. Consult the reference for NXmonitor for field names. If additional monitors exist within your experiment, they will be stored as additional NXmonitor groups at entry level.

Consult the documentation for NXentry in order to find out under which names to store information such as titles, user names, experiment times etc.

A more elaborate example of this process can be found in the following section on creating an application definition.

**1.3.3 Creating a NXDL Specification**

An NXDL specification for a NeXus file is required if you desire to standardize NeXus files from various sources. Another name for a NXDL description is application definition. A NXDL specification can be used to verify NeXus files to conform to the standard encapsulated in the application definition. The process for constructing a NXDL specification is similar to the one described above for the construction of NeXus files.

One easy way to describe how to store data in the NeXus class structure and to create a NXDL specification is to work through an example. Along the way, we will describe some key decisions that influence our particular choices of metadata selection and data organization. So, on with the example ...

**Application Definition Steps**

With all this introductory stuff out of the way, let us look at the process required to define an application definition:

1. *Think!* hard about what has to go into the data file.
2. *Map* the required fields into the NeXus hierarchy
3. *Describe* this map in a NXDL file
4. *Standardize* your definition through communication with the NIAC
Step 1: Think! hard about data

This is actually the hard bit. There are two things to consider:

1. What has to go into the data file?
2. What is the normal plot for this type of data?

For the first part, one of the NeXus guiding principles gives us - Guidance! “A NeXus file must contain all the data necessary for standard data analysis.”

Not more and not less for an application definition. Of course the definition of standard data for analysis or a standard plot depends on the science and the type of data being described. Consult senior scientists in the field about this if you are unsure. Perhaps you must call an international meeting with domain experts to haggle that out. When considering this, people tend to put in everything which might come up. This is not the way to go.

A key test question is: Is this data item necessary for common data analysis? Only these necessary data items belong in an application definition.

The purpose of an application definition is that an author of upstream software who consumes the file can expect certain data items to be there at well defined places. On the other hand if there is a development in your field which analyzes data in a novel way and requires more data to do it, then it is better to err towards the side of more data.

Now for the case of WONI, the standard data analysis is either Rietveld refinement or profile analysis. For both purposes, the kind of radiation used to probe the sample (for WONI, neutrons), the wavelength of the radiation, the monitor (which tells us how long we counted) used to normalize the data, the counts and the two theta angle of each detector element are all required. Usually, it is desirable to know what is being analyzed, so some metadata would be nice: a title, the sample name and the sample temperature. The data typically being plotted is two theta against counts, as shown in Example Powder Diffraction Plot from (fictional) WONI at HYNES above. Summarizing, the basic information required from WONI is given next.

- title of measurement
- sample name
- sample temperature
- counts from the incident beam monitor
- type of radiation probe
- wavelength ($\lambda$) of radiation incident on sample
- angle ($2\theta$ or two theta) of detector elements
- counts for each detector element

If you start to worry that this is too little information, hold on, the section on Using an Application Definition (Using an Application Definition) will reveal the secret how to go from an application definition to a practical file.

Step 2: Map Data into the NeXus Hierarchy

This step is actually easier than the first one. We need to map the data items which were collected in Step 1 into the NeXus hierarchy. A NeXus file hierarchy starts with an NXentry group. At this stage it is advisable to pull up the base class definition for NXentry and study it. The first thing you might notice is that NXentry contains a field named title. Reading the documentation, you quickly realize that this is a good place to store our title. So the first mapping has been found.

```
title = /NXentry/title
```
Note: In this example, the mapping descriptions just contain the path strings into the NeXus file hierarchy with the class names of the groups to use. As it turns out, this is the syntax used in NXDL link specifications. How convenient!

Another thing to notice in the NXentry base class is the existence of a group of class NXsample. This looks like a great place to store information about the sample. Studying the NXsample base class confirms this view and there are two new mappings:

```plaintext
1. sample name = /NXentry/NXsample/name
2. sample temperature = /NXentry/NXsample/temperature
```

Scanning the NXentry base class further reveals there can be a NXmonitor group at this level. Looking up the base class for NXmonitor reveals that this is the place to store our monitor information.

```plaintext
monitor = /NXentry/NXmonitor/data
```

For the other data items, there seem to be no solutions in NXentry. But each of these data items describe the instrument in more detail. NeXus stores instrument descriptions in the /NXentry/NXinstrument branch of the hierarchy. Thus, we continue by looking at the definition of the NXinstrument base class. In there we find further groups for all possible instrument components. Looking at the schematic of WONI (The (fictional) WONI example powder diffractometer), we realize that there is a source, a monochromator and a detector. Suitable groups can be found for these components in NXinstrument and further inspection of the appropriate base classes reveals the following further mappings:

```plaintext
1. probe = /NXentry/NXinstrument/NXsource/probe
2. wavelength = /NXentry/NXinstrument/NXcrystal/wavelength
3. two theta of detector elements = /NXentry/NXinstrument/NXdetector/polar_angle
4. counts for each detector element = /NXentry/NXinstrument/NXdetector/data
```

Thus we mapped all our data items into the NeXus hierarchy! What still needs to be done is to decide upon the content of the NXdata group in NXentry. This group describes the data necessary to make a quick plot of the data. For WONI this is counts versus two theta. Thus we add this mapping:

```plaintext
1. two theta of detector elements = /NXentry/NXdata/polar_angle
2. counts for each detector element = /NXentry/NXdata/data
```

The full mapping of WONI data into NeXus is documented in the next table:

<table>
<thead>
<tr>
<th>WONI data</th>
<th>NeXus path</th>
</tr>
</thead>
<tbody>
<tr>
<td>title of measurement</td>
<td>/NXentry/title</td>
</tr>
<tr>
<td>sample name</td>
<td>/NXentry/NXsample/name</td>
</tr>
<tr>
<td>sample temperature</td>
<td>/NXentry/NXsample/temperature</td>
</tr>
<tr>
<td>monitor</td>
<td>/NXentry/NXmonitor/data</td>
</tr>
<tr>
<td>type of radiation probe</td>
<td>/NXentry/NXinstrument/NXsource/probe</td>
</tr>
<tr>
<td>wavelength of radiation incident on sample</td>
<td>/NXentry/NXinstrument/NXcrystal/wavelength</td>
</tr>
<tr>
<td>two theta of detector elements</td>
<td>/NXentry/NXinstrument/NXdetector/polar_angle</td>
</tr>
<tr>
<td>counts for each detector element</td>
<td>/NXentry/NXinstrument/NXdetector/data</td>
</tr>
<tr>
<td>two theta of detector elements</td>
<td>/NXentry/NXdata/polar_angle</td>
</tr>
<tr>
<td>counts for each detector element</td>
<td>/NXentry/NXdata/data</td>
</tr>
</tbody>
</table>

Looking at this table, one might get concerned that the two theta and counts data is stored in two places and thus duplicated. Stop worrying, this problem is solved at the NeXus API level. Typically NXdata will only hold links to the corresponding data items in /NXentry/NXinstrument/NXdetector.

In this step problems might occur. The first is that the base class definitions contain a bewildering number of parameters. This is on purpose: the base classes serve as dictionaries which define names for everything which possibly can occur. You do not have to give all that information. The key question is, as already said, What is required for typical

1.3. Constructing NeXus Files and Application Definitions 49
data analysis for this type of application? You might also be unsure how to correctly store a particular data item. In such a case, contact the NIAC for help. Another problem which can occur is that you require to store information for which there is no name in one of the existing base classes or you have a new instrument component for which there is no base class altogether. In such a case, please feel free to contact the NIAC with a suggestion for an extension of the base classes in question.

**Step 3: Describe this map in a NXDL file**

This is even easier. Some XML editing is necessary. Fire up your XML editor of choice and open a file. If your XML editor supports XML schema while editing XML, it is worth to load nxdl.xsd. Now your XML editor can help you to create a proper NXDL file. As always, the start is an empty template file. This looks like the XML code below.

**Note:** This is just the basic XML for a NXDL definition. It is advisable to change some of the documentation strings.

### NXDL template file

```xml
<?xml version="1.0" encoding="UTF-8"?>
<!--
# NeXus - Neutron and X-ray Common Data Format
#
# Copyright (C) 2008-2012 NeXus International Advisory Committee (NIAC)
#
# This library is free software; you can redistribute it and/or
# modify it under the terms of the GNU Lesser General Public
# License as published by the Free Software Foundation; either
# version 3 of the License, or (at your option) any later version.
#
# This library is distributed in the hope that it will be useful,
# but WITHOUT ANY WARRANTY; without even the implied warranty of
# MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE. See the GNU
# Lesser General Public License for more details.
#
# You should have received a copy of the GNU Lesser General Public
# License along with this library; if not, write to the Free Software
# Foundation, Inc., 59 Temple Place, Suite 330, Boston, MA 02111-1307 USA
#
# For further information, see http://www.nexusformat.org
-->
<definition name="NX__template__" extends="NXobject" type="group"
    category="application"
    xmlns="http://definition.nexusformat.org/nxdl/3.1"
    xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance"
    xsi:schemaLocation="http://definition.nexusformat.org/nxdl/3.1 ../nxdl.xsd"
    version="1.0b"
>
    <doc>template for a NXDL application definition</doc>
</definition>
```

For example, copy and rename the file to NXwoni.nxdl.xml. Then, locate the XML root element definition and change the name attribute (the XML shorthand for this attribute is /definition/@name) to NXwoni. Change the doc as well. Also consider keeping track of /definition/@version as suits your development of this NXDL file.

The next thing which needs to be done is adding groups into the definition. A group is defined by some XML, as in this example:
The type is the actual NeXus base class this group belongs to. Optionally a name attribute may be given (default is data).

Next, one needs to include data items too. The XML for such a data item looks similar to this:

```xml
<field name="polar_angle" type="NX_FLOAT units="NX_ANGLE">
   <doc>Link to polar angle in /NXentry/NXinstrument/NXdetector</doc>
   <dimensions rank="1">
      <dim index="1" value="ndet"/>
   </dimensions>
</field>
```

The meaning of the name attribute is intuitive, the type can be looked up in the relevant base class definition. A field definition can optionally contain a doc element which contains a description of the data item. The dimensions entry specifies the dimensions of the data set. The size attribute in the dimensions tag sets the rank of the data, in this example: rank="1". In the dimensions group there must be rank dim fields. Each dim tag holds two attributes: index determines to which dimension this tag belongs, the 1 means the first dimension. The value attribute then describes the size of the dimension. These can be plain integers, variables, such as in the example ndet or even expressions like tof+1.

Thus a NXDL file can be constructed. The full NXDL file for the WONI example is given in Full listing of the WONI Application Definition. Clever readers may have noticed the strong similarity between our working example NXwoni and NXmonopd since they are essentially identical. Give yourselves a cookie if you spotted this.

**Step 4: Standardize with the NIAC**

Basically you are done. Your first application definition for NeXus is constructed. In order to make your work a standard for that particular application type, some more steps are required:

- Send your application definition to the NIAC for review
- Correct your definition per the comments of the NIAC
- Cure and use the definition for a year
- After a final review, it becomes the standard

The NIAC must review an application definition before it is accepted as a standard. The one year curation period is in place in order to gain practical experience with the definition and to sort out bugs from Step 1. In this period, data shall be written and analyzed using the new application definition.

**Full listing of the WONI Application Definition**

```xml
<?xml version="1.0" encoding="UTF-8"?>
<?xml-stylesheet type="text/xsl" href="nxdlformat.xsl" ?>
<!--
# NeXus - Neutron and X-ray Common Data Format
#
# Copyright (C) 2008-2012 NeXus International Advisory Committee (NIAC)
#
# This library is free software; you can redistribute it and/or
# modify it under the terms of the GNU Lesser General Public
# License as published by the Free Software Foundation; either
# version 3 of the License, or (at your option) any later version.
```
# This library is distributed in the hope that it will be useful,
# but WITHOUT ANY WARRANTY; without even the implied warranty of
# MERCHANTABILITY or FITNESS FOR A PARTICULAR PURPOSE. See the GNU
# Lesser General Public License for more details.
#
# You should have received a copy of the GNU Lesser General Public
# License along with this library; if not, write to the Free Software
# Foundation, Inc., 59 Temple Place, Suite 330, Boston, MA 02111-1307 USA
#
# For further information, see http://www.nexusformat.org

<definition name="NXmonopd" extends="NXobject" type="group"
category="application"
xmlns="http://definition.nexusformat.org/nxdl/@NXDL_RELEASE@"
xmns:xsi="http://www.w3.org/2001/XMLSchema-instance"
xsi:schemaLocation="http://definition.nexusformat.org/nxdl/@NXDL_RELEASE@ ../nxdl.xsd"
version="1.0b">

<doc>
Monochromatic Neutron and X-Ray Powder Diffraction. Instrument
definition for a powder diffractometer at a monochromatic neutron
or X-ray beam. This is both suited for a powder diffractometer
with a single detector or a powder diffractometer with a position
sensitive detector.
</doc>

<group type="NXentry" name="entry">
  <field name="title"/>
  <field name="start_time" type="NX_DATE_TIME"/>
  <field name="definition">
    <doc>
      Official NeXus NXDL schema to which this file conforms
    </doc>
    <enumeration>
      <item value="NXmonopd"/>
    </enumeration>
  </field>
</group>

<group type="NXinstrument">
  <group type="NXsource">
    <field name="type"/>
    <field name="name"/>
    <field name="probe">
      <enumeration>
        <item value="neutron"/>
        <item value="x-ray"/>
        <item value="electron"/>
      </enumeration>
    </field>
  </group>
  <group type="NXcrystal">
    <field name="wavelength" type="NX_FLOAT" units="NX_WAVELENGTH">
      <doc>
        Optimum diffracted wavelength
      </doc>
      <dimensions rank="1">
        <dim index="1" value="i"/>
      </dimensions>
    </field>
  </group>
  <group type="NXdetector">
    <field name="polar_angle" type="NX_FLOAT" axis="1">
      <doc>
        where ndet = number of detectors
      </doc>
    </field>
  </group>
</group>
<dimensions rank="1">
    <dim index="1" value="ndet" />
</dimensions>

<field name="data" type="NX_INT" signal="1">
    
    <doc>
        Detector signal (usually counts) are already corrected for detector efficiency
    </doc>
    
    <dimensions rank="1">
        <dim index="1" value="ndet" />
    </dimensions>
</field>

<group type="NXsample">
    <field name="name">
        
        <doc>Descriptive name of sample</doc>
    </field>

    <field name="rotation_angle" type="NX_FLOAT" units="NX_ANGLE">
        
        <doc>
            Optional rotation angle for the case when the powder diagram has been obtained through an omega-2theta scan like from a traditional single detector powder diffractometer
        </doc>
    </field>
</group>

<group type="NXmonitor">
    <field name="mode">
        
        <doc>Count to a preset value based on either clock time (timer) or received monitor counts (monitor).
        </doc>

        <enumeration>
            <item value="monitor"/>
            <item value="timer"/>
        </enumeration>
    </field>

    <field name="preset" type="NX_FLOAT">
        
        <doc>Preset value for time or monitor</doc>
    </field>

    <field name="integral" type="NX_FLOAT" units="NX_ANY">
        
        <doc>Total integral monitor counts</doc>
    </field>
</group>

<group type="NXdata">
    <link name="polar_angle" target="/NXentry/NXinstrument/NXdetector/polar_angle">
        
        <doc>Link to polar angle in /NXentry/NXinstrument/NXdetector</doc>
    </link>

    <link name="data" target="/NXentry/NXinstrument/NXdetector/data">
        
        <doc>Link to data in /NXentry/NXinstrument/NXdetector</doc>
    </link>
</group>

</definition>
Using an Application Definition

The application definition is like an interface for your data file. In practice files will contain far more information. For this, the extendable capability of NeXus comes in handy. More data can be added, and upstream software relying on the interface defined by the application definition can still retrieve the necessary information without any changes to their code.

NeXus application definitions only standardize classes. You are free to decide upon names of groups, subject to them matching regular expression for NeXus name attributes (see the regular expression pattern for NXDL group and field names in the Naming Conventions section). Note the length limit of 63 characters imposed by HDF5. Please use sensible, descriptive names and separate multi worded names with underscores.

Something most people wish to add is more metadata, for example in order to index files into a database of some sort. Go ahead, do so, if applicable, scan the NeXus base classes for standardized names. For metadata, consider to use the NXarchive definition. In this context, it is worth to mention that a practical NeXus file might adhere to more than one application definition. For example, WONI data files may adhere to both the NXmonopd and NXarchive definitions. The first for data analysis, the second for indexing into the database.

Often, instrument scientists want to store the complete state of their instrument in data files in order to be able to find out what went wrong if the data is unsatisfactory. Go ahead, do so, please use names from the NeXus base classes.

Site policy might require you to store the names of all your bosses up to the current head of state in data files. Go ahead, add as many NXuser classes as required to store that information. Knock yourselves silly over this.

Your Scientific Accounting Department (SAD) may ask of you the preposterous; to store billing information into data files. Go ahead, do so if your judgment allows. Just do not expect the NIAC to provide base classes for this and do not use the prefix NX for your classes.

In most cases, NeXus files will just have one NXentry class group. But it may be required to store multiple related data sets of the results of data analysis into the same data file. In this case create more entries. Each entry should be interpretable standalone, i.e. contain all the information of a complete NXentry class. Please keep in mind that groups or data items which stay constant across entries can always be linked in.

1.3.4 Processed Data

Data reduction and analysis programs are encouraged to store their results in NeXus data files. As far as the necessary, the normal NeXus hierarchy is to be implemented. In addition, processed data files must contain a NXprocess group. This group, that documents and preserves data provenance, contains the name of the data processing program and the parameters used to run this program in order to achieve the results stored in this entry. Multiple processing steps must have a separate entry each.

1.4 Strategies for storing information in NeXus data files

NeXus may appear daunting, at first, to use. The number of base classes is quite large as well as is the number of application definitions. This chapter describes some of the strategies that have been recommended for how to store information in NeXus data files.

When we use the term storing, some might be helped if they consider this as descriptions for how to classify their data. It is intended for this chapter to grow, with the addition of different use cases as they are presented for suggestions.

1.4.1 Strategies: The simplest case(s)

Perhaps the simplest case might be either a step scan with two or more columns of data. Another simple case might be a single image acquired by an area detector. In either of these hypothetical cases, the situation is so simple that there
is little addition information available to be described (for whatever reason).

**Step scan with two or more data columns**

Consider the case where we wish to store the data from a step scan. This case may involve two or more related 1-D arrays of data to be saved, each having the same length. For our hypothetical case, we’ll have these positioners as arrays:

<table>
<thead>
<tr>
<th>positioner arrays</th>
<th>detector arrays</th>
</tr>
</thead>
<tbody>
<tr>
<td>ar, ay, dy</td>
<td>I0, I00, time, Epoch, photodiode</td>
</tr>
</tbody>
</table>

### 1.4.2 Strategies: The wavelength

*Where should the wavelength of my experiment be written?* This is one of the *Frequently Asked Questions*. The canonical location to store wavelength has been:

/NXentry/NXinstrument/NXcrystal/wavelength

More recently, this location makes more sense to many:

/NXentry/NXinstrument/NXmonochromator/wavelength

**NXcrystal** describes a crystal monochromator or analyzer. Recently, scientists with monochromatic radiation not defined by a crystal, such as from an electron-beam undulator or a neutron helical velocity selector, were not satisfied with creating a fictitious instance of a crystal just to preserve the wavelength from their instrument. Thus, the addition of the **NXmonochromator** base class to NeXus, which also allows “energy” to be specified if one is so inclined.

**Note:** See the *Class path specification* section for a short discussion of the difference between the HDF5 path and the NeXus symbolic class path.

### 1.4.3 Strategies: The next case

The **NIAC: The NeXus International Advisory Committee** welcomes suggestions for additional sections in this chapter.

### 1.5 Verification and validation of files

The intent of verification and validation of files is to ensure, in an unbiased way, that a given file conforms to the relevant specifications. NeXus uses various automated tools to validate files. These tools include conversion of content from HDF to XML and transformation (via XSLT) from XML format to another such as NXDL, XSD, and Schematron. This chapter will first provide an overview of the process, then define the terms used in validation, then describe how multiple base classes or application definitions might apply to a given NeXus data file, and then describe the various validation techniques in more detail. Validation does not check that the data content of the file is sensible; this requires scientific interpretation based on the technique.

Validation is useful to anyone who manipulates or modifies the contents of NeXus files. This includes scientists/users, instrument staff, software developers, and those who might mine the files for metadata. First, the scientist or user of the data must be certain that the information in a file can be located reliably. The instrument staff or software developer must be confident the information they have written to the file has been located and formatted properly. At some time, the content of the NeXus file may contribute to a larger body of work such as a metadata catalog for a scientific instrument, a laboratory, or even an entire user facility.
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

1.5.1 Overview

NeXus files adhere to a set of rules and can be tested against these rules for compliance. The rules are implemented using standard tools and can themselves be tested to verify compliance with the standards for such definitions. Validation includes the testing of both NeXus data files and the NXDL specifications that describe the rules.

The rules for writing NeXus data files are different than the rules for writing NeXus class definitions. To validate a NeXus data file, these two rule sets must eventually merge, as shown in the next figure. The data file (either HDF4, HDF5, or XML) is first converted into an internal format to facilitate validation, including data types, array dimensions, naming, and other items. Most of the data is not converted since data validation is non-trivial. Also note that the units are not validated. All the NXDL files are converted into a single Schematron file (again, internal use for validation) only when NXDL revisions are checked into the NeXus definitions repository as NXDL changes are not so frequent.

Figure 1.9: Flowchart of the NeXus validation process.

**NeXus data files**  NeXus data files (also known as NeXus data file instances) are validated to ensure the various parts of the data file are arranged according to the governing NXDL specifications used in that file instance.

**Note:** Since NeXus has several rules that are quite difficult to apply in either XSD or Schematron, direct validation of data files using standard tools is not possible. To validate NeXus data files, it is necessary to use nxvalidate.

**NeXus Definition Language (NXDL) specification files**  NXDL files are validated to ensure they adhere to the rules for writing NeXus base classes and application definitions.

1.5.2 Definitions of these terms

Let’s be clear about some terms used in this section.

**HDF**  Hierarchical Data Format from The HDF Group. NeXus data files using HDF may be stored in either version 4 (HDF4) or version 5 (HDF5). New NeXus HDF files should only use HDF5. The preferred file extensions (but not required) include .hdf, .h5, .nxs, and .nx5.

**NXDL**  NeXus Definition Language files define the specifications for NeXus base classes, application definitions, and contributed classes and definitions. It is fully described in the NXDL chapter.

**Schematron**  Schematron is an alternative to XSD and is used to validate the content and structure of an XML file. NeXus uses Schematron internally to validate data files.

14 http://www.schematron.com
**Validation**  File validation is the comparison of file contents, in an unbiased way, with the set of rules that define the structure of such files.

**XML**  The eXtensible Markup Language (XML)\(^{15}\) is a standard business tool for the exchange of information. It is broadly supported by a large software library in many languages. NeXus uses XML for several purposes: data files, NXDL definitions, rules, and XSLT transformations.

**XSD**  XML files are often defined by a set of rules (or *schema*). A common language used to implement these rules is XML Schema (XSD)\(^{16}\) Fundamentally, all XML, XSD, XSLT, and Schematron files are XML.

**XSLT**  XML files can be flexible enough to convert from one set of rules to another. An example is when one company wishes to exchange catalog or production information with another. The XML StyL-sheet Transformation (XSLT)\(^{17}\) language is often used to describe each direction of the conversion of the XML files between the two rule sets.

### 1.5.3 NeXus data files may use multiple base classes or application definitions

NeXus data files may have more than one data set or may have multiple instances of just about any base class or even application definitions. The NeXus data file validation is prepared to handle this without any special effort by the provider of the data file.

### 1.5.4 Validation techniques

File validation is the process to determine if a given file is prepared consistent with a set of guidelines or rules. In NeXus, there are several different types of files. First, of course, is the data file yet it can be provided in one of several forms: HDF4, HDF5, or XML. Specifications for data files are provided by one or (usually) more NeXus definition files (NXDL, for short). These NXDL files are written in XML and validated by the NXDL specification which is written in the XML Schema (XSD) language. Thus, automated file verification is available for data files, definition files, and the rules for definition files.

**Validation of NeXus data files**

Each NeXus data file can be validated against the NXDL rules. (The full suite of NXDL specifications is converted into Schematron rules by an XSLT transformation and then combined into a single file. It is not allowed to have a NeXus base class and also an application definition with the same name since one will override the other in the master Schematron file) The validation is done using Schematron and the NXvalidate program. Schematron was selected, rather than XML Schema (XSD), to permit established rules for NeXus files, especially the rule allowing the nodes within `NXentry` to appear in any order.

The validation process is mainly checking file structure (presence or absence of groups/fields) - it is usually impossible to check the actual data itself, other than confirm that it is of the correct data type (string, float etc.). The only exception is when the NXDL specification is either a fixed value or an enumeration - in which case the data is checked.

During validation, the NeXus data file instance (either HDF or XML) is first converted into an XML file in a form that facilitates validation (e.g. with large numeric data removed). Then the XML file is validated by Schematron against the `schema/all.sch` file.

---

15 http://www.w3schools.com/xml
16 http://www.w3schools.com/schema
17 http://www.w3schools.com/xsl/
Validation of NeXus Definition Language (NXDL) specification files

Each NXDL file must be validated against the rules that define how NXDL files are to be arranged. The NXDL rules are specified in the form of XML Schema (XSD).

Standard tools (validating editor or command line or support library) can be used to validate any NXDL file. Here’s an example using `xmllint` from a directory that contains `nxdl.xsd`, `nxdlTypes.xsd`, and `applications/NXsas.nxdl.xml`:

Use of `xmllint` to validate a NXDL specification.

`xmllint --noout --schema nxdl.xsd applications/NXsas.nxdl.xml`

Validation of the NXDL rules

NXDL rules are specified using the rules of XML Schema (XSD). The XSD syntax of the rules is validated using standard XML file validation tools: either a validating editor (such as oXygen, xmlSpy, or eclipse) or common UNIX/Linux command line tools

Use of `xmllint` to validate the NXDL rules.

`xmllint --valid nxdl.xsd`

The validating editor method is used by the developers while the `xmllint` command line tool is the automated method used by the NeXus definitions subversion repository.

Validation of XSLT files

XSLT transformations are validated using standard tools such as a validating editor or `xmllint`.

Transformation of NXDL files to Schematron

Schematron ¹ is a rule-based language that allows very specific validation of an XML document. Its advantages over using XSD schema are that:

- more specific pattern-based rules based on data content can be written
- full XSLT/XPath expression syntax available for writing validation tests
- error messages can be customised and thus more meaningful
- It is easier to validate documents when entities can occur in any order.

XSD does provide a mechanism for defining a class structure and inheritance, so its usage within NeXus in addition to schematron has not been ruled out. But for a basic validation of file content, schematron looks best.

The NXDL definition files are converted into a set of Schematron rules using the `xslt/nxdl2sch.xsl` XSLT stylesheet. The NeXus instance file (either in XML, HDF4, or HDF5) is turned into a reduced XML validation file. This file is very similar to a pure NeXus XML file, but with additional metadata for dimensions and also with most of the actual numeric data removed.

The validation process then compares the set of Schematron rules against the reduced XML validation file. Schematron itself is implemented as a set of XSLT transforms. NeXus includes the Schematron files, as well as the Java based XSLT engine saxon.
The Java based nxvalidate GUI can be run to validate files. Currently, the structure of the file is validated (i.e. valid names are used at the correct points), but this will be extended to array dimensions and link targets. Error messages are printed about missing mandatory fields, and informational messages are printed about fields that are neither optional or mandatory (in case they are a typing error). Even non-standard names must comply with a set of rules (e.g. no spaces are allowed in names). Enumerations are checked that they conform to an allowed value. The data type is checked and the units will also be checked.

### 1.6 Frequently Asked Questions

This is a list of commonly asked questions concerning the NeXus data format.

1. **Is it Nexus, NeXus or NeXuS?**

   NeXus is correct. It is a format for data from Neutron and X-ray facilities, hence those first letters are capitalised. The format is also used for muon experiments, but there is no \textit{mu} (or m) in NeXus and no s in muon. So the s stays in lower case.

2. **How many facilities use NeXus?**

   This is not easy to say, not all facilities using NeXus actively participate in the committee. Some facilities have reported their adoption status on the Facilities Wiki page. Please have a look at this list. Keep in mind that it is not complete.

3. **NeXus files are binary? This is crazy! How am I supposed to see my data?**

   NeXus files are not per se binary. If you use the XML backend the data are stored in a relatively human readable form (see this example). This backend however is only recommended for very small data sets. With the multidimensional data that is routinely recorded on many modern instruments it is very difficult anyway to retrieve useful information on a VT100 terminal. If you want to try, for example \texttt{nxbrowse} is a utility provided by the NeXus community that can be very helpful to those who want to inspect their files and avoid graphical applications. For larger data volumes the binary backends used with the appropriate tools are by far superior in terms of efficiency and speed and most users happily accept that after having worked with supersized “human readable” files for a while.

4. **What on-disk file format should I choose for my data?**

   HDF5 is the default file container to use for NeXus data. It is the recommended format for all applications. HDF4 is still supported as a on disk format for NeXus but for new installations preference should be given to HDF5. The XML backend is available for special use cases. Choose this option with care considering the space and speed implications.

5. **Why are the NeXus classes so complicated? I’ll never store all that information**

   The NeXus classes are essentially glossaries of terms. If you need to store a piece of information, consult the class definitions to see if it has been defined. If so, use it. It is not compulsory to include every item that has been defined in the base class if it is not relevant to your experiment. On the other hand, a NeXus application definition lists a smaller set of compulsory items that should allow other researchers or software to analyze your data. You should really follow the application definition that corresponds to your experiment to take full advantage of NeXus.

6. **I don’t like NeXus. It seems much faster and simpler to develop my own file format. Why should I even consider NeXus?**

   If you consider using an efficient on disk storage format, HDF5 is a better choice than most others. It is fast and efficient and well supported in all mainstream programming languages and a fair share of popular analysis packages. The format is so widely used and backed by a big organisation that it will continue to be supported for the foreseeable future. So if you are going to use HDF5 anyway, why not use the NeXus definition to lay out the data in a standardised way? The NeXus community
spent years trying to get the standard right and while you will not agree with every single choice they made in the past, you should be able to store the data you have in a quite reasonable way. If you do not comply with NeXus, chances are most people will perceive your format as different but not necessarily better than NeXus by any large measure. So it may not be worth the effort. Seriously.

If you encounter any problems because the classes are not sufficient to describe your configuration, please contact the NIAC Executive Secretary explaining the problem, and post a suggestion at the relevant class wiki page. Or raise the problem in one of the mailing lists. The NIAC is always willing to consider new proposals.

7. I want to contribute an application definition. How do I go about it?

Read the NXDL Tutorial in Creating a NXDL Specification and have a try. You can ask for help on the mailing lists. Once you have a definition that is working well for at least your case, you can submit it to the NIAC for acceptance as a standard. The procedures for acceptance are defined in the NIAC constitution.

8. What is the purpose of NXdata?

NXdata contains links to the data stored elsewhere in the NXentry. It identifies the default plottable data. This is one of the basic motivations (see Simple plotting) for the NeXus standard. The choice of the name NXdata is historic and does not really reflect its function.

9. How do I identify the plottable data?

See the section: Find the plottable data.

10. How can I specify reasonable axes for my data?

See the section: Linking Multi Dimensional Data with Axis Data.

11. Why aren’t NXsample and NXmonitor groups stored in the NXinstrument group?

A NeXus file can contain a number of NXentry groups, which may represent different scans in an experiment, or sample and calibration runs, etc. In many cases, though by no means all, the instrument has the same configuration so that it would be possible to save space by storing the NXinstrument group once and using multiple links in the remaining NXentry groups. It is assumed that the sample and monitor information would be more likely to change from run to run, and so should be stored at the top level.

12. Specifications are complicated and often provide too much information for what I need. Where can I find some good example data files?

There are a few checked into the definitions repository. At the moment the selection is quite limited and not very representative. This repository will be edited as more example files become available.

13. Can I use a NXDL specification to parse a NeXus data file?

This should be possible as there is nothing in the NeXus specifications to prevent this but it is not implemented in NAPI. You would need to implement it for yourself.

14. Why do I need to specify the NAPItype? My programming language does not need that information and I don’t care about C and colleagues. Can I leave it out?

NAPItype is necessary. When implementing the NeXus-XML API we strived to make this as general as HDF and reasonably efficient for medium sized datasets. This is why we store arrays as a large bunch of numbers in C-storage order. And we need the NAPItype to figure out the dimensions of the dataset.

15. Do I have to use the NAPI subroutines? Can’t I read (or write) the NeXus data files with my own routines?

Refer to the most recent version of the NIAC constitution on the NIAC wiki: http://wiki.nexusformat.org/NIAC#Constitution
You are not required to use the NAPI to write valid NeXus data files. It is possible to avoid the NAPI to write and read valid NeXus data files. But, the programmer who chooses this path must have more understanding of how the NeXus HDF or XML data file is written. Validation of data files written without the NAPI is strongly encouraged.

16. I’m using links to place data in two places. Which one should be the data and which one is the link?

**Note:** NeXus uses HDF5 hard links

In HDF, a hard link points to a data object. A soft link points to a directory entry. Since NeXus uses hard links, there is no need to distinguish between two (or more) directory entries that point to the same data.

Both places have pointers to the actual data. That is the way hard links work in HDF5. There is no need for a preference to either location. NeXus defines a `target` attribute to label one directory entry as the source of the data (in this, the link `target`). This has value in only a few situations such as when converting the data from one format to another. By identifying the original in place, duplicate copies of the data are not converted.

17. **If I write my data according to the current specification for NXsas** (substitute any other application definition), will other software be able to read my data?

Yes. NXsas, like other Application Definitions, defines and names the minimum information required for analysis or data processing. As long as all the information required by the specification is present, analysis software should be able to process the data. If other information is also present, there is no guarantee that small-angle scattering analysis software will notice.

18. Where do I store the wavelength of my experiment?

See the Strategies: The wavelength section.

19. Where do I store metadata about my experiment?

See the Where to Store Metadata section.
EXAMPLES OF WRITING AND READING NEXUS DATA FILES

Simple examples of reading and writing NeXus data files are provided in the NeXus Introduction chapter and also in the NAPI: NeXus Application Programmer Interface (frozen) chapter. Here, three examples are provided showing how to write a NeXus data file without using the NAPI.

2.1 Code Examples that use the NAPI

Various examples are given that show how to read and write NeXus data files using the NAPI: NeXus Application Programmer Interface (frozen).

2.1.1 Example NeXus programs using NAPI

NAPI Simple 2-D Write Example (C, F77, F90)

Code examples are provided in this section that write 2-D data to a NeXus HDF5 file in C, F77, and F90 languages using the NAPI.

The following code reads a two-dimensional set counts with dimension scales of t and phi using local routines, and then writes a NeXus file containing a single NXentry group and a single NXdata group. This is the simplest data file that conforms to the NeXus standard. The same code is provided in C, F77, and F90 versions. Compare these code examples with Example NeXus C programs using native HDF5 commands.

NAPI C Example: write simple NeXus file

```c
#include "napi.h"

int main()
{
  int counts[50][1000], n_t=1000, n_p=50, dims[2], i;
  float t[1000], phi[50];
  NXhandle file_id;

  /*
  * Read in data using local routines to populate phi and counts
  * for example you may create a getdata() function and call
  * ...
  */
```
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

Chapter 2. Examples of writing and reading NeXus data files

NAPI F77 Example: write simple NeXus file

Note: The F77 interface is no longer being developed.
+ (file_id, 'user', 'Joe Bloggs', 10, NX_CHAR)

!Open top-level NXentry group
status = NXmakegroup (file_id, 'Entry1', 'NXentry')
status = NXopengroup (file_id, 'Entry1', 'NXentry')

!Open NXdata group within NXentry group
status = NXmakegroup (file_id, 'Data1', 'NXdata')
status = NXopengroup (file_id, 'Data1', 'NXdata')

!Output time channels
status = NXmakedata (file_id, 'time_of_flight', NX_FLOAT32, 1, n_t)
status = NXopendata (file_id, 'time_of_flight')
status = NXputdata (file_id, t)

status = NXputcharattr (file_id, 'units', 'microseconds', 12, NX_CHAR)
status = NXclosedata (file_id)

!Output detector angles
status = NXmakedata (file_id, 'polar_angle', NX_FLOAT32, 1, n_p)
status = NXopendata (file_id, 'polar_angle')
status = NXputdata (file_id, phi)

status = NXputcharattr (file_id, 'units', 'degrees', 7, NX_CHAR)
status = NXclosedata (file_id)

!Output data
dims(1) = n_t
dims(2) = n_p
status = NXmakedata (file_id, 'counts', NX_INT32, 2, dims)
status = NXopendata (file_id, 'counts')
status = NXputdata (file_id, counts)
status = NXputattr (file_id, 'signal', 1, 1, NX_INT32)
status = NXputattr (file_id, 'axes', 'polar_angle:time_of_flight', 26, NX_CHAR)
status = NXclosedata (file_id)

!Close NXdata and NXentry groups and close file
status = NXclosegroup (file_id)
status = NXclosegroup (file_id)
status = NXclose (file_id)

stop
end

NAPI F90 Example: write simple NeXus file

program WRITEDATA
use NXUmodule
type(NXhandle) :: file_id
integer, pointer :: counts(:, :)
real, pointer :: t(:, ), phi(:, )

!Use local routines to allocate pointers and fill in data
call getlocaldata (t, phi, counts)

!Open output file
if (NXopen("NXfile.nxs", NXACC_CREATE, file_id) /= NX_OK) stop
if (NXUwriteglobals (file_id, user="Joe Bloggs") /= NX_OK) stop

!Set compression parameters
if (NXUsetcompress (file_id, NX_COMP_LZW, 1000) /= NX_OK) stop

2.1. Code Examples that use the NAPI
!Open top-level NXentry group
  if (NXUwritegroup (file_id, "Entry1", "NXentry") /= NX_OK) stop
!Open NXdata group within NXentry group
  if (NXUwritegroup (file_id, "Data1", "NXdata") /= NX_OK) stop
!Output time channels
  if (NXUwritedata (file_id, "time_of_flight", t, "microseconds") /= NX_OK) stop
!Output detector angles
  if (NXUwritedata (file_id, "polar_angle", phi, "degrees") /= NX_OK) stop
!Output data
  if (NXUwritedata (file_id, "counts", counts, "counts") /= NX_OK) stop
  if (NXputattr (file_id, "signal", 1) /= NX_OK) stop
  if (NXputattr (file_id, "axes", "polar_angle:time_of_flight") /= NX_OK) stop
!Close NXdata group
  if (NXclosegroup (file_id) /= NX_OK) stop
!Close NXentry group
  if (NXclosegroup (file_id) /= NX_OK) stop
!Close NeXus file
  if (NXclose (file_id) /= NX_OK) stop
end program WRITEDATA

NAPI Python Simple 3-D Write Example

A single code example is provided in this section that writes 3-D data to a NeXus HDF5 file in the Python language using the NAPI. The data file may be retrieved from the repository of NeXus data file examples:

data http://svn.nexusformat.org/definitions/examplendata/simple3D.h5

The data to be written to the file is a simple three-dimensional array (2 x 3 x 4) of integers. The single dataset is intended to demonstrate the order in which each value of the array is stored in a NeXus HDF5 data file.

NAPI Python Example: write simple NeXus file

#!/usr/bin/python
import sys
import nx
import numpy
nf = nx.open("simple3D.h5", "w5")
nf.makegroup("entry","NXentry")
nf.opengroup("entry","NXentry")
nf.makegroup("data","NXdata")
nf.opengroup("data","NXdata")
a = numpy.zeros((2,3,4),dtype=numpy.int)
val = 0
for i in range(2):
  for j in range(3):
    for k in range(4):
      a[i,j,k] = val
      val = val + 1
nf.makedata("test","int32",[2,3,4])
View a NeXus HDF5 file using *h5dump*

For the purposes of an example, it is instructive to view the content of the NeXus HDF5 file produced by the above program. Since HDF5 is a binary file format, we cannot show the contents of the file directly in this manual. Instead, we first view the content by showing the output from the *h5dump* tool provided as part of the HDF5 tool kit:

```bash
h5dump simple3D.h5
```

NAPI Python Example: *h5dump* output of NeXus HDF5 file

```python
HDF5 "simple3D.h5" {
GROUP "/" {
  ATTRIBUTE "NeXus_version" {
    DATATYPE H5T_STRING {
      STRSIZE 5;
      STRPAD H5T_STR_NULLTERM;
      CSET H5T_CSET_ASCII;
      CType H5T_C_S1;
    }
    DATASPACE SCALAR
    DATA { (0): "4.1.0" }
  }
  ATTRIBUTE "file_name" {
    DATATYPE H5T_STRING {
      STRSIZE 11;
      STRPAD H5T_STR_NULLTERM;
      CSET H5T_CSET_ASCII;
      CType H5T_C_S1;
    }
    DATASPACE SCALAR
    DATA { (0): "simple3D.h5" }
  }
  ATTRIBUTE "HDF5_Version" {
    DATATYPE H5T_STRING {
      STRSIZE 5;
      STRPAD H5T_STR_NULLTERM;
      CSET H5T_CSET_ASCII;
      CType H5T_C_S1;
    }
    DATASPACE SCALAR
    DATA { (0): "1.10.0" }
  }
}
```
DATA {
  (0): "1.6.6"
}

ATTRIBUTE "file_time" {
  DATATYPE H5T_STRING {
    STRSIZE 24;
    STRPAD H5T_STR_NULLTERM;
    CSET H5T_CSET_ASCII;
    CTYPE H5T_C_S1;
  }
  DATASPACE SCALAR
  DATA {
    (0): "2011-11-18 17:26:27+0100"
  }
}

GROUP "entry" {
  ATTRIBUTE "NX_class" {
    DATATYPE H5T_STRING {
      STRSIZE 7;
      STRPAD H5T_STR_NULLTERM;
      CSET H5T_CSET_ASCII;
      CTYPE H5T_C_S1;
    }
    DATASPACE SCALAR
    DATA {
      (0): "NXentry"
    }
  }
  GROUP "data" {
    ATTRIBUTE "NX_class" {
      DATATYPE H5T_STRING {
        STRSIZE 6;
        STRPAD H5T_STR_NULLTERM;
        CSET H5T_CSET_ASCII;
        CTYPE H5T_C_S1;
      }
      DATASPACE SCALAR
      DATA {
        (0): "NXdata"
      }
    }
  }
}

DATASET "test" {
  DATATYPE H5T_STD_I32LE
  DATASPACE SIMPLE { ( 2, 3, 4 ) / ( 2, 3, 4 ) }
  DATA {
    (0,0,0): 0, 1, 2, 3,
    (0,1,0): 4, 5, 6, 7,
    (0,2,0): 8, 9, 10, 11,
    (1,0,0): 12, 13, 14, 15,
    (1,1,0): 16, 17, 18, 19,
    (1,2,0): 20, 21, 22, 23
  }
  ATTRIBUTE "signal" {
    DATATYPE H5T_STD_I32LE
    DATASPACE SCALAR
    DATA {
      (0): 1
    }
  }
}
View a NeXus HDF5 file using h5toText.py

The output of h5dump contains a lot of structural information about the HDF5 file that can distract us from the actual content we added to the file. Next, we show the output from a custom Python tool (h5toText.py) that we describe in a later section (h5toText support module) of this chapter. This tool was developed to show the actual data content of an HDF5 file that we create.

NAPI Python Example: h5toText output of NeXus HDF5 file

```python
simple3D.h5: NeXus data file
  @NeXus_version = 4.1.0
  @file_name = simple3D.h5
  @HDF5_Version = 1.6.6
  @file_time = 2011-11-18 17:26:27+0100
entry: NXentry
  @NX_class = NXentry
data: NXdata
  @NX_class = NXdata
test: NX_INT32[2,3,4] = __array
  @signal = 1
  __array = [
    [1, 1, 2, 3]
    [4, 5, 6, 7]
    [8, 9, 10, 11]
  ]
  [
    [12, 13, 14, 15]
    [16, 17, 18, 19]
    [20, 21, 22, 23]
  ]
```

2.2 Code Examples that do not use the NAPI

Sometimes, for whatever reason, it is necessary to write or read NeXus files without using the routines provided by the NAPI: NeXus Application Programmer Interface (frozen). Each example in this section is written to support just one of the low-level file formats supported by NeXus (HDF4, HDF5, or XML).

2.2.1 Example NeXus C programs using native HDF5 commands

C-language code examples are provided for writing and reading NeXus-compliant files using the native HDF5 interfaces. These examples are derived from the simple NAPI examples for writing and reading given in the Introduction.
chapter. Compare these code examples with *Example NeXus programs using NAPI*.

**Writing a simple NeXus file using native HDF5 commands in C**

```c
#include <hdf5.h>
#include <stdlib.h>
#include <string.h>

#define LENGTH 400

int main(int argc, char *argv[])
{
    float two_theta[LENGTH];
    int counts[LENGTH], i, rank, signal;

    /* HDF-5 handles */
    hid_t fid, fapl, gid, atts, atttype, attid;
    hid_t datatype, dataspace, dataprop, dataid;
    hsize_t dim[1], maxdim[1];

    /* create some data: nothing NeXus or HDF-5 specific */
    for(i = 0; i < LENGTH; i++)
    {
        two_theta[i] = 10. + 0.1 * i;
        counts[i] = (int)(1000 * ((float)random()/(float)RAND_MAX));
    }
    dim[0] = LENGTH;
    maxdim[0] = LENGTH;
    rank = 1;
```

NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

NeXus programs using NAPI.
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

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* open the file. The file attribute forces normal file
* closing behaviour down HDF-5’s throat
*/
fapl = H5Pcreate(H5P_FILE_ACCESS);
H5Pset_fclose_degree(fapl,H5F_CLOSE_STRONG);
fid = H5Fcreate("NXfile.h5", H5F_ACC_TRUNC, H5P_DEFAULT,fapl);
H5Pclose(fapl);

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/*
* create scan:NXentry
*/
gid = H5Gcreate(fid, (const char *)"scan",0);
/*
* store the NX_class attribute. Notice that you
* have to take care to close those hids after use
*/
atts = H5Screate(H5S_SCALAR);
atttype = H5Tcopy(H5T_C_S1);
H5Tset_size(atttype, strlen("NXentry"));
attid = H5Acreate(gid,"NX_class", atttype, atts, H5P_DEFAULT);
H5Awrite(attid, atttype, (char *)"NXentry");
H5Sclose(atts);
H5Tclose(atttype);
H5Aclose(attid);

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/*
* same thing for data:Nxdata in scan:NXentry.
* A subroutine would be nice to have here.......
*/
gid = H5Gcreate(fid, (const char *)"/scan/data",0);
atts = H5Screate(H5S_SCALAR);
atttype = H5Tcopy(H5T_C_S1);
H5Tset_size(atttype, strlen("NXdata"));
attid = H5Acreate(gid,"NX_class", atttype, atts, H5P_DEFAULT);
H5Awrite(attid, atttype, (char *)"NXdata");
H5Sclose(atts);
H5Tclose(atttype);
H5Aclose(attid);

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/*
* store the counts dataset
*/
dataspace = H5Screate_simple(rank,dim,maxdim);
datatype = H5Tcopy(H5T_NATIVE_INT);
dataprop = H5Pcreate(H5P_DATASET_CREATE);
dataid = H5Dcreate(gid,(char *)"counts",datatype,dataspace,dataprop);
H5Dwrite(dataid, datatype, H5S_ALL, H5S_ALL, H5P_DEFAULT, counts);
H5Sclose(dataspace);
H5Tclose(datatype);
H5Pclose(dataprop);
/*
* set the signal=1 attribute
*/
atts = H5Screate(H5S_SCALAR);
atttype = H5Tcopy(H5T_NATIVE_INT);
H5Tset_size(atttype,1);
attid = H5Acreate(dataid,"signal", atttype, atts, H5P_DEFAULT);

2.2. Code Examples that do not use the NAPI

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```c
int signal = 1;
H5Awrite(attid, atttype, &signal);
H5Sclose(atts);
H5Tclose(atttype);
H5Aclose(attid);
H5Dclose(dataid);

/*
 * store the two_theta dataset
 */
dataspace = H5Screate_simple(rank, dim, maxdim);
datatype = H5Tcopy(H5T_NATIVE_FLOAT);
dataprop = H5Pcreate(H5P_DATASET_CREATE);
dataid = H5Dcreate(gid, (char *)"two_theta", datatype, dataspace, dataprop);
H5Dwrite(dataid, datatype, H5S_ALL, H5S_ALL, H5P_DEFAULT, two_theta);
H5Sclose(dataspace);
H5Tclose(datatype);
H5Pclose(dataprop);

/*
 * set the units attribute
 */
atttype = H5Tcopy(H5T_C_S1);
H5Tset_size(atttype, strlen("degrees"));
atts = H5Screate(H5S_SCALAR);
attid = H5Acreate(dataid, "units", atttype, atts, H5P_DEFAULT);
H5Awrite(attid, atttype, (char *)"degrees");
H5Sclose(atts);
H5Tclose(atttype);
H5Aclose(attid);

/*
 * set the target attribute for linking
 */
atttype = H5Tcopy(H5T_C_S1);
H5Tset_size(atttype, strlen("/scan/data/two_theta"));
atts = H5Screate(H5S_SCALAR);
attid = H5Acreate(dataid, "target", atttype, atts, H5P_DEFAULT);
H5Awrite(attid, atttype, (char *)"/scan/data/two_theta");
H5Sclose(atts);
H5Tclose(atttype);
H5Aclose(attid);

H5Dclose(dataid);

/*
 * make a link in /scan to /scan/data/two_theta, thereby renaming two_theta to hugo
 */
H5Glink(fid, H5G_LINK_HARD, "/scan/data/two_theta", "/scan/hugo");

/*
 * close the file
 */
H5Fclose(fid);
```

Chapter 2. Examples of writing and reading NeXus data files
Reading a simple NeXus file using native HDF5 commands in C

```c
/**
 * Reading example for reading NeXus files with plain
 * HDF-5 API calls. This reads out counts and two_theta
 * out of the file generated by nxh5write.
 * WARNING: I left out all error checking in this example.
 * In production code you have to take care of those errors
 * Mark Koennecke, October 2011
 */
#include <hdf5.h>
#include <stdlib.h>

int main(int argc, char *argv[])
{
    float *two_theta = NULL;
    int *counts = NULL, rank, i;
    hid_t fid, dataid, fapl;
    hsize_t *dim = NULL;
    hid_t datatype, dataspace, memdataspace;

    /* Open file, thereby enforcing proper file close
     * semantics
     */
    fapl = H5Pcreate(H5P_FILE_ACCESS);
    H5Pset_fclose_degree(fapl,H5F_CLOSE_STRONG);
    fid = H5Fopen("NXfile.h5", H5F_ACC_RDONLY,fapl);
    H5Pclose(fapl);

    /* open and read the counts dataset */
    dataid = H5Dopen(fid,"/scan/data/counts");
    dataspace = H5Dget_space(dataid);
    rank = H5Sget_simple_extent_ndims(dataspace);
    dim = malloc(rank*sizeof(hsize_t));
    H5Sget_simple_extent_dims(dataspace, dim, NULL);
    counts = malloc(dim[0]*sizeof(int));
    memdataspace = H5Tcopy(H5T_NATIVE_INT32);
    H5Dread(dataid,memdataspace,H5S_ALL, H5S_ALL,H5P_DEFAULT, counts);
    H5Dclose(dataid);
    H5Sclose(dataspace);
    H5Tclose(memdataspace);

    /* open and read the two_theta data set */
    dataid = H5Dopen(fid,"/scan/data/two_theta");
    dataspace = H5Dget_space(dataid);
    rank = H5Sget_simple_extent_ndims(dataspace);
    dim = malloc(rank*sizeof(hsize_t));
    H5Sget_simple_extent_dims(dataspace, dim, NULL);
    two_theta = malloc(dim[0]*sizeof(float));
    memdataspace = H5Tcopy(H5T_NATIVE_FLOAT);
    H5Dread(dataid,memdataspace,H5S_ALL, H5S_ALL,H5P_DEFAULT, two_theta);
}
```

2.2. Code Examples that do not use the NAPI
H5Dclose(dataid);
H5Sclose(dataspace);
H5Tclose(memdataspace);
H5Fclose(fid);

for (i = 0; i < dim[0]; i++) {
    printf("%8.2f %10d\n", two_theta[i], counts[i]);
}

2.2.2 Python Examples using h5py

One way to gain a quick familiarity with NeXus is to start working with some data. For at least the first few examples in this section, we have a simple two-column set of 1-D data, collected as part of a series of alignment scans by the APS USAXS instrument during the time it was stationed at beam line 32ID. We will show how to write this data using the Python language and the h5py package (using h5py calls directly rather than using the NeXus NAPI). The actual data to be written was extracted (elsewhere) from a spec data file and read as a text block from a file by the Python source code. Our examples will start with the simplest case and add only mild complexity with each new case since these examples are meant for those who are unfamiliar with NeXus.

The data shown plotted in the next figure will be written to the NeXus HDF5 file using the only two required NeXus objects NXentry and NXdata in the first example and then minor variations on this structure in the next two examples. The data model is identical to the one in the Introduction chapter except that the names will be different, as shown below:

![data structure diagram]

Figure 2.1: data structure, (from Introduction)

our h5py example

```
/entry:NXentry
  /mr_scan:NXdata
    /mr : float64[31]
    /I00 : int32[31]
```

1. h5py: http://code.google.com/p/h5py
2. SPEC: http://certif.com/spec.html
2.2. Code Examples that do not use the NAPI
Writing the simplest data using h5py

These two examples show how to write the simplest data (above). One example writes the data directly to the NXdata group while the other example writes the data to NXinstrument/NXdetector/data and then creates a soft link to that data in NXdata.

h5py example writing the simplest NeXus data file

In this example, the 1-D scan data will be written into the simplest possible NeXus HDF5 data file, containing only the required NeXus components. NeXus requires at least one NXentry group at the root level of an HDF5 file. The NXentry group contains all the data and associated information that comprise a single measurement. NeXus also requires that each NXentry group must contain at least one NXdata group. NXdata is used to describe the plottable data in the NXentry group. The simplest place to store data in a NeXus file is directly in the NXdata group, as shown in the next figure.

In the above figure, the data file (writer_1_3_h5py.hdf5) contains a hierarchy of items, starting with an NXentry named entry. (The full HDF5 path reference, /entry in this case, is shown to the right of each component in the data structure.) The next h5py code example will show how to build an HDF5 data file with this structure. Starting with the numerical data described above, the only information written to the file is the absolute minimum information NeXus requires. In this example, you can see how the HDF5 file is created, how Data Groups and datasets (Data Fields) are created, and how Data Attributes are assigned. Note particularly the NX_class attribute on each HDF5 group that describes which of the NeXus Base Class Definitions is being used. When the next Python program (writer_1_3_h5py.py) is run from the command line (and there are no problems), the writer_1_3_h5py.hdf5 file is generated.

```python
#!/usr/bin/env python
'''
Writes the simplest NeXus HDF5 file using h5py
according to the example from Figure 1.3
in the Introduction chapter
'''
```
import h5py
import numpy

INPUT_FILE = 'input.dat'
HDF5_FILE = 'writer_1_3_h5py.hdf5'

#---------------------------
tthData, countsData = numpy.loadtxt(INPUT_FILE).T
f = h5py.File(HDF5_FILE, "w")  # create the HDF5 NeXus file
# since this is a simple example, no attributes are used at this point
nxentry = f.create_group('Scan')
nxentry.attrs['NX_class'] = 'NXentry'
nndata = nxentry.create_group('data')
nndata.attrs['NX_class'] = 'NXdata'
tth = nndata.create_dataset("two_theta", data=tthData)
tth.attrs['units'] = "degrees"

counts = nndata.create_dataset("counts", data=countsData)
counts.attrs['units'] = "counts"
counts.attrs['signal'] = 1

f.close()  # be CERTAIN to close the file

We wish to make things a bit simpler for ourselves when creating the common structures we use in our data files. To help, we gather together some of the common concepts such as create a file, create a NeXus group, create a dataset and start to build a helper library. (See mylib support module for more details.) Here, we call it my_lib. Applying it to the simple example above, our code only becomes a couple lines shorter! (Let’s hope the library starts to help in larger or more complicated projects.) Here’s the revision that replaces direct calls to numpy and h5py with calls to our library. It generates the file writer_1_3.hdf5.

#!/usr/bin/env python

"""
Writes the simplest NeXus HDF5 file using a simple helper library with h5py and numpy calls according to the example from Figure 1.3 in the Introduction chapter
"""

import my_lib

INPUT_FILE = 'input.dat'
HDF5_FILE = 'writer_1_3.hdf5'

#---------------------------
tthData, countsData = my_lib.get2ColumnData(INPUT_FILE)
f = my_lib.makeFile(HDF5_FILE)
# since this is a simple example, no attributes are used at this point
nxentry = my_lib.makeGroup(f, 'Scan', 'NXentry')
nndata = my_lib.makeGroup(nxentry, 'data', 'NXdata')

2.2. Code Examples that do not use the NAPI
```python
my_lib.makeDataset(nxdata, "two_theta", tthData, units='degrees')
my_lib.makeDataset(nxdata, "counts", countsData,
    units='counts', signal=1, axes='two_theta')
f.close()  # be CERTAIN to close the file
```

One of the tools provided with the HDF5 support libraries is the h5dump command, a command-line tool to print out the contents of an HDF5 data file. With no better tool in place (the output is verbose), this is a good tool to investigate what has been written to the HDF5 file. View this output from the command line using h5dump writer_1_3.hdf5. Compare the data contents with the numbers shown above. Note that the various HDF5 data types have all been decided by the h5py support package.

**Note:** The only difference between this file and one written using the NAPI is that the NAPI file will have some additional, optional attributes set at the root level of the file that tells the original file name, time it was written, and some version information about the software involved.
41 STRPAD H5T_STR_NULLPAD;
42 CSET H5T_CSET_ASCII;
43 CType H5T_C_S1;
44 }
45 DATASPACE SCALAR
46 DATA {
47 (0): "counts"
48 }
49 }
50 ATTRIBUTE "signal" {
51 DATATYPE H5T_STRING {
52 STRSIZE 1;
53 STRPAD H5T_STR_NULLPAD;
54 CSET H5T_CSET_ASCII;
55 CType H5T_C_S1;
56 }
57 DATASPACE SCALAR
58 DATA {
59 (0): "1"
60 }
61 }
62 ATTRIBUTE "axes" {
63 DATATYPE H5T_STRING {
64 STRSIZE 9;
65 STRPAD H5T_STR_NULLPAD;
66 CSET H5T_CSET_ASCII;
67 CType H5T_C_S1;
68 }
69 DATASPACE SCALAR
70 DATA {
71 (0): "two_theta"
72 }
73 }
74 }
75 DATASET "two_theta" {
76 DATATYPE H5T_IEEE_F64LE
77 DATASPACE SIMPLE { (31) / (31) }
78 DATA {
79 (0): 17.9261, 17.9259, 17.9258, 17.9256, 17.9254, 17.9252,
80 (6): 17.9251, 17.9249, 17.9247, 17.9246, 17.9244, 17.9243,
81 (12): 17.9241, 17.9239, 17.9237, 17.9236, 17.9234, 17.9232,
82 (18): 17.9231, 17.9229, 17.9228, 17.9226, 17.9224, 17.9222,
83 (24): 17.9221, 17.9219, 17.9217, 17.9216, 17.9214, 17.9213,
84 (30): 17.9211
85 }
86 ATTRIBUTE "units" {
87 DATATYPE H5T_STRING {
88 STRSIZE 7;
89 STRPAD H5T_STR_NULLPAD;
90 CSET H5T_CSET_ASCII;
91 CType H5T_C_S1;
92 }
93 DATASPACE SCALAR
94 DATA {
95 (0): "degrees"
96 }
97 }
98 }
Since the output of h5dump is verbose, a tool (see h5toText support module) was created to print out the structure of HDF5 data files. This tool provides a simplified view of the NeXus file. It is run with a command like this: python h5toText.py h5dump writer_1_3.hdf5. Here is the output:

writer_1_3.hdf5:NeXus data file
  Scan: NXentry
    @NX_class = NXentry
  data: NXdata
    @NX_class = NXdata
    counts: NX_INT32[31] = __array
      @units = counts
      @signal = 1
      @axes = two_theta
        __array = [1037, 1318, 1704, ... , 1321]
  two_theta: NX_FLOAT64[31] = __array
    @units = degrees
    __array = [17.926079999999999, 17.925909999999998, 17.925750000000001, ... , 17.92108]

As the data files in these examples become more complex, you will appreciate the information density provided by the h5toText.py tool.

h5py example writing a simple NeXus data file with links

Building on the previous example, we wish to identify our measured data with the detector on the instrument where it was generated. In this hypothetical case, since the detector was positioned at some angle two_theta, we choose to store both datasets, two_theta and counts, in a NeXus group. One appropriate NeXus group is NXdetector. This group is placed in a NXinstrument group which is placed in a NXentry group. Still, NeXus requires a NXdata group. Rather than duplicate the same data already placed in the detector group, we choose to link to those datasets from the NXdata group. (Compare the next figure with Linking in a NeXus file in the NeXus Design chapter of the NeXus User Manual.) The NeXus Design chapter provides a figure (Linking in a NeXus file) with a small variation from our previous example, placing the measured data within the /entry/instrument/detector group. Links are made from that data to the /entry/data group.

![Figure 2.4: h5py example showing linking in a NeXus file](image)

Figure 2.4: h5py example showing linking in a NeXus file

The Python code to build an HDF5 data file with that structure (using numerical data from the previous example) is
shown below.

```python
#!/usr/bin/env python
'''
Writes a simple NeXus HDF5 file using h5py with links
according to the example from Figure 2.1 in the Design chapter
'''

import my_lib

INPUT_FILE = 'input.dat'
HDF5_FILE = 'writer_2_1.hdf5'

#---------------------------
tthData, countsData = my_lib.get2ColumnData(INPUT_FILE)
f = my_lib.makeFile(HDF5_FILE)  # create the HDF5 NeXus file

nxentry = my_lib.makeGroup(f, 'entry', 'NXentry')
nxinstrument = my_lib.makeGroup(nxentry, 'instrument', 'NXinstrument')
nxdetector = my_lib.makeGroup(nxinstrument, 'detector', 'NXdetector')
tth = my_lib.makeDataset(nxdetector, "two_theta", tthData, units='degrees')
counts = my_lib.makeDataset(nxdetector, "counts", countsData,
    units='counts', signal=1, axes='two_theta')

nxdata = my_lib.makeGroup(nxentry, 'data', 'NXdata')
my_lib.makeLink(nxdetector, tth, nxdata.name+'/two_theta')
my_lib.makeLink(nxdetector, counts, nxdata.name+'/counts')

f.close()  # be CERTAIN to close the file
```

It is interesting to compare the output of the h5dump of the data file writer_2_1.hdf5 with our Python instructions.

```bash
HDF5 "writer_2_1.hdf5" {
GROUP "/" {
    GROUP "entry" {
        ATTRIBUTE "NX_class" {
            DATATYPE H5T_STRING {
                STRSIZE 7;
                STRPAD H5T_STR_NULLPAD;
                CSET H5T_CSET_ASCII;
                CTYPE H5T_C_S1;
            }
            DATASPACE SCALAR
            DATA {
                (0): "NXentry"
            }
        }
    }
    GROUP "data" {
        ATTRIBUTE "NX_class" {
            DATATYPE H5T_STRING {
                STRSIZE 6;
                STRPAD H5T_STR_NULLPAD;
                CSET H5T_CSET_ASCII;
                CTYPE H5T_C_S1;
            }
        }
    }
}
```
DATASPACE SCALAR
DATA {
(0): "NXdata"
}

DATASET "counts" {
DATATYPE H5T_STD_I32LE
DATASPACE SIMPLE { ( 31 ) / ( 31 ) }
DATA {
(0): 1037, 1318, 1704, 2857, 4516, 9998, 23819, 31662, 40458,
(9): 49087, 56514, 63499, 66802, 66863, 66599, 66206, 65747,
(17): 65250, 64129, 63044, 60796, 56795, 51550, 43710, 29315,
(25): 19782, 12992, 6622, 4198, 2248, 1321
}
ATTRIBUTE "units" {
DATATYPE H5T_STRING {
STRSIZE 6;
STRPAD H5T_STR_NULLPAD;
CSET H5T_CSET_ASCII;
CTYPE H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "counts"
}
}

ATTRIBUTE "signal" {
DATATYPE H5T_STRING {
STRSIZE 1;
STRPAD H5T_STR_NULLPAD;
CSET H5T_CSET_ASCII;
CTYPE H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "1"
}
}

ATTRIBUTE "axes" {
DATATYPE H5T_STRING {
STRSIZE 9;
STRPAD H5T_STR_NULLPAD;
CSET H5T_CSET_ASCII;
CTYPE H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "two_theta"
}
}

ATTRIBUTE "target" {
DATATYPE H5T_STRING {
STRSIZE 33;
STRPAD H5T_STR_NULLPAD;
CSET H5T_CSET_ASCII;
CTYPE H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "/entry/instrument/detector/counts"
}

DATASET "two_theta" {
DATATYPE H5T_IEEE_F64LE
DATASPACE SIMPLE { ( 31 ) / ( 31 ) }
DATA {
(0): 17.9261, 17.9259, 17.9258, 17.9256, 17.9254, 17.9252,
(6): 17.9251, 17.9249, 17.9247, 17.9246, 17.9244, 17.9243,
(12): 17.9241, 17.9239, 17.9237, 17.9236, 17.9234, 17.9232,
(18): 17.9231, 17.9229, 17.9228, 17.9226, 17.9224, 17.9222,
(24): 17.9221, 17.9219, 17.9217, 17.9216, 17.9214, 17.9213,
(30): 17.9211
}
ATTRIBUTE "units" {
DATATYPE H5T_STRING {
    STRSIZE 7;
    STRPAD H5T_STR_NULLPAD;
    CSET H5T_CSET_ASCII;
    CType H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "degrees"
}
}
ATTRIBUTE "target" {
DATATYPE H5T_STRING {
    STRSIZE 36;
    STRPAD H5T_STR_NULLPAD;
    CSET H5T_CSET_ASCII;
    CType H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "/entry/instrument/detector/two_theta"
}
}

GROUP "instrument" {
ATTRIBUTE "NX_class" {
DATATYPE H5T_STRING {
    STRSIZE 12;
    STRPAD H5T_STR_NULLPAD;
    CSET H5T_CSET_ASCII;
    CType H5T_C_S1;
}
DATASPACE SCALAR
DATA {
(0): "NXinstrument"
}
}
GROUP "detector" {
ATTRIBUTE "NX_class" {
DATATYPE H5T_STRING {

Look carefully! It appears from the output of h5dump that the actual data for two_theta and counts has moved into the NXdata group at HDF5 path /entry/data! But we stored that data in the NXdetector group at /entry/instrument/detector. This is normal for h5dump output.

A bit of explanation is necessary at this point. The data is not stored in either HDF5 group directly. Instead, HDF5 creates a DATA storage element in the file and posts a reference to that DATA storage element as needed. An HDF5 hard link requests another reference to that same DATA storage element. The h5dump tool describes in full that DATA storage element the first time (alphabetically) it is called. In our case, that is within the NXdata group. The next time it is called, within the NXdetector group, h5dump reports that a hard link has been made and shows the HDF5 path to the description.

NeXus recognizes this behavior of the HDF5 library and adds an additional structure when building hard links, the target attribute, to preserve the original location of the data. Not that it actually matters. The h5toText.py tool knows about the additional NeXus target attribute and shows the data to appear in its original location, in the NXdetector group.

```
writer_2_1.hdf5:NeXus data file
entry:NXentry
  @NX_class = NXentry
data:NXdata
    @NX_class = NXdata
    counts --> /entry/instrument/detector/counts
two_theta --> /entry/instrument/detector/two_theta
instrument:NXinstrument
  @NX_class = NXinstrument
detector:NXdetector
    @NX_class = NXdetector
counts:NX_INT32[31] = __array
      @units = counts
      @signal = 1
      @axes = two_theta
      @target = /entry/instrument/detector/counts
      __array = [1037, 1318, 1704, ..., 1321]
two_theta:NX_FLOAT64[31] = __array
      @units = degrees
      @target = /entry/instrument/detector/two_theta
```
Complete `h5py` example writing and reading a NeXus data file

Writing the HDF5 file using `h5py`

In the main code section of `BasicWriter.py`, a current time stamp is written in the format of ISO 8601. For simplicity of this code example, we use a text string for the time, rather than computing it directly from Python support library calls. It is easier this way to see the exact type of string formatting for the time. When using the Python `datetime` package, one way to write the time stamp is:

```python
timestamp = 'T'.join(str(datetime.datetime.now()).split())
```

The data (mr is similar to “two_theta” and I00 is similar to “counts”) is collated into two Python lists. We use our `my_lib` support to read the file and parse the two-column format.

The new HDF5 file is opened (and created if not already existing) for writing, setting common NeXus attributes in the same command from our support library. Proper HDF5+NeXus groups are created for `/entry:NXentry/mr_scan:NXdata`. Since we are not using the NAPI, our support library must create and set the NX_class attribute on each group.

Note: We want to create the desired structure of `/entry:NXentry/mr_scan:NXdata/`. First, our support library calls `nxentry = f.create_group("entry")` to create the NXentry group called `entry` at the root level. Then, it calls `nxdata = nxentry.create_group("mr_scan")` to create the NXentry group called `entry` as a child of the NXentry group.

Next, we create a dataset called `title` to hold a title string that can appear on the default plot.

Next, we create datasets for `mr` and `I00` using our support library. The data type of each, as represented in `numpy`, will be recognized by `h5py` and automatically converted to the proper HDF5 type in the file. A Python dictionary of attributes is given, specifying the engineering units and other values needed by NeXus to provide a default plot of this data. By setting `signal=1` as an attribute on `I00`, NeXus recognizes `I00` as the default y axis for the plot. The `axes="mr"` connects the dataset to be used as the x axis.

Finally, we must remember to call `f.close()` or we might corrupt the file when the program quits.

### `BasicWriter.py`: Write a NeXus HDF5 file using Python with `h5py`

```python
#!/usr/bin/env python
'''Writes a NeXus HDF5 file using h5py and numpy'''

import h5py  # HDF5 support
import numpy
import my_lib  # uses h5py

print "Write a NeXus HDF5 file using h5py and numpy"

print "Write a NeXus HDF5 file using h5py and numpy"

file_name = "prj_test.nexus.hdf5"
timestamp = "2010-10-18T17:17:04-0500"

# load data from two column format
data = numpy.loadtxt('input.dat').T
mr_arr = data[0]
I00_arr = numpy.asarray(data[1], 'int32')
```
# create the HDF5 NeXus file
f = my_lib.makeFile(fileName, file_name=fileName,
    file_time=timestamp,
    instrument="APS USAXS at 32ID-B",
    creator="BasicWriter.py",
    NeXus_version="4.3.0",
    HDF5_Version=h5py.version.hdf5_version,
    h5py_version=h5py.version.version)

nxentry = my_lib.makeGroup(f, "entry", "NXentry")
my_lib.makeDataset(nxentry, 'title', data='1-D scan of I00 v. mr')

nxdata = my_lib.makeGroup(nxentry, "mr_scan", "NXdata")
my_lib.makeDataset(nxdata, "mr", mr_arr, units='degrees',
    long_name='USAXS mr (degrees)')
my_lib.makeDataset(nxdata, "I00", i00_arr, units='counts',
    signal=1, # Y axis of default plot
    axes='mr', # name "mr" as X axis
    long_name='USAXS I00 (counts)')

f.close() # be CERTAIN to close the file

print "wrote file:", fileName

Reading the HDF5 file using h5py

The file reader, BasicReader.py, is very simple since the bulk of the work is done by h5py. Our code opens the HDF5
we wrote above, prints the HDF5 attributes from the file, reads the two datasets, and then prints them out as columns.
As simple as that. Of course, real code might add some error-handling and extracting other useful stuff from the file.

Note: See that we identified each of the two datasets using HDF5 absolute path references (just using the group and
dataset names). Also, while coding this example, we were reminded that HDF5 is sensitive to upper or lowercase.
That is, I00 is not the same is i00.

BasicReader.py: Read a NeXus HDF5 file using Python with h5py

#!/usr/bin/env python
'''Reads NeXus HDF5 files using h5py and prints the contents'''
import h5py # HDF5 support

fileName = "prj_test.nexus.hdf5"
f = h5py.File(fileName, "r")
for item in f.attrs.keys():
    print item + ":", f.attrs[item]
mr = f["/entry/mr_scan/mr"]
i00 = f["/entry/mr_scan/I00"]
for i in range(len(mr)):
    print "%d %g %d" % (i, mr[i], i00[i])
f.close()
Output from BasicReader.py is shown next.

```
file_name: prj_test.nexus.hdf5
file_time: 2010-10-18T17:17:04-0500
creator: BasicWriter.py
HDF5_Version: 1.8.5
NeXus_version: 4.3.0
h5py_version: 1.2.1
instrument: APS USAXS at 32ID-B
#   mr  I00
 0 17.9261  1037
 1 17.9259  1318
 2 17.9258  1704
 3 17.9256  2857
 4 17.9254  4516
 5 17.9252  9998
 6 17.9251 23819
 7 17.9249 31662
 8 17.9247 40458
 9 17.9246 49087
10 17.9244 56514
11 17.9243 63499
12 17.9241 66802
13 17.9239 66863
14 17.9237 66599
15 17.9236 66206
16 17.9234 65747
17 17.9232 65250
18 17.9231 64129
19 17.9229 63044
20 17.9228 60796
21 17.9226 56795
22 17.9224 51550
23 17.9222 43710
24 17.9221 29315
25 17.9219 19782
26 17.9217 12992
27 17.9216 6622
28 17.9214 4198
29 17.9213 2248
30 17.9211 1321
```

Validating the HDF5 file

Now we have an HDF5 file that contains our data. What makes this different from a NeXus data file? A NeXus file has a specific arrangement of groups and datasets in an HDF5 file.

To test that our HDF5 file conforms to the NeXus standard, we use the Java-version of NXvalidate. Referring to the next figure, we compare our HDF5 file with the rules for generic data files (all.nxdl.xml). The only items that have been flagged are the “non-standard field names” mr and I00. Neither of these two names is specifically named in the NeXus NXDL definition for the NXdata base class. As we’ll see shortly, this is not a problem.

---

3 generic NeXus data files: NeXus data files for which no application-specific NXDL applies
Figure 2.5: NeXus validation of our HDF5 file

Note: Note that NXvalidate shows only the first data field for \(mr\) and \(I00\).

Plotting the HDF5 file

Now that we are certain our file conforms to the NeXus standard, let’s plot it using the NeXpy client tool. To help label the plot, we added the long_name attributes to each of our datasets. We also added metadata to the root level of our HDF5 file similar to that written by the NAPI. It seemed to be a useful addition. Compare this with plot of our \(mr\) scan and note that the horizontal axis of this plot is mirrored from that above. This is because the data is stored in the file in descending \(mr\) order and NeXpy has plotted it that way by default.

Links to Data in External HDF5 Files

HDF5 files may contain links to data (or groups) in other files. This can be used to advantage to refer to data in existing HDF5 files and create NeXus-compliant data files. Here, we show such an example, using the same counts vs two_theta data from the examples above.

file: external_angles.hdf5

Take for example, the structure of external_angles.hdf5, a simple HDF5 data file that contains just the two_theta angles in an HDF5 dataset at the root level of the file. Although this is a valid HDF5 data file, it is not a valid NeXus data file:

\[\text{file: external_angles.hdf5} \]

\[\text{Take for example, the structure of external_angles.hdf5, a simple HDF5 data file that contains just the two_theta angles in an HDF5 dataset at the root level of the file. Although this is a valid HDF5 data file, it is not a valid NeXus data file.} \]

\[\text{NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1} \]

\[\text{Note that NXvalidate shows only the first data field for mr and I00.} \]
Figure 2.6: plot of our mr_scan using NeXpy

```python
angles: float64[31] = [17.926079999999999, ..., 17.92108]
@units = degrees
```

file: external_counts.hdf5

The data in the file external_angles.hdf5 might be referenced from another HDF5 file (such as external_counts.hdf5) by an HDF5 external link. Here is an example of the structure:

```python
entry: NXentry
  instrument: NXinstrument
  detector: NXdetector
  counts: NX_INT32[31] = [1037, ..., 1321]
    @units = counts
    @signal = 1
    @axes = two_theta
  two_theta --> file="external_angles.hdf5", path="/angles"
```

Note: The file external_counts.hdf5 is not a complete NeXus file since it does not contain an NXdata group containing a dataset with signal=1 attribute.

file: external_master.hdf5

A valid NeXus data file could be created that refers to the data in these files without making a copy of the data files themselves.

---

5 see these URLs for further guidance on HDF5 external links: [http://www.hdfgroup.org/HDF5/doc/2.5.0/RM/RM_H5L.html#Link-CreateExternal](http://www.hdfgroup.org/HDF5/doc/2.5.0/RM/RM_H5L.html#Link-CreateExternal), [http://www.h5py.org/docs-1.3/guide/group.html#external-links](http://www.h5py.org/docs-1.3/guide/group.html#external-links)

2.2. Code Examples that do not use the NAPI
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

Note: It is necessary for all these files to be located together in the same directory for the HDF5 external file links to work properly.

To be a valid NeXus file, it must contain a `NXentry` group containing a `NXdata` group containing only one dataset with the attribute `signal=1`. For the files above, it is simple to make a master file that links to the data we desire, from structure that we create. In `external_counts.hdf5` above, see that the required attribute `signal=1` is already present. Here is `external_master.hdf5`, an example:

```
entry: NXEntry
  instrument --> file="external_counts.hdf5", path="/entry/instrument"
data: NXdata
  counts --> file="external_counts.hdf5", path="/entry/instrument/detector/counts"
two_theta --> file="external_angles.hdf5", path="/angles"
```

source code: externalExample.py

Here is the complete code of a Python program, using `h5py` to write a NeXus-compliant HDF5 file with links to data in other HDF5 files.

```
#!/usr/bin/env python

# Import necessary libraries
import my_lib

# Define input file names
FILE_INPUT = 'input.dat'
FILE_HDF5_MASTER = 'external_master.hdf5'
FILE_HDF5_ANGLES = 'external_angles.hdf5'
FILE_HDF5_COUNTS = 'external_counts.hdf5'

# Get some data
tht_data, counts_data = my_lib.get2ColumnData(FILE_INPUT)

# Put the angle data in an external (non-NeXus) HDF5 data file
f = my_lib.makeFile(FILE_HDF5_ANGLES)
angles = my_lib.makeDataset(f, 'angles', tht_data, units='degrees')
f.close()  # Be CERTAIN to close the file

# Put the detector counts in an external NeXus HDF5 data file
f = my_lib.makeFile(FILE_HDF5_COUNTS)
entry = my_lib.makeGroup(f, 'entry', 'NXEntry')
instrument = my_lib.makeGroup(entry, 'instrument', 'NXInstrument')
detector = my_lib.makeGroup(instrument, 'detector', 'NXDetector')
counts = my_lib.makeDataset(detector, 'counts', counts_data, units='counts', signal=1, axes='two_theta')

# Make a link since "two_theta" has not been stored here
my_lib.makeExternalLink(f, FILE_HDF5_ANGLES,
```

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35
36 f.close()
37
38 # create a master NeXus HDF5 file
39 f = my_lib.makeFile(FILE_HDF5_MASTER)
40 nxentry = my_lib.makeGroup(f, 'entry', 'NXentry')
41 nxdata = my_lib.makeGroup(nxentry, 'data', 'NXdata')
42 my_lib.makeExternalLink(f, FILE_HDF5_ANGLES,
43 'angles', nxdata.name+'/two_theta')
44 my_lib.makeExternalLink(f, FILE_HDF5_COUNTS,
45 '/entry/instrument/detector/counts',
46 nxdata.name+'/counts')
47 my_lib.makeExternalLink(f, FILE_HDF5_COUNTS,
48 '/entry/instrument',
49 nxentry.name+'/instrument')
50 f.close()

Python Helper Modules for h5py Examples

Two additional Python modules were used to describe these h5py examples. The source code for each is given here. The first is a library we wrote that helps us create standard NeXus components using h5py. The second is a tool that helps us inspect the content and structure of HDF5 files.

mylib support module

The examples in this section make use of a small helper library that calls h5py to create the various NeXus data components of Data Groups, Data Fields, Data Attributes, and Links. In a smaller sense, this subroutine library (mylib) fills the role of the NAPI for writing the data using h5py.

```python
#!/usr/bin/env python
'''
mylib: routines to support reading & writing NeXus HDF5 files using h5py
'''

import h5py  # HDF5 support
import numpy  # in this case, provides data structures

def makeFile(filename, **attr):
    """
    create and open an empty NeXus HDF5 file using h5py
    ""
    Any named parameters in the call to this method will be saved as attributes of the root of the file.
    Note that **attr is a dictionary of named parameters.
    :param str filename: valid file name
    :param attr: optional keywords of attributes
    :return: h5py file object
    ""
    obj = h5py.File(filename, "w")
    addAttributes(obj, attr)
    return obj

def makeGroup(parent, name, nxclass, **attr):
    """
```
create a NeXus group

Any named parameters in the call to this method
will be saved as attributes of the group.
Note that **attr is a dictionary of named parameters.

:param obj parent: parent group
:param str name: valid NeXus group name
:param str nxclass: valid NeXus class name
:param attr: optional keywords of attributes
:return: h5py group object

```python
obj = parent.create_group(name)
obj.attrs["NX_class"] = nxclass
addAttributes(obj, attr)
return obj
```

def makeDataset(parent, name, data = None, **attr):
    '''
    create and write data to a dataset in the HDF5 file hierarchy
    Any named parameters in the call to this method
    will be saved as attributes of the dataset.
    '''
    if data == None:
        obj = parent.create_dataset(name)
    else:
        obj = parent.create_dataset(name, data=data)
    addAttributes(obj, attr)
    return obj

def makeLink(parent, sourceObject, targetName):
    '''
    create an internal NeXus (hard) link in an HDF5 file
    '''
    if not 'target' in sourceObject.attrs:
        # NeXus link, NOT an HDF5 link!
        sourceObject.attrs["target"] = str(sourceObject.name)
    parent._id.link(sourceObject.name, targetName, h5py.h5g.LINK_HARD)

def makeExternalLink(hdf5FileObject, sourceFile, sourcePath, targetPath):
    '''
    create an external link from sourceFile, sourcePath to targetPath in hdf5FileObject
    '''
    if not os.path.isabs(sourcePath):
        sourcePath = os.path.join(os.path.dirname(sourceFile), sourcePath)
    hdf5FileObject._id.link(sourceFile, sourcePath, targetPath, h5py.h5g.LINK_HARD)
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.. note::
   Since the object retrieved is in a different file, its ".file" and ".parent" properties will refer to objects in that file, not the file in which the link resides.

.. see:: http://www.h5py.org/docs-1.3/guide/group.html#external-links

This routine is provided as a reminder how to do this simple operation.

```python
hdf5FileObject[targetPath] = h5py.ExternalLink(sourceFile, sourcePath)
```

```python
def addAttributes(parent, attr):
    
    add attributes to an h5py data item
    
    :param obj parent: h5py parent object
    :param dict attr: dictionary of attributes
    
    if attr and type(attr) == type({}):  # attr is a dictionary of attributes
        for k, v in attr.items():
            parent.attrs[k] = v
```

```python
def get2ColumnData(fileName):
    
    read two-column data from a file, first column is float, second column is integer
    
    buffer = numpy.loadtxt(fileName).T
    xArr = buffer[0]
    yArr = numpy.asarray(buffer[1], 'int32')
    return xArr, yArr
```

h5toText support module

The module h5toText reads an HDF5 data file and prints out the structure of the groups, datasets, attributes, and links in that file. There is a command-line option to print out more or less of the data in the dataset arrays.

```python
#!/usr/bin/env python

# Print the structure of an HDF5 file to stdout

import h5py
import os
import sys
import getopt

class H5toText(object):
    
    ```
Example usage showing default display:

```python
mc = H5toText(filename)
mc.array_items_shown = 5
mc.report()
'''
filename = None
requested_filename = None
isNeXus = False
array_items_shown = 5

def __init__(self, filename, makeReport = False):
    ''' Constructor '''
    self.requested_filename = filename
    if os.path.exists(filename):
        self.filename = filename
        self.isNeXus = self.testIsNeXus()
        if makeReport:
            self.report()

def report(self):
    ''' reporter '''
    if self.filename == None: return
    f = h5py.File(self.filename, 'r')
    txt = self.filename
    if self.isNeXus:
        txt += ':NeXus data file'
    self.showGroup(f, txt, indentation = '')
    f.close()

def testIsNeXus(self):
    '''
test if the selected HDF5 file is a NeXus file

    At this time, the code only tests for the existence of
    the NXentry group. The tests should be extended to require
    a NXdata group and a single dataset containing signal=1 attribute.
    '''
    result = False
    try:
        f = h5py.File(self.filename, 'r')
        for value in f.itervalues():
            #print str(type(value))
            if '.Group' not in str(type(value)):
                continue
            #print value.attrs.keys()
            if 'NX_class' not in value.attrs:
                continue
            v = value.attrs['NX_class']
            #print type(v), v, type("a string")
            possible_types = [{"<type 'numpy.string_'>", },
            possible_types.append("<type 'str'>")
            if str(type(v)) not in possible_types:
                continue
            if str(v) == str('NXentry'):
                # TODO: apply more tests
                # for group NXdata
                # and signal=1 attribute on only one dataset
```
result = True
break
f.close()
except:
    pass
return result

def showGroup(self, obj, name, indentation = " "):  
    '''print the contents of the group'''
    nxclass = ""
    if 'NX_class' in obj.attrs:
        class_attr = obj.attrs['NX_class']
        nxclass = ":" + str(class_attr)
    print indentation + name + nxclass
    self.showAttributes(obj, indentation)
    # show datasets and links next
groups = []
    for itemname in sorted(obj):
        linkref = obj.get(itemname, getlink=True)
        if '.ExternalLink' in str(type(linkref)):
            # if the external file is not present, cannot know if
            # link target is a dataset or a group or another link
            fmt = '%s %s --> file="%s", path="%s"'
        else:
            classref = obj.get(itemname, getclass=True)
            value = obj.get(itemname)
        if '.File' in str(classref) or '.Group' in str(classref):
            groups.append(value)
        elif '.Dataset' in str(classref):
            self.showDataset(value, itemname, indentation +"
        else:
            msg = "unidentified $s: $s, $s"
            itemname, repr(classref), repr(linkref)
            raise Exception, msg
    # then show things that look like groups
    for value in groups:
        itemname = value.name.split('/')[1]
        self.showGroup(value, itemname, indentation +"

def showAttributes(self, obj, indentation = " "):  
    '''print any attributes'''
    for name, value in obj.attrs.iteritems():
        print "$s $s = %s" % (indentation, name, str(value))

def showDataset(self, dset, name, indentation = " "):  
    '''print the contents and structure of a dataset'''
    shape = dset.shape
    if self.isNeXus:
        if "target" in dset.attrs:
            if dset.attrs['target'] != dset.name:
                print "$s$ = %s" % (indentation, name, 
                dset.attrs['target'])
        return
txType = self.getType(dset)
txShape = self.getShape(dset)
    if shape == (1,):
        value = "%s = %s" % str(dset[0])
        print "$s$ = %s" % (indentation, name, txType,
def getType(self, obj):
    ''' get the storage (data) type of the dataset '''
    t = str(obj.dtype)
    if t[0:2] == '|S':
        t = 'char[' + t[2:]
    if self.isNeXus:
        t = 'NX_' + t.upper()
    return t

def getShape(self, obj):
    ''' return the shape of the HDF5 dataset '''
    s = obj.shape
    l = []
    for dim in s:
        l.append(str(dim))
    if l == ['1']:
        result = """  
    else:
        result = "[" + ','.join(l) + "]"
    return result

def formatArray(self, obj, indentation = ' '):
    ''' nicely format an array up to rank=5 '''
    shape = obj.shape
    r = ""
    if len(shape) in (1, 2, 3, 4, 5):
        r = self.formatNdArray(obj, indentation + ' ')
    if len(shape) > 5:
        r = "### no arrays for rank > 5 ###"
    return r

def decideNumShown(self, n):
    ''' determine how many values to show '''
    if self.array_items_shown != None:
        if n > self.array_items_shown:
            n = self.array_items_shown - 2
        return n

def formatNdArray(self, obj, indentation = ' '):
    ''' return a list of lower-dimension arrays, nicely formatted '''
    shape = obj.shape
    rank = len(shape)
    if not rank in (1, 2, 3, 4, 5): return None
    n = self.decideNumShown(shape[0])
    r = []

def showAttributes(dset, indentation = ' '):
    txShape = dset.attrs['nv_shape']
    txType = dset.attrs['nv_type']
    txValue = dset.attrs['nv_value']
    print "%s%s = %s" % (indentation, name, value)
for i in range(n):
    if rank == 1: item = obj[i]
    if rank == 2: item = self.formatNdArray(obj[i, :])
    if rank == 3: item = self.formatNdArray(obj[i, :, :],
                                                    indentation + ' ')
    if rank == 4: item = self.formatNdArray(obj[i, :, :, :],
                                                    indentation + ' ')
    if rank == 5: item = self.formatNdArray(obj[i, :, :, :, :],
                                                    indentation + ' ')
    r.append( item )
    if n < shape[0]:
        # skip over most
        r.append("...")
        # get the last one
    if rank == 1: item = obj[-1]
    if rank == 2: item = self.formatNdArray(obj[-1, :])
    if rank == 3: item = self.formatNdArray(obj[-1, :, :],
                                                    indentation + ' ')
    if rank == 4: item = self.formatNdArray(obj[-1, :, :, :],
                                                    indentation + ' ')
    if rank == 5: item = self.formatNdArray(obj[-1, :, :, :, :],
                                                    indentation + ' ')
    r.append( item )

    if rank == 1:
        s = str( r )
    else:
        s = "\n" + indentation + ' '
        s += ("\n" + indentation + ' ').join(r)
        s += "\n" + indentation + ""
    return s

def do_filelist(filelist, limit=5):
    '''
    interpret the structure of a list of HDF5 files
    :
    :param [str] filelist: one or more file names to be interpreted
    :param int limit: maximum number of array items to be shown (default = 5)
    '''
    for item in filelist:
        mc = H5toText(item)
        mc.array_items_shown = limit
        mc.report()

def do_test():
    limit = 3
    filelist = []
    filelist.append(‘th02c_ps02_1_master.h5’)
    filelist.append(‘external_angles.hdf5’)
    filelist.append(‘external_counts.hdf5’)
    filelist.append(‘external_master.hdf5’)
    filelist.append(‘../Create/example1.hdf5’)
    filelist.append(‘../Create/example2.hdf5’)
    filelist.append(‘../Create/example3.hdf5’)
    filelist.append(‘../Create/example4.hdf5’)
    filelist.append(‘../../../NeXus/definitions/trunk/manual/examples/h5py/prj_test.nexus.hdf5’)
    filelist.append(‘../../../NeXus/definitions/exampledata/code/hdf5/dmc01.h5’)

2.2. Code Examples that do not use the NAPI
filelist.append('../../../NeXus/definitions/exampledata/code/hdf5/dmc02.h5')
filelist.append('../../../NeXus/definitions/exampledata/code/hdf5/focus2007n001335.hdf')
filelist.append('../../../NeXus/definitions/exampledata/code/hdf5/NXtest.h5')
filelist.append('../../../NeXus/definitions/exampledata/code/hdf5/sans2009n012333.hdf')
filelist.append('../Create/simple5.nxs')
filelist.append('../Create/bad.h5')

do_filelist(filelist, limit)

def main():
    """standard command-line interface""
    try:
        opts, args = getopt.getopt(sys.argv[1:], "n:")
    except:
        print
        print "usage: ", sys.argv[0], " [-n ##] HDF5_file_name [another_HDF5_file_name]"
        print " -n ## : limit number of displayed array items to ## (must be 3 or more or 'None')"
        print
        for item in opts:
            if item[0] == "-n":
                if item[1].lower() == "none":
                    limit = None
                else:
                    limit = int(item[1])
                do_filelist(args)

    if __name__ == '__main__':
        if len(sys.argv) > 1:
            main()
        else:
            do_test()

2.2.3 Viewing 2-D Data from LRMECS

The IPNS LRMECS instrument stored data in NeXus HDF4 data files. One such example is available from the repository of NeXus data file examples. For this example, we will start with a conversion of that original data file into HDF5 format.

HDF4  http://svn.nexusformat.org/definitions/exampledata/IPNS/LRMECS/lrcs3701.nxs

HDF5  http://svn.nexusformat.org/definitions/exampledata/IPNS/LRMECS/lrcs3701.nx5

This dataset contains two histograms with 2-D images (148x750 and 148x32) of 32-bit integers. First, we use the h5dump tool to investigate the header content of the file (not showing any of the data).

Visualize Using h5dump

Here, the output of the command:

```
    h5dump -H lrcs3701.nx5
```

has been edited to only show the first NXdata group (/Histogram1/data):
LRMECS lrcs3701 data: h5dump output

1 HDF5 "C:\Users\Pete\Documents\eclipse\NeXus\definitions\exampledata\IPNS\LRMECS\lrcs3701.nx5" {
2  GROUP "/Histogram1/data" {
3    DATASET "data" {
4      DATATYPE H5T_STD_I32LE
5      DATASPACE SIMPLE { ( 148, 750 ) / ( 148, 750 ) }
6    }
7    DATASET "polar_angle" {
8      DATATYPE H5T_IEEE_F32LE
9      DATASPACE SIMPLE { ( 148 ) / ( 148 ) }
10   }
11   DATASET "time_of_flight" {
12      DATATYPE H5T_IEEE_F32LE
13      DATASPACE SIMPLE { ( 751 ) / ( 751 ) }
14   }
15   DATASET "title" {
16      DATATYPE H5T_STRING {
17        STRSIZE 44;
18        STRPAD H5T_STR_NULLTERM;
19        CSET H5T_CSET_ASCII;
20        CTYPEN H5T_STRING;
21      }
22      DATASPACE SIMPLE { ( 1 ) / ( 1 ) }
23   }
24 }
25 }

Visualize Using HDFview

For many, the simplest way to view the data content of an HDF5 file is to use the HDFview program (http://www.hdfgroup.org/hdf-java-html/hdfview) from The HDF Group. After starting HDFview, the data file may be loaded by dragging it into the main HDF window. On opening up to the first NXdata group /Histogram1/data (as above), and then double-clicking the dataset called: data, we get our first view of the data.

The data may be represented as an image by accessing the Open As menu from HDFview (on Windows, right click the dataset called data and select the Open As item, consult the HDFview documentation for different platform instructions). Be sure to select the Image radio button, and then (accepting everything else as a default) press the Ok button.

Note: In this image, dark represents low intensity while white represents high intensity.

LRMECS lrcs3701 data: image

Visualize Using IgorPro

Another way to visualize this data is to use a commercial package for scientific data visualization and analysis. One such package is IgorPro from http://www.wavemetrics.com

IgorPro provides a browser for HDF5 files that can open our NeXus HDF5 and display the image. Follow the instructions from WaveMetrics to install the HDF5 Browser package: http://www.wavemetrics.com/products/igorpro/dataaccess/hdf5.htm
Figure 2.7: LRMECS lrcs3701 data: HDFview

Figure 2.8: LRMECS lrcs3701 data: HDFview Open As dialog
You may not have to do this step if you have already installed the HDF5 Browser. IgorPro will tell you if it is not installed properly. To install the HDF5 Browser, first start IgorPro. Next, select from the menus and submenus: Data; Load Waves; Packages; Install HDF5 Package as shown in the next figure. IgorPro may direct you to perform more activities before you progress from this step.

Next, open the HDF5 Browser by selecting from the menus and submenus: Data; Load Waves; New HDF5 Browser as shown in the next figure.

Next, click the Open HDF5 File button and open the NeXus HDF5 file lrcs3701.nxs. In the lower left Groups panel, click the data dataset. Also, under the panel on the right called Load Dataset Options, choose No Table as shown. Finally, click the Load Dataset button (in the Datasets group) to display the image.

Note: In this image, dark represents low intensity while white represents high intensity. The image has been rotated for easier representation in this manual.
Figure 2.10: LRMECS lrcs3701 data: *IgorPro* install HDF5 Browser
Figure 2.11: LRMECS lrcs3701 data: *IgorPro HDFBrowser* dialog

2.2. Code Examples that do not use the NAPI
Figure 2.12: LRMECS lrcs3701 data: IgorPro HDFBrowser dialog

Figure 2.13: LRMECS lrcs3701 data: IgorPro Image
3.1 Introduction to NeXus definitions

While the design principles of NeXus are explained in the *NeXus: User Manual*, this Reference Documentation specifies all allowed *base classes* and all standardized *application definitions*. Furthermore, it also contains *contributed* definitions of new base classes or application definitions that are currently under review.

Base class definitions and application definitions have basically the same structure, but different semantics: Base class definitions define the *complete* set of terms that *might* be used in an instance of that class. Application definitions define the *minimum* set of terms that *must* be used in an instance of that class.

Base classes and application definitions are specified using a domain-specific XML scheme, the *NXDL: The NeXus Definition Language*.

3.1.1 Overview of NeXus definitions

For each class definition, the documentation is derived from content provided in the NXDL specification.

The documentation for each class consists of:

1. **short table:**
   - the current version of the NXDL specification used for the class
   - the category of the class (base class / application definition / contributed definition)
   - The NeXus class extended by this class. Most NeXus base classes only extend the base class definition (NXDL).
   - any other base classes (groups) cited by this class
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

2. **symbol list:** keywords used to designate array dimensions. At present, this list is not guaranteed to be complete (some array dimension names appear only in the description column of the class member table, and not here)

3. **source:** a link to the authoritative NXDL source

4. **tree outline:** hierarchical list of members.

5. **member table:** list of top-level members with natural-language annotations.

6. **supplementary member tables as needed:** member tables of subgroups.

### 3.1.2 Tree outlines

A compact listing of the basic structure (groups, fields, dimensions, attributes, and links) is prepared for each NXDL specification. Indentation shows nested structure. **Attributes** are prepended with the @ symbol. **Links** use the characters --&gt; to represent the path to the intended source of the information.

### 3.1.3 Member tables

Member tables provide basic information about each field or group in the class. An example of the varieties of specifications are given in the following table using items found in various NeXus base classes.

<table>
<thead>
<tr>
<th>Name</th>
<th>Type</th>
<th>Units</th>
<th>Description (and Occurrences)</th>
</tr>
</thead>
<tbody>
<tr>
<td>program_name</td>
<td>NX_CHAR</td>
<td></td>
<td>Name of program used to generate this file</td>
</tr>
<tr>
<td>@version</td>
<td>NX_CHAR</td>
<td></td>
<td>Program version number&lt;br&gt;Occurences: 1 : default</td>
</tr>
<tr>
<td>@configuration</td>
<td>NX_CHAR</td>
<td></td>
<td>configuration of the program</td>
</tr>
<tr>
<td>thumbnail</td>
<td>NXnote</td>
<td></td>
<td>A small image that is representative of the entry. An example of this is a 640x480 JPEG image automatically produced by a low resolution plot of the NXdata.</td>
</tr>
<tr>
<td>@mime_type</td>
<td>NX_CHAR</td>
<td></td>
<td>expected: mime_type=&quot;image/*&quot;</td>
</tr>
<tr>
<td>NXgeometry</td>
<td></td>
<td></td>
<td>describe the geometry of this class</td>
</tr>
<tr>
<td>distance</td>
<td>NX_FLOAT</td>
<td>NX_LENGTH</td>
<td>Distance from sample</td>
</tr>
<tr>
<td>mode</td>
<td>“Single Bunch”</td>
<td>“Multi Bunch”</td>
<td>source operating mode</td>
</tr>
<tr>
<td>target_material</td>
<td>Ta</td>
<td>W</td>
<td>depleted_U</td>
</tr>
</tbody>
</table>

The columns in the table are described as follows:

**Name (and attributes)** Name of the data field. Since name needs to be restricted to valid program variable names, no “-” characters can be allowed. Name must satisfy both HDF and XML naming.

1. NameStartChar ::= _ | a..z | A..Z

2. NameChar ::= NameStartChar | 0..9

3. Name ::= NameStartChar (NameChar)*

4. Or, as a regular expression: [_a-zA-Z][_a-zA-Z0-9]*

5. equivalent regular expression: [_a-zA-Z][_][k_]*
Attributes, identified with a leading “@” symbol (@) and belong with the preceding field or group, are additional metadata used to define this field or group. In the example above, the `program_name` element has two attributes: `version` (required) and `configuration` (optional) while the `thumbnail` element has one attribute: `mime_type` (optional).

For groups, the name may not be declared in the NXDL specification. In such instances, the `value shown in parentheses` in the `Name and Attributes` column is a suggestion, obtained from the group by removing the “NX” prefix. See `NXentry` for examples.

**Type** Type of data to be represented by this variable. The type is one of those specified in *NXDL: The NeXus Definition Language*. In the case where the variable can take only one value from a known list, the list of known values is presented, such as in the `target_material` field above: Ta | W | depleted_U | enriched_U | Hg | Pb | C. Selections with included whitespace are surrounded by quotes. See the example above for usage.

For fields, the data type may not be specified in the NXDL file. The default data type is `NX_CHAR` and this is shown in parentheses in the `Type` column. See `NXdata` for examples.

**Units** Data units, given as character strings, must conform to the NeXus units standard. See the `NeXus units` section for details.

**Description (and Occurrences)** A simple text description of the data field. No markup or formatting is allowed. The absence of `Occurrences` in the item description signifies that both `minOccurs` and `maxOccurs` have the default values. If the number of occurrences of an item are specified in the NXDL (through `@minOccurs` and `@maxOccurs` attributes), they will be reported in the `Description` column similar to the example shown above. Default values for occurrences are shown in the following table. The `NXDL element type` is either a group (such as a NeXus base class), a field (that specifies the name and type of a variable), or an attribute of a field or group. The number of times an item can appear ranges between `minOccurs` and `maxOccurs`. A default `minOccurs` of zero means the item is optional. For attributes, `maxOccurs` cannot be greater than 1.

<table>
<thead>
<tr>
<th>NXDL element type</th>
<th>minOccurs</th>
<th>maxOccurs</th>
</tr>
</thead>
<tbody>
<tr>
<td>group</td>
<td>0</td>
<td>unbounded</td>
</tr>
<tr>
<td>field</td>
<td>0</td>
<td>unbounded</td>
</tr>
<tr>
<td>attribute</td>
<td>0</td>
<td>1</td>
</tr>
</tbody>
</table>

### 3.2 NXDL: The NeXus Definition Language

Information in NeXus data files is arranged by a set of rules. These rules facilitate the exchange of data between scientists and software by standardizing common terms such as the way engineering units are described and the names for common things and the way that arrays are described and stored.

The set of rules for storing information in NeXus data files is declared using the NeXus Definition Language. NXDL itself is governed by a set of rules (a schema) that should simplify learning the few terms in NXDL. In fact, the NXDL rules, written as an XML Schema, are machine-readable using industry-standard and widely-available software tools for XML files such as `xsltproc` and `xmllint`. This chapter describes the rules and terms from which NXDL files are constructed.

#### 3.2.1 Introduction

NeXus Definition Language (NXDL) files allow scientists to define the nomenclature and arrangement of information in NeXus data files. These NXDL files can be specific to a scientific discipline such as tomography or small-angle scattering, specific analysis or data reduction software, or even to define another component (base class) used to design and build NeXus data files.
In addition to this chapter and the Tutorial chapter, look at the set of NeXus NXDL files to learn how to read and write NXDL files. These files are available from the NeXus definitions repository and are most easily viewed on GitHub: https://github.com/nexusformat/definitions in the base_classes, applications, and contributed directories. The rules (expressed as XML Schema) for NXDL files may also be viewed from this URL. See the files nxdl.xsd for the main XML Schema and nxdltypes.xsd for the listings of allowed data types and categories of units allowed in NXDL files.

NXDL files can be checked (validated) for syntax and content. With validation, scientists can be certain their definitions will be free of syntax errors. Since NXDL is based on the XML standard, there are many editing programs available to ensure that the files are well-formed. There are many standard tools such as xmlLint and xsltproc that can process XML files. Further, NXDL files are backed by a set of rules (an XML Schema) that define the language and can be used to check that an NXDL file is both correct by syntax and valid by the NeXus rules.

NXDL files are machine-readable. This enables their automated conversion into schema files that can be used, in combination with other NXDL files, to validate NeXus data files. In fact, all of the tables in the Class Definitions Chapter have been generated directly from the NXDL files.

The language of NXDL files is intentionally quite small, to provide only that which is necessary to describe scientific data structures (or to establish the necessary XML structures). Rather than have scientists prepare XML Schema files directly, NXDL was designed to reduce the jargon necessary to define the structure of data files. The two principle objects in NXDL files are: group and field. Documentation (doc) is optional for any NXDL component. Either of these objects may have additional attributes that contribute simple metadata.

The Class Definitions Chapter lists the various classes from which a NeXus file is constructed. These classes provide the glossary of items that could, in principle, be stored in a standard-conforming NeXus file (other items may be inserted into the file if the author wishes, but they won’t be part of the standard). If you are going to include a particular piece of metadata, refer to the class definitions for the standard nomenclature. However, to assist those writing data analysis software, it is useful to provide more than a glossary; it is important to define the required contents of NeXus files that contain data from particular classes of neutron, X-ray, or muon instrument.

**NXDL Elements and Data Types**

The documentation in this section has been obtained directly from the NXDL Schema file: nxdl.xsd. First, the basic elements are defined in alphabetical order. Attributes to an element are indicated immediately following the element and are preceded with an “@” symbol, such as @attribute. Then, the common data types used within the NXDL specification are defined. Pay particular attention to the rules for validItemName and validNXClassName.

**NXDL Elements**

*attribute* An attribute element can only be a child of a field or group element. It is used to define attribute elements to be used and their data types and possibly an enumeration of allowed values.

For more details, see: attributeType

*definition* A definition element can only be used at the root level of an NXDL specification. Note: Due to the large number of attributes of the definition element, they have been omitted from the figure below.

For more details, see: definition, definitionType, and definitionTypeAttr

---

1 For example XML Copy Editor (http://xml-copy-editor.sourceforge.net/)
2 http://en.wikipedia.org/wiki/XML#Well-formedness_and_error-handling
Figure 3.1: Graphical representation of the NXDL attribute element

**dimensions**  The dimensions element describes the shape of an array. It is used only as a child of a field element. For more details, see: dimensionsType

**doc**  A doc element can be a child of most NXDL elements. In most cases, the content of the doc element will also become part of the NeXus manual. For more details, see: docType

**enumeration**  An enumeration element can only be a child of a field or attribute element. It is used to restrict the available choices to a predefined list, such as to control varieties in spelling of a controversial word (such as metre vs. meter). For more details, see: enumerationType

**field**  The field element provides the value of a named item. Many different attributes are available to further define the field. Some of the attributes are not allowed to be used together (such as axes and axis); see the documentation of each for details. It is used only as a child of a group element. For more details, see: fieldType

3.2. NXDL: The NeXus Definition Language
Figure 3.2: Graphical representation of the NXDL definition element
3.2. NXDL: The NeXus Definition Language

Figure 3.3: Graphical representation of the NXDL dimensions element

Figure 3.4: Graphical representation of the NXDL doc element
Figure 3.5: Graphical representation of the NXDL `enumeration` element

Figure 3.6: Graphical representation of the NXDL `field` element
group  A group element can only be a child of a definition or group element. It describes a common level of organization in a NeXus data file, similar to a subdirectory in a file directory tree.
For more details, see: groupType

link  A link element can only be a child of a field or group element. It describes the path to the original source of the parent field or group.
For more details, see: linkType

symbols  A symbols element can only be a child of a definition element. It defines the array index symbols to be used when defining arrays as field elements with common dimensions and lengths.
For more details, see: symbolsType

NXDL Data Types (internal)

Data types that define the NXDL language are described here. These data types are defined in the XSD Schema (nxdl.xsd) and are used in various parts of the Schema to define common structures or to simplify a complicated entry. While the data types are not intended for use in NXDL specifications, they define structures that may be used in NXDL specifications.

attributeType  Any new group or field may expect or require some common attributes.
(This data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

Attributes of attributeType

@name
   Name of the attribute (unique within the enclosing group).

@type
   Type of the attribute. For group specifications, the class name. For field or attribute specifications, the NXDL data type.

Elements of attributeType

doc
   Description of this attribute. This documentation will go into the manual.

enumeration
   An enumeration specifies the values to be used.

definition  A definition element is the group at the root of every NXDL specification. It may only appear at the root of an NXDL file and must only appear once for the NXDL to be well-formed.
Figure 3.7: Graphical representation of the NXDL group element
Figure 3.8: Graphical representation of the NXDL `link` element

Figure 3.9: Graphical representation of the NXDL `symbols` element
**definitionType**  
A definition is the root element of every NXDL definition. It may only appear at the root of an NXDL file and must only appear once for the NXDL to be well-formed.

The `definitionType` defines the documentation, attributes, fields, and groups that will be used as children of the `definition` element. Could contain these elements:

- attribute
- doc
- field
- group
- link

Note that a definition element also includes the definitions of the `basicComponent` data type. (The `definitionType` data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

**Attributes of definitionType**

@category

NXDL base definitions define the dictionary of terms to use for these components. All terms in a base definition are optional. NXDL application definitions define what is required for a scientific interest. All terms in an application definition are required. NXDL contributed definitions may be considered either base or applications. Contributed definitions `<emphasis>must</emphasis>` indicate their intended use, either as a base class or as an application definition.

@extends

The `extends` attribute allows this definition to subclass from another NXDL, otherwise `extends="NXobject"` should be used.

@ignoreExtraAttributes

Only validate known attributes; do not not warn about unknowns. The `ignoreExtraAttributes` attribute is a flag to the process of validating NeXus data files. By setting `ignoreExtraAttributes="true"`, presence of any undefined attributes in this class will not generate warnings during validation. Normally, validation will check all the attributes against their definition in the NeXus base classes and application definitions. Any items found that do not match the definition in the NXDL will generate a warning message.

The `ignoreExtraAttributes` attribute should be used sparingly!

@ignoreExtraFields

Only validate known fields; do not not warn about unknowns. The `ignoreExtraFields` attribute is a flag to the process of validating NeXus data files. By setting `ignoreExtraFields="true"`, presence of any undefined fields in this class will not generate warnings during validation. Normally, validation will check all the fields against their definition in the NeXus base classes and application definitions. Any items found that do not match the definition in the NXDL will generate a warning message.

The `ignoreExtraFields` attribute should be used sparingly!
@ignoreExtraGroups

Only validate known groups; do not not warn about unknowns. The ignoreExtraGroups attribute is a flag to the process of validating NeXus data files. By setting ignoreExtraGroups="true", presence of any undefined groups in this class will not generate warnings during validation. Normally, validation will check all the groups against their definition in the NeXus base classes and application definitions. Any items found that do not match the definition in the NXDL will generate a warning message.

The ignoreExtraGroups attribute should be used sparingly!

@name

The name of this NXDL file (without the file extensions). The name must be unique amongst all the NeXus base class, application, and contributed definitions. For the class to be adopted by the NIAC, the first two letters must be “NX” (in uppercase). Any other use must not begin with “NX” in any combination of upper or lower case.

@restricts

The restricts attribute is a flag to the data validation. When restricts="1", any non-standard component found (and checked for validity against this NXDL specification) in a NeXus data file will be flagged as an error. If the restricts attribute is not present, any such situations will produce a warning.

@svnid

(2014-08-19: deprecated since switch to GitHub version control) The identifier string from the subversion revision control system. This reports the time stamp and the revision number of this file. (Updated automatically, unlike the version attribute.)

@type

Must be type="group"

@version

Version of this NXDL definition. Each NXDL specification may have a different version to facilitate software maintenance. This value is modified by the person who edits this file when this NXDL specification has changed significantly (in a way that downstream software should be aware).

Elements of definitionType

symbols

Use a symbols list to define each of the mnemonics that represent the length of each dimension in a vector or array.

Groups under definitionType

In addition to an optional symbols list, a definition may contain any of the items allowed in a group.
**definitionTypeAttr** Prescribes the allowed values for `definition type` attribute. (This data type is used internally in the NXDL schema to define a data type.)

The value may be any one from this list only:

- `group`
- `definition`

**dimensionsType** dimensions of a data element in a NeXus file (This data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

**Attributes of dimensionsType**

@rank


**Elements of dimensionsType**

dim

Specify the parameters for each index of the `dimensions` element with a `dim` element. The number of `dim` entries should be equal to the `rank` of the array. For example, these terms describe a 2-D array with lengths `(nsurf, nwl)`:

```
<dimensions rank="2">
  <dim index="1" value="nsurf"/>
  <dim index="2" value="nwl"/>
</dimensions>
```

The `value` attribute is used by NXDL and also by the NeXus data file validation tools to associate and coordinate the same array length across multiple fields in a group.

@incr

The dimension specification is related to the `refindex` axis within the `ref` field by an offset of `incr`. Requires `ref` and `refindex` attributes to be present.

@index

Number or symbol indicating which axis (subscript) is being described, ranging from 1 up to `rank` (rank of the data structure). For example, given an array `A[i, j, k].index="1"` would refer to the `i` axis (subscript). (`NXdata uses index="0" to indicate a situation when the specific index is not known a priori.`)

@ref

The dimension specification is the same as that in the `ref` field, specified either by a relative path, such as `polar_angle` or `../Qvec` or absolute path, such as `/entry/path/to/follow/to/ref/field.`
@refindex

The dimension specification is the same as the refindex axis within the ref field. Requires ref attribute to be present.

@value

Integer length (number of values), or mnemonic symbol representing the length of this axis.

doc

Documentation might be necessary to describe how the parts of the dimensions element are to be used.

docType

NXDL allows for documentation on most elements using the doc element. The documentation is useful in several contexts. The documentation will be rendered in the manual. Documentation, is provided as tooltips by some XML editors when editing NXDL files. Simple documentation can be typed directly in the NXDL:

```xml
<field name="name">
  <doc>Descriptive name of sample</doc>
</field>
```

This is suitable for basic descriptions that do not need extra formatting such as a bullet-list or a table. For more advanced control, use the rules of restructured text, such as in the NXdetector specification. Refer to examples in the NeXus base classNXDL files such as NXdata.

Could contain these elements:

- any

(This data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

enumerationType

An enumeration restricts the values allowed for a specification. Each value is specified using an item element, such as: `<item value="Synchrotron X-ray Source"/>`. Could contain these elements:

- doc
- item

(This data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

```xml
<field name="mode">
  <doc>source operating mode</doc>
  <enumeration>
    <item value="Single Bunch"><doc>for storage rings</doc></item>
    <item value="Multi Bunch"><doc>for storage rings</doc></item>
    <!-- other sources could add to this -->
  </enumeration>
</field>
```

Elements of enumerationType

item

One of the prescribed values. Use the value attribute.
@value

The value of value of an enumItem is defined as an attribute rather than a name.

doc

Individual items can be documented but this documentation might not be printed in the NeXus Reference Guide.

fieldType

A field declares a new element in the component being defined. A field is synonymous with the HDF4 SDS (Scientific Data Set) and the HDF5 dataset terms. Could contain these elements:

- attribute
- dimensions
- doc
- enumeration

Note that a field element also includes the definitions of the basicComponent data type. (The fieldType data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

@axes

Presence of the axes attribute means this field is an ordinate.

This attribute contains a colon (or comma in legacy files) delimited list of the names of independent axes when plotting this field. Each name in this list must exist as a field in the same group. <!– perhaps even discourage use of square brackets in axes attribute? –> (Optionally, the list can be enclosed by square brackets but this is not common.) The regular expression for this rule is:

[A-Za-z_][\w_]*([ :][A-Za-z_][\w_]*)*

@axis

NOTE: Use of this attribute is discouraged. It is for legacy support. You should use the axes attribute instead.

Presence of the axis attribute means this field is an abcissa.

The attribute value is an integer indicating this field as an axis that is part of the data set. The data set is a field with the attribute signal=1 in the same group. The value can range from 1 up to the number of independent axes (abcissae) in the data set.

A value of axis=1 indicates that this field contains the data for the first independent axis. For example, the X axis in an XY data set.

A value of axis=2 indicates that this field contains the data for the second independent axis. For example, the Y axis in a 2-D data set.

A value of axis=3 indicates that this field contains the data for the third independent axis. For example, the Z axis in a 3-D data set.

A field with an axis attribute should not have a signal attribute.
@interpretation

This instructs the consumer of the data what the last dimensions of the data are. It allows plotting software to work out the natural way of displaying the data.

For example a single-element, energy-resolving, fluorescence detector with 512 bins should have interpretation="spectrum". If the detector is scanned over a 512 x 512 spatial grid, the data reported will be of dimensions: 512 x 512 x 512. In this example, the initial plotting representation should default to data of the same dimensions of a 512 x 512 pixel image detector where the images where taken at 512 different pressure values.

In simple terms, the allowed values mean:

- scaler = 0-D data to be plotted
- spectrum = 1-D data to be plotted
- image = 2-D data to be plotted
- vertex = 3-D data to be plotted

@long_name

Descriptive name for this field (may include whitespace and engineering units). Often, the long_name (when defined) will be used as the axis label on a plot.

@maxOccurs

Defines the maximum number of times this element may be used. Its value is confined to zero or greater. Must be greater than or equal to the value for the “minOccurs” attribute. A value of “unbounded” is allowed.

@minOccurs

 Defines the minimum number of times this element may be used. Its value is confined to zero or greater. Must be less than or equal to the value for the “maxOccurs” attribute. A value of “unbounded” is allowed.

@offset

The stride and offset attributes are used together to index the array of data items in a multi-dimensional array. They may be used as an alternative method to address a data array that is not stored in the standard NeXus method of “C” order.

The offset attribute determines the starting coordinates of the data array for each dimension.


The offset attribute contains a comma-separated list of integers. (In addition to the required comma delimiter, whitespace is also allowed to improve readability.) The number of items in the list is equal to the rank of the data being stored. The value of each item is the offset in the array of the first data item of that subscript of the array.

@primary

Integer indicating the priority of selection of this field for plotting (or visualization) as an axis.

Presence of the primary attribute means this field is an abcissa.
@signal

Presence of the signal attribute means this field is an ordinate.

Integer marking this field as plottable data (ordinates). The value indicates the priority of selection or interest. Some facilities only use signal=1 while others use signal=2 to indicate plottable data of secondary interest. Higher numbers are possible but not common and interpretation is not standard.

A field with a signal attribute should not have an axis attribute.

@stride

The stride and offset attributes are used together to index the array of data items in a multi-dimensional array. They may be used as an alternative method to address a data array that is not stored in the standard NeXus method of “C” order.

The stride list chooses array locations from the data array with each value in the stride list determining how many elements to move in each dimension. Setting a value in the stride array to 1 moves to each element in that dimension of the data array, while setting a value of 2 in a location in the stride array moves to every other element in that dimension of the data array. A value in the stride list may be positive to move forward or negative to step backward. A value of zero will not step (and is of no particular use).


The stride attribute contains a comma-separated list of integers. (In addition to the required comma delimiter, whitespace is also allowed to improve readability.) The number of items in the list is equal to the rank of the data being stored. The value of each item is the spacing of the data items in that subscript of the array.

@type

Defines the type of the element as allowed by the NAPI (NeXus Application Programmer Interface). See elsewhere for the complete list of allowed NAPI types.

@units

String describing the engineering units. The string should be appropriate for the value and should conform to the NeXus rules for units. Conformance is not validated at this time.

attribute

attributes to be used with this field

dimensions

dimensions of a data element in a NeXus file

enumeration

A field can specify which values are to be used
**groupType**  A group element refers to the definition of an existing NX object or a locally-defined component. Could contain these elements:

- attribute
- doc
- field
- group
- link

Note that a group element also includes the definitions of the basicComponent data type. (The groupType data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)

**Attributes of groupType**

@maxOccurs

Maximum number of times this group is allowed to be present within its parent group. Note each group must have a name attribute that is unique among all group and field declarations within a common parent group.

@minOccurs

Minimum number of times this group is allowed to be present within its parent group. Note each group must have a name attribute that is unique among all group and field declarations within a common parent group.

@name

A particular scientific application may expect a name of a group element. It is helpful but not required to specify the name attribute in the NXDL file. It is suggested to always specify a name to avoid ambiguity. It is also suggested to derive the name from the type, using an additional number suffix as necessary. For example, consider a data file with only one NXentry. The suggested default name would be entry. For a data file with two or more NXentry groups, the suggested names would be entry1, entry2, ... Alternatively, a scientific application such as small-angle scattering might require a different naming procedure; two different NXaperture groups might be given the names beam-defining slit and scatter slit.

@type

The type attribute must contain the name of a NeXus base class, application definition, or contributed definition.

linkType  A link to another item. Use a link to avoid needless repetition of information. (This data type is used internally in the NXDL schema to define elements and attributes to be used by users in NXDL specifications.)
@target

A link to another item, represented as an absolute path to an existing field such as
/NXentry/NXinstrument/NXdetector/polar_angle. Could contain these elements:

- doc

Matching regular expression:

(/[a-zA-Z_][\w_]*(:[a-zA-Z_][\w_]*)?)+

symbolsType  Each symbol has a name and optional documentation. Please provide documentation that indicates
what each symbol represents. For example:

<symbols>
  <symbol name="nsurf"><doc>number of reflecting surfaces</doc></symbol>
  <symbol name="nwl"><doc>number of wavelengths</doc></symbol>
</symbols>

Elements of symbolsType

doc

Describe the purpose of this list of symbols. This documentation will go into the manual.

symbol

When multiple field elements share the same dimensions, such as the dimension scales associated with
plottable data in an NXdata group, the length of each dimension written in a NeXus data file should be
something that can be tested by the data file validation process.

@name

Mnemonic variable name for this array index symbol.

doc

Describe the purpose of the parent symbol. This documentation will go into the manual.

basicComponent  A basicComponent defines the allowed name format and attributes common to all field and
group specifications. (This data type is used internally in the NXDL schema to define elements and attributes to be
used by users in NXDL specifications.)

Attributes of basicComponent

@deprecated

The presence of the deprecated attribute indicates to the data file validation process that an advisory
message (specified as the content of the deprecated attribute) will be reported. Future versions of the
NXDL file might not define (or even re-use) this component. For example:

deprecated="as of release MAJOR.MINOR"

Note: because deprecated is an attribute, the XML rules do not permit it to have any element content.
@name

The `name` attribute is the identifier string for this entity. It is required that `name` must be unique within the enclosing `group`. The rule (`validItemName`) is defined to only allow names that can be represented as valid variable names in most computer languages.

Elements of basicComponent

doc

Describe this `basicComponent` and its use. This documentation will go into the manual.

`validItemName`  Used for allowed names of elements and attributes. Need to be restricted to valid program variable names. Note: This means no “-” or “.” characters can be allowed and you cannot start with a number. HDF4 had a 64 character limit on names (possibly including NULL) and NeXus enforces this via the `NX_MAXNAMELEN` variable with a 64 character limit (which may be 63 on a practical basis if one considers a NULL terminating byte). (This data type is used internally in the NXDL schema to define a data type.)

The value may be any `xs:token` that also matches the regular expression:

```
[A-Za-z_][\w_]*
```

`validNXClassName`  Used for allowed names of NX class types (e.g. NXdetector) not the instance (e.g. bank1) which is covered by `validItemName`. (This data type is used internally in the NXDL schema to define a data type.)

The value may be any `nx:validItemName` that also matches the regular expression:

```
NX.+
```

`validTargetName`  This is a valid link target - currently it must be an absolute path made up of valid names with the `/` character delimiter. But we may want to consider allowing “..” (parent of directory) at some point. If the `name` attribute is helpful, then use it in the path with the syntax of `name:type` as in these examples:

```
/NXentry/NXinstrument/analyzer:NXcrystal/ef
/NXentry/NXinstrument/monochromator:NXcrystal/ei
/NX_other
```

Must also consider use of `name` attribute in resolving link targets. (This data type is used internally in the NXDL schema to define a data type.)

From the HDF5 documentation (http://www.hdfgroup.org/HDF5/doc/UG/UG_frame09Groups.html):

Note that relative path names in HDF5 do not employ the ‘‘../’’ notation, the UNIX notation indicating a parent directory, to indicate a parent group.

Thus, if we only consider the case of `name:type`, the matching regular expression syntax is written: `/([a-zA-Z_][\w_]*(:[a-zA-Z_][\w_]*))*`. Note that HDF5 also permits relative path names, such as: GroupA/GroupB/Dataset1 but this is not permitted in the matching regular expression and not supported in NAPI.

The value may be any `xs:token` that also matches the regular expression:

```
(/[a-zA-Z_][\w_]*(:[a-zA-Z_][\w_]*))*
```
nonNegativeUnbounded  A nonNegativeUnbounded allows values including all positive integers, zero, and the string unbounded. (This data type is used internally in the NXDL schema to define a data type.)

The xs:string data type The xs:string data type can contain characters, line feeds, carriage returns, and tab characters. See http://www.w3schools.com/Schema/schema_dtypes_string.asp for more details.

The xs:token data type The xs:string data type is derived from the xs:string data type. The xs:token data type also contains characters, but the XML processor will remove line feeds, carriage returns, tabs, leading and trailing spaces, and multiple spaces. See http://www.w3schools.com/Schema/schema_dtypes_string.asp for more details.

NXDL: Data Types and Units

Data Types allowed in NXDL specifications

Data types for use in NXDL describe the expected type of data for a NeXus field. These terms are very broad. More specific terms are used in actual NeXus data files that describe size and array dimensions. In addition to the types in the following table, the NAPI type is defined when one wishes to permit a field with any of these data types.

<table>
<thead>
<tr>
<th>Data Type</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>ISO8601</td>
<td>ISO 8601 date and time representation (<a href="http://www.w3.org/TR/NOTE-datetime">http://www.w3.org/TR/NOTE-datetime</a>)</td>
</tr>
<tr>
<td>NX_BINARY</td>
<td>any representation of binary data - if text, line terminator is [CR][LF]</td>
</tr>
<tr>
<td>NX_BOOLEAN</td>
<td>true/false value ( true</td>
</tr>
<tr>
<td>NX_CHAR</td>
<td>any string representation</td>
</tr>
<tr>
<td>NX_DATE_TIME</td>
<td>alias of ISO8601</td>
</tr>
<tr>
<td>NX_FLOAT</td>
<td>any representation of a floating point number</td>
</tr>
<tr>
<td>NX_INT</td>
<td>any representation of an integer number</td>
</tr>
<tr>
<td>NX_NUMBER</td>
<td>any valid NeXus number representation</td>
</tr>
<tr>
<td>NX_POSINT</td>
<td>any representation of a positive integer number (greater than zero)</td>
</tr>
<tr>
<td>NX_UINT</td>
<td>any representation of an unsigned integer number (includes zero)</td>
</tr>
</tbody>
</table>

Unit Categories allowed in NXDL specifications

Unit categories in NXDL specifications describe the expected type of units for a NeXus field. They should describe valid units consistent with the NeXus units section. The values for unit categories are restricted (by an enumeration) to the following table.

<table>
<thead>
<tr>
<th>Unit Category</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>NX_ANGLE</td>
<td>example: degrees or radians or arcminutes or</td>
</tr>
<tr>
<td>NX_ANY</td>
<td>usage: things like logs that aren’t picky on units</td>
</tr>
<tr>
<td>NX_AREA</td>
<td>example: m2 or barns</td>
</tr>
<tr>
<td>NX_CHARGE</td>
<td>example: pC or C</td>
</tr>
<tr>
<td>NX_CROSS_SECTION</td>
<td>example: barns</td>
</tr>
<tr>
<td>NX_CURRENT</td>
<td>example: A</td>
</tr>
<tr>
<td>NX_DIMENSIONLESS</td>
<td>for fields where the units cancel out, example: “” or mm/mm (NOTE: not the same as NX_UNITLESS)</td>
</tr>
<tr>
<td>NX_EMITTANCE</td>
<td>emittance (length * angle) of a radiation source, example: nm*rad</td>
</tr>
</tbody>
</table>
NX_ENERGY  example: J or keV
NX_FLUX example: s-1 cm-2
NX_FREQUENCY example: Hz
NX_LENGTH example: m
NX_MASS example: g
NX_MASS_DENSITY example: g cm-3
NX_MOLECULAR_WEIGHT example: g mol-1
NX_PERIOD (alias to NX_TIME) period of pulsed source, example: microseconds
NX_PER_AREA example: cm-2
NX_PER_LENGTH example: cm-1
NX_POWER example: W
NX_PRESSURE example: Pa
NX_PULSES (alias to NX_NUMBER) clock pulses
NX_SCATTERING_LENGTH_DENSITY example: cm-2
NX_SOLID_ANGLE example: sr | steradian
NX_TEMPERATURE example: K
NX_TIME example: s
NX_TIME_OF_FLIGHT (alias to NX_TIME) example: s
NX_UNITLESS for fields that don’t have a unit (e.g. hkl) so that they don’t inherit the wrong units
  (NOTE: not the same as NX_DIMENSIONLESS)
NX_VOLTAGE example: V
NX_VOLUME example: m3
NX_WAVELENGTH example: Angstrom
NX_WAVENUMBER units for Q, example: Angstrom-1 or nm-1

3.3 Base Class Definitions

A description of each NeXus base class definition is given. NeXus base class definitions define the complete set of terms that might be used in an instance of that class. Consider the base classes as a set of components that are used to construct a data file.

3.3.1 NXaperture

Status:
  base class, extends NXobject, version 1.0

Description:
  Template of a beamline aperture.

Symbols:
Structure:

- **material**: `NX_CHAR`
  Absorbing material of the aperture

- **description**: `NX_CHAR`
  Description of aperture

- **(geometry)**: `NXgeometry`
  Location and shape of aperture

- **(geometry)**: `NXgeometry`
  Location and shape of each blade

- **(note)**: `NXnote`
  Describe an additional information in a note*

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXaperture.nxdl.xml

### 3.3.2 NXattenuator

**Status:**

- Base class, extends `NXobject`, version 1.0

**Description:**

Description of a device that reduces the intensity of a beam by attenuation. If uncertain whether to use `NXfilter` (band-pass filter) or `NXattenuator` (reduces beam intensity), then choose `NXattenuator`.

**Symbols:**

- No symbol table

**Groups cited:** none

**Structure:**

- **distance**: `NX_FLOAT` `{units=NX_LENGTH}`
  Distance from sample

- **type**: `NX_CHAR`
  Type or composition of attenuator, e.g. polythene

- **thickness**: `NX_FLOAT` `{units=NX_LENGTH}`
  Thickness of attenuator along beam direction

- **scattering_cross_section**: `NX_FLOAT` `{units=NX_CROSS_SECTION}`
  Scattering cross section (coherent+incoherent)

- **absorption_cross_section**: `NX_FLOAT` `{units=NX_CROSSSECTION}`
  Absorption cross section

- **attenuator_transmission**: `NX_FLOAT` `{units=NX_DIMENSIONLESS}`
The nominal amount of the beam that gets through (transmitted intensity)/(incident intensity)

**status**: `NX_CHAR`

In or out or moving of the beam

Any of these values: `in`, `out`, `moving`

**@time**: `NX_DATE_TIME`

time stamp for this observation

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXattenuator.nxdl.xml

### 3.3.3 NXbeam

**Status**: base class, extends `NXobject`, version 1.0

**Description**: Template of the state of the neutron or X-ray beam at any location. It will be referenced by beamline component groups within the NXinstrument group or by the NXsample group. Note that variables such as the incident energy could be scalar values or arrays. This group is especially valuable in storing the results of instrument simulations in which it is useful to specify the beam profile, time distribution etc. at each beamline component. Otherwise, its most likely use is in the NXsample group in which it defines the results of the neutron scattering by the sample, e.g., energy transfer, polarizations.

**Symbols**: No symbol table

**Groups cited**: `NXdata`

**Structure**:

- **distance**: `NX_FLOAT {units=NX_LENGTH}`
  Distance from sample

- **incident_energy[i]**: `NX_FLOAT {units=NX_ENERGY}`
  Energy on entering beamline component

- **final_energy[i]**: `NX_FLOAT {units=NX_ENERGY}`
  Energy on leaving beamline component

- **energy_transfer[i]**: `NX_FLOAT {units=NX_ENERGY}`
  Energy change caused by beamline component

- **incident_wavelength[i]**: `NX_FLOAT {units=NX_WAVELENGTH}`
  Wavelength on entering beamline component

- **incident_wavelength_spread[i]**: `NX_FLOAT {units=NX_WAVELENGTH}`
  Wavelength spread FWHM on entering component

- **incident_beam_divergence[2, j]**: `NX_FLOAT {units=NX_ANGLE}`
  Divergence of beam entering this component

- **final_wavelength[i]**: `NX_FLOAT {units=NX_WAVELENGTH}`
  Wavelength on leaving beamline component
incident_polarization[2, j]: NX_FLOAT {units=NX_ANY}
  Polarization vector on entering beamline component
final_polarization[2, j]: NX_FLOAT {units=NX_ANY}
  Polarization vector on leaving beamline component
final_wavelength_spread[i]: NX_FLOAT {units=NX_WAVELENGTH}
  Wavelength spread FWHM of beam leaving this component
final_beam_divergence[2, j]: NX_FLOAT {units=NX_ANGLE}
  Divergence FWHM of beam leaving this component
flux[i]: NX_FLOAT {units=NX_FLUX}
  flux incident on beam plane area
(data): NXdata
  Distribution of beam with respect to relevant variable e.g. wavelength. This is mainly useful for simulations which need to store plottable information at each beamline component.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXbeam.nxdl.xml

3.3.4 NXbeam_stop

Status:
  base class, extends NXobject, version 1.0

Description:
  A class for a beamstop. Beamstops and their positions are important for SANS and SAXS experiments.

Symbols:
  No symbol table

Groups cited: NXgeometry

Structure:
  description: NX_CHAR
    description of beamstop
    Any of these values: circular | rectangular
  size: NX_FLOAT {units=NX_LENGTH}
    size of beamstop
  x: NX_FLOAT {units=NX_LENGTH}
    x position of the beamstop in relation to the detector
  y: NX_FLOAT {units=NX_LENGTH}
    y position of the beamstop in relation to the detector
  distance_to_detector: NX_FLOAT {units=NX_LENGTH}
    distance of the beamstop to the detector
  status: NX_CHAR
Any of these values: in | out

(geometry): NXgeometry

engineering shape, orientation and position of the beam stop.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXbeam_stop.nxdl.xml

3.3.5 NXbending_magnet

Status:

base class, extends NXobject, version 1.0

Description:

description for a bending magnet

Symbols:

No symbol table

Groups cited: NXdata, NXgeometry

Structure:

critical_energy: NX_FLOAT {units=NX_ENERGY}
bending_radius: NX_FLOAT {units=NX_LENGTH}
magnetic_field: NX_FLOAT {units=NX_CURRENT}

strength of magnetic field of dipole magnets

accepted_photon_beam_divergence: NX_FLOAT {units=NX_LENGTH}

An array of four numbers giving X+, X-, Y+ and Y- half divergence

source_distance_x: NX_FLOAT {units=NX_LENGTH}

Distance of source point from particle beam waist in X (horizontal) direction.

source_distance_y: NX_FLOAT {units=NX_LENGTH}

Distance of source point from particle beam waist in Y (vertical) direction.

divergence_x_plus: NX_FLOAT {units=NX_ANGLE}

Accepted photon beam divergence in X+ (horizontal outboard) direction. Note that divergence_x_plus+divergence_x_minus is the total horizontal beam divergence.

divergence_x_minus: NX_FLOAT {units=NX_ANGLE}

Accepted photon beam divergence in X- (horizontal inboard) direction. Note that divergence_x_plus+divergence_x_minus is the total horizontal beam divergence.

divergence_y_plus: NX_FLOAT {units=NX_ANGLE}

Accepted photon beam divergence in Y+ (vertical upward) direction. Note that divergence_y_plus+divergence_y_minus is the total vertical beam divergence.

divergence_y_minus: NX_FLOAT {units=NX_ANGLE}

Accepted photon beam divergence in Y- (vertical downward) direction. Note that divergence_y_plus+divergence_y_minus is the total vertical beam divergence.

spectrum: NXdata

3.3. Base Class Definitions
bending magnet spectrum

(\textit{geometry}): \textit{NXgeometry}

“Engineering” position of bending magnet

\textbf{Source:} Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXbending_magnet.nxdl.xml

\section*{3.3.6 \textit{NXcapillary}}

\textbf{Status:}

base class, extends \textit{NXobject}, version 1.0

\textbf{Description:}

This is a dictionary of field names to use for describing a capillary as used in X-ray beamlines. Based on information provided by Gerd Wellenreuther.

\textbf{Symbols:}

No symbol table

\textbf{Groups cited:} \textit{NXdata}

\textbf{Structure:}

\textit{type}: \textit{NX_CHAR}

Type of the capillary

Any of these values:

\begin{itemize}
\item single\textunderscore bounce
\item polycapillary
\item conical\textunderscore capillary
\end{itemize}

\textit{manufacturer}: \textit{NX_CHAR}

The manufacturer of the capillary. This is actually important as it may have an impact on performance.

\textit{maximum\textunderscore incident\textunderscore angle}: \textit{NX_FLOAT} \{units=\textit{NX\textunderscore ANGLE}\}

\textit{accepting\textunderscore aperture}: \textit{NX_FLOAT} \{units=\textit{NX\textunderscore ANGLE}\}

\textit{working\textunderscore distance}: \textit{NX_FLOAT} \{units=\textit{NX\textunderscore LENGTH}\}

\textit{focal\textunderscore size}: \textit{NX_FLOAT}

The focal size in FWHM

\textit{gain}: \textit{NXdata}

The gain of the capillary as a function of energy

\textit{transmission}: \textit{NXdata}

The transmission of the capillary as a function of energy

\textbf{Source:} Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXcapillary.nxdl.xml
3.3.7 NXcharacterization

Status:
  base class, extends NXobject, version 1.0

Description:
  Note: This base class may be removed in future releases of NXDL. If you have a use for this base class, please provide a description of your intended use to the NIAC (nexus-committee@nexusformat.org).

Symbols:
  No symbol table

Groups cited: none

Structure:
  @source: NX_CHAR
  If missing, the source file is the current file
  @location: NX_CHAR
  @mime_type: NX_CHAR
  If missing, the source file is NAPI readable
definition: NX_CHAR
  @version: NX_CHAR
  @URL: NX_CHAR

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXcharacterization.nxdl.xml

3.3.8 NXcollection

Status:
  base class, extends NXobject, version 1.0

Description:
  Use NXcollection to gather together any set of terms. The original suggestion is to use this as a container class for the description of a beamline.

  For NeXus validation, NXcollection will always generate a warning since it is always an optional group. Anything (groups, fields, or attributes) placed in an NXcollection group will not be validated.

Symbols:
  No symbol table

Groups cited: none

Structure:
  beamline: NX_CHAR
  name of the beamline for this collection

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXcollection.nxdl.xml
3.3.9 NXcollimator

Status:
base class, extends NXobject, version 1.0

Description:
Template of a beamline collimator.

Symbols:
No symbol table

Groups cited: NXgeometry, NXlog

Structure:

- **type**: NX_CHAR
  - Any of these values: Soller | radial | oscillating | honeycomb

- **soller_angle**: NX_FLOAT {units=NX_ANGLE}
  - Angular divergence of Soller collimator

- **divergence_x**: NX_FLOAT {units=NX_ANGLE}
  - Divergence of collimator in local x direction

- **divergence_y**: NX_FLOAT {units=NX_ANGLE}
  - Divergence of collimator in local y direction

- **frequency**: NX_FLOAT {units=NX_FREQUENCY}
  - Frequency of oscillating collimator

- **blade_thickness**: NX_FLOAT {units=NX_LENGTH}
  - Blade thickness

- **blade_spacing**: NX_FLOAT {units=NX_LENGTH}
  - Blade spacing

- **absorbing_material**: NX_CHAR
  - Name of absorbing material

- **transmitting_material**: NX_CHAR
  - Name of transmitting material

- **(geometry)**: NXgeometry
  - Position, shape and size

- **frequency_log**: NXlog
  - Log of frequency

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXcollimator.nxdl.xml
3.3.10 NXcrystal

Status:

base class, extends NXobject, version 1.0

Description:

Template of a crystal monochromator or analyzer. Permits double bent monochromator comprised of multiple segments with anisotropic Gaussian mosaic.

If curvatures are set to zero or are absent, array is considered to be flat.

Scattering vector is perpendicular to surface. Crystal is oriented parallel to beam incident on crystal before rotation, and lies in vertical plane.

Symbols:

These symbols will be used below to coordinate dimensions with the same lengths.

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>n_comp</td>
<td>number of different unit cells to be described</td>
</tr>
<tr>
<td>i</td>
<td>number of wavelengths</td>
</tr>
</tbody>
</table>

Groups cited: NXdata, NXgeometry, NXlog, NXshape

Structure:

usage: NX_CHAR

How this crystal is used. Choices are in the list.

Any of these values:

- Bragg: reflection geometry
- Laue: The chemical formula specified using CIF conventions. Abbreviated version of CIF standard: * Only recognized element symbols may be used. * Each element symbol is followed by a ‘count’ number. A count of ‘1’ may be omitted. * A space or parenthesis must separate each cluster of (element symbol + count). * Where a group of elements is enclosed in parentheses, the multiplier for the group must follow the closing parentheses. That is, all element and group multipliers are assumed to be printed as subscripted numbers. * Unless the elements are ordered in a manner that corresponds to their chemical structure, the order of the elements within any group or moiety depends on whether or not carbon is present. * If carbon is present, the order should be: C, then H, then the other elements in alphabetical order of their symbol. If carbon is not present, the elements are listed purely in alphabetic order of their symbol. This is the Hill system used by Chemical Abstracts. See, for example: http://www.iucr.org/__data/iucr/cif/standard/cifstd15.html, http://www.cas.org/training/stneasytips/subinforformula1.html, or http://www.indiana.edu/~cheminfo/courses/471cnfs.html.

<table>
<thead>
<tr>
<th>Symbol</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>type</td>
<td>NX_CHAR</td>
</tr>
</tbody>
</table>

Type or material of monochromating substance. Chemical formula can be specified separately. Use the “reflection” field to indicate the (hkl) orientation. Use the “d_spacing” field to record the lattice plane spacing.

This field was changed (2010-11-17) from an enumeration to a string since common usage showed a wider variety of use than a simple list. These are the items in the list at the time of the change: PG (Highly Oriented Pyrolytic Graphite) | Ge | Si | Cu | Fe3Si | CoFe | Cu2MnAl (Heusler) | Multilayer | Diamond.

chemical_formula: NX_CHAR
The chemical formula specified using CIF conventions. Abbreviated version of CIF standard:

- Only recognized element symbols may be used.
- Each element symbol is followed by a ‘count’ number. A count of ‘1’ may be omitted.
- A space or parenthesis must separate each cluster of (element symbol + count).
- Where a group of elements is enclosed in parentheses, the multiplier for the group must follow the closing parentheses. That is, all element and group multipliers are assumed to be printed as subscripted numbers.
- Unless the elements are ordered in a manner that corresponds to their chemical structure, the order of the elements within any group or moiety depends on whether or not carbon is present.
- If carbon is present, the order should be: C, then H, then the other elements in alphabetical order of their symbol. If carbon is not present, the elements are listed purely in alphabetical order of their symbol.
- This is the Hill system used by Chemical Abstracts.

**order_no:** `NX_INT`
A number which describes if this is the first, second... nth crystal in a multi crystal monochromator

**cut_angle:** `NX_FLOAT` {units=`NX_ANGLE`}
Cut angle of reflecting Bragg plane and plane of crystal surface

**space_group:** `NX_CHAR`
Space group of crystal structure

**unit_cell[n_comp, 6]:** `NX_FLOAT` {units=`NX_LENGTH`}
Unit cell parameters (lengths and angles)

**unit_cell_a:** `NX_FLOAT` {units=`NX_LENGTH`}
Unit cell lattice parameter: length of side a

**unit_cell_b:** `NX_FLOAT` {units=`NX_LENGTH`}
Unit cell lattice parameter: length of side b

**unit_cell_c:** `NX_FLOAT` {units=`NX_LENGTH`}
Unit cell lattice parameter: length of side c

**unit_cell_alpha:** `NX_FLOAT` {units=`NX_ANGLE`}
Unit cell lattice parameter: angle alpha

**unit_cell_beta:** `NX_FLOAT` {units=`NX_ANGLE`}
Unit cell lattice parameter: angle beta

**unit_cell_gamma:** `NX_FLOAT` {units=`NX_ANGLE`}
Unit cell lattice parameter: angle gamma

**unit_cell_volume:** `NX_FLOAT` {units=`NX_VOLUME`}
Volume of the unit cell

**orientation_matrix[3, 3]:** `NX_FLOAT`
Orientation matrix of single crystal sample using Busing-Levy convention

wavelength[i]: NX_FLOAT {units=NX_WAVELENGTH}
Optimum diffracted wavelength

d_spacing: NX_FLOAT {units=NX_LENGTH}
spacing between crystal planes of the reflection

scattering_vector: NX_FLOAT {units=NX_WAVENUMBER}
Scattering vector, Q, of nominal reflection

reflection[3]: NX_INT {units=NX_UNITLESS}
Miller indices (hkl) values of nominal reflection

thickness: NX_FLOAT {units=NX_LENGTH}
Thickness of the crystal. (Required for Laue orientations - see “usage” field)

density: NX_NUMBER {units=NX_MASS_DENSITY}
mass density of the crystal

segment_width: NX_FLOAT {units=NX_LENGTH}
Horizontal width of individual segment

segment_height: NX_FLOAT {units=NX_LENGTH}
Vertical height of individual segment

segment_thickness: NX_FLOAT {units=NX_LENGTH}
Thickness of individual segment

segment_gap: NX_FLOAT {units=NX_LENGTH}
Typical gap between adjacent segments

segment_columns: NX_FLOAT {units=NX_LENGTH}
number of segment columns in horizontal direction

segment_rows: NX_FLOAT {units=NX_LENGTH}
number of segment rows in vertical direction

mosaic_horizontal: NX_FLOAT {units=NX_ANGLE}
horizontal mosaic Full Width Half Maximum

mosaic_vertical: NX_FLOAT {units=NX_ANGLE}
vertical mosaic Full Width Half Maximum

curvature_horizontal: NX_FLOAT {units=NX_ANGLE}
Horizontal curvature of focusing crystal

curvature_vertical: NX_FLOAT {units=NX_ANGLE}
Vertical curvature of focusing crystal

is_cylindrical: NX_BOOLEAN
Is this crystal bent cylindrically?

cylindrical_orientation_angle: NX_NUMBER {units=NX_ANGLE}
If cylindrical: cylinder orientation angle

**polar_angle[i]**: *NX_FLOAT* \{units=*NX_ANGLE*\}

Polar (scattering) angle at which crystal assembly is positioned. Note: some instrument geometries call this term 2theta.

**azimuthal_angle[i]**: *NX_FLOAT* \{units=*NX_ANGLE*\}

Azimuthal angle at which crystal assembly is positioned

**bragg_angle[i]**: *NX_FLOAT* \{units=*NX_ANGLE*\}

Bragg angle of nominal reflection

**temperature**: *NX_FLOAT* \{units=*NX_TEMPERATURE*\}

average/nominal crystal temperature

**temperature_coefficient**: *NX_FLOAT* \{units=*NX_ANY*\}

how lattice parameter changes with temperature

(geometries): *NXshape*

Position of crystal

**temperature_log**: *NXlog*

log file of crystal temperature

**reflectivity**: *NXdata*

crystal reflectivity versus wavelength

**transmission**: *NXdata*

crystal transmission versus wavelength

**shape**: *NXshape*

A NXshape group describing the shape of the crystal arrangement

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXcrystal.nxdl.xml

### 3.3.11 NXdata

**Status:**

base class, extends *NXobject*, version 1.0

**Description:**

(required) *NXdata* is a template of plottable data and their dimension scales. It is mandatory that there is at least one *NXdata* group in each NXentry group. Note that the variable and data can be defined with different names. The signal and axes attribute of the data item define which items are plottable data and which are dimension scales.

- Each *NXdata* group will consist of only one data set containing plottable data and their standard deviations.
- This data set may be of arbitrary rank up to a maximum of *NX_MAXRANK=32*.
- The plottable data will be identified by the attribute: signal=1
- The plottable data will identify the dimension scale specification(s) in the axes attribute.
If available, the standard deviations of the data are to be stored in a data set of the same rank and dimensions, with the name errors.

- For each data dimension, there should be a one-dimensional array of the same length.
- These one-dimensional arrays are the dimension scales of the data, i.e. the values of the independent variables at which the data is measured, such as scattering angle or energy transfer.

There are two methods of linking each data dimension to its respective dimension scale.

The preferred (and recommended) method uses the axes attribute to specify the names of each dimension scale.

The older method uses the axis attribute on each dimension scale to identify with an integer the axis whose value is the number of the dimension.

NXdata is used to implement one of the basic motivations in NeXus, to provide a default plot for the data of this NXentry. The actual data might be stored in another group and (hard) linked to the NXdata group.

Symbols:

These symbols will be used below to coordinate datasets with the same shape.

- dataRank: rank of the data field
- n: length of the variable field
- nx: length of the x field
- ny: length of the y field
- nz: length of the z field

Groups cited: none

Structure:

variable[n]: NX_NUMBER

Dimension scale defining an axis of the data. Client is responsible for defining the dimensions of the data. The name of this field may be changed to fit the circumstances. Standard NeXus client tools will use the attributes to determine how to use this field.

@long_name: NX_CHAR

Axis label

@distribution: NX_BOOLEAN

0|false: single value, 1|true: multiple values

@first_good: NX_INT

Index of first good value

@last_good: NX_INT

Index of last good value

@axis: NX_POSINT

Index (positive integer) identifying this specific set of numbers. N.B. The axis attribute is the old way of designating a link. Do not use the axes attribute with the axis attribute. The axes attribute is now preferred.

variable_errors[n]: NX_NUMBER
Errors (uncertainties) associated with axis variable Client is responsible for defining the dimensions of the data. The name of this field may be changed to fit the circumstances but is matched with the variable field with _errors appended.

data[n]: NX_NUMBER
This field contains the data values to be used as the NeXus plottable data. Client is responsible for defining the dimensions of the data. The name of this field may be changed to fit the circumstances. Standard NeXus client tools will use the attributes to determine how to use this field.

@signal: NX_POSINT
Plottable (independent) axis, indicate index number. Only one field in a NXdata group may have the signal=1 attribute. Do not use the signal attribute with the axis attribute.

@axes: NX_CHAR
Defines the names of the dimension scales (independent axes) for this data set as a colon-delimited array. NOTE: The axes attribute is the preferred method of designating a link. Do not use the axes attribute with the axis attribute.

@uncertainties: NX_CHAR
Specify the names of the errors (uncertainties) of the dependent axes as plottable data. NOTE: The errors attribute uses the same syntax as the axes attribute.

@long_name: NX_CHAR
data label

errors[n]: NX_NUMBER
Standard deviations of data values - the data array is identified by the attribute signal=1. The errors array must have the same dimensions as data. Client is responsible for defining the dimensions of the data.

scaling_factor: NX_FLOAT
The elements in data are usually float values really. For efficiency reasons these are usually stored as integers after scaling with a scale factor. This value is the scale factor. It is required to get the actual physical value, when necessary.

offset: NX_FLOAT
An optional offset to apply to the values in data.

x[nx]: NX_FLOAT {units=NX_ANY}
This is an array holding the values to use for the x-axis of data. The units must be appropriate for the measurement.

y[ny]: NX_FLOAT {units=NX_ANY}
This is an array holding the values to use for the y-axis of data. The units must be appropriate for the measurement.

z[nz]: NX_FLOAT {units=NX_ANY}
This is an array holding the values to use for the z-axis of data. The units must be appropriate for the measurement.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXdata.nxdl.xml
### 3.3.12 NXdetector

**Status:**

base class, extends `NXobject`, version 1.1

**Description:**

Template of a detector, detector bank, or multidetector.

**Symbols:**

These symbols will be used below to coordinate datasets with the same shape.

- `np`: number of scan points (only present in scanning measurements)
- `i`: number of detector pixels in the first (X, slowest) direction
- `j`: number of detector pixels in the second (Y, faster) direction
- `k`: number of detector pixels in the third (Z, if necessary, fastest) direction
- `tof`: number of bins in the time-of-flight histogram

**Groups cited:** `NXcharacterization`, `NXcollection`, `NXdata`, `NXdetector_module`, `NXgeometry`, `NXnote`

**Structure:**

```plaintext
time_of_flight[tof+1]: NX_FLOAT {units=NX_TIME_OF_FLIGHT}
    Total time of flight
    @axis: NX_POSINT
        Obligatory value: 3
    @primary: NX_POSINT
        Obligatory value: 1
    @long_name: NX_CHAR
        Axis label
    @link: NX_CHAR
        absolute path to location in NXdetector

raw_time_of_flight[tof+1]: NX_INT {units=NX_PULSES}
    In DAQ clock pulses
    @frequency: NX_NUMBER
        Clock frequency in Hz

detector_number[i, j]: NX_INT
    Identifier for detector

data[np, i, j, tof]: NX_NUMBER {units=NX_ANY}
    Data values from the detector.
    @signal: NX_POSINT
        Obligatory value: 1
    @axes: NX_CHAR
        [number of scan points,x_offset?,y_offset?,time_of_flight?]```
@long_name: NX_CHAR
Title of measurement
@check_sum: NX_INT
Integral of data as check of data integrity
@link: NX_CHAR
absolute path to location in NXdetector
data_error[np, i, j, tof]: NX_NUMBER {units=NX_ANY}
The best estimate of the uncertainty in the data value. Where possible, this should be the
standard deviation, which has the same units as the data.
@units: NX_CHAR
@link: NX_CHAR
absolute path to location in NXdetector
x_pixel_offset[i, j]:NX_FLOAT {units=NX_LENGTH}
Offset from the detector center in x-direction. Can be multidimensional when needed.
@axis: NX_POSINT
Obligatory value: 1
@primary: NX_POSINT
Obligatory value: 1
@long_name: NX_CHAR
Axis label
@link: NX_CHAR
absolute path to location in NXdetector
y_pixel_offset[i, j]:NX_FLOAT {units=NX_LENGTH}
Offset from the detector center in the y-direction. Can be multidimensional when different
values are required for each pixel.
@axis: NX_POSINT
Obligatory value: 2
@primary: NX_POSINT
Obligatory value: 1
@long_name: NX_CHAR
Axis label
distance[np, i, j]: NX_FLOAT {units=NX_LENGTH}
TODO: need documentation
polar_angle[np, i, j]: NX_FLOAT {units=NX_ANGLE}
TODO: need documentation
azimuthal_angle[np, i, j]: NX_FLOAT {units=NX_ANGLE}
TODO: need documentation
**description**: *NX_CHAR*

name/manufacturer/model/etc. information

**local_name**: *NX_CHAR*

Local name for the detector

**solid_angle[i, j]**: *NX_FLOAT* {units=*NX_SOLID_ANGLE*}

Solid angle subtended by the detector at the sample

**x_pixel_size[i, j]**: *NX_FLOAT* {units=*NX_LENGTH*}

Size of each detector pixel. If it is scalar all pixels are the same size.

**y_pixel_size[i, j]**: *NX_FLOAT* {units=*NX_LENGTH*}

Size of each detector pixel. If it is scalar all pixels are the same size

**dead_time[np, i, j]**: *NX_FLOAT* {units=*NX_TIME*}

Detector dead time

**gas_pressure[i, j]**: *NX_FLOAT* {units=*NX_PRESSURE*}

Detector gas pressure

**detection_gas_path**: *NX_FLOAT* {units=*NX_LENGTH*}

maximum drift space dimension

**crate[i, j]**: *NX_INT*

Crate number of detector

@**local_name**: *NX_CHAR*

Equivalent local term

**slot[i, j]**: *NX_INT*

Slot number of detector

@**local_name**: *NX_CHAR*

Equivalent local term

**input[i, j]**: *NX_INT*

Input number of detector

@**local_name**: *NX_CHAR*

Equivalent local term

**type**: *NX_CHAR*

Description of type such as He3 gas cylinder, He3 PSD, scintillator, fission chamber, proportion counter, ion chamber, ccd, pixel, image plate, CMOS, ...

**calibration_date**: *NX_DATE_TIME*

date of last calibration (geometry and/or efficiency) measurements

**layout**: *NX_CHAR*

How the detector is represented

Any of these values: *point* | *linear* | *area*
count_time[np]: NX_NUMBER {units=NX_TIME}
    Elapsed actual counting time

sequence_number[nBrightFrames]: NX_CHAR
    In order to properly sort the order of the images taken in (for example) a tomography experi-
    ment, a sequence number is stored with each image.

beam_center_x: NX_FLOAT {units=NX_LENGTH}
    This is the x position where the direct beam would hit the detector. This is a length, not a pixel
    position, and can be outside of the actual detector.

beam_center_y: NX_FLOAT {units=NX_LENGTH}
    This is the y position where the direct beam would hit the detector. This is a length, not a pixel
    position, and can be outside of the actual detector.

frame_start_number: NX_INT
    This is the start number of the first frame of a scan. In PX one often scans a couple of frames on
    a give sample, then does something else, then returns to the same sample and scans some more
    frames. Each time with a new data file. This number helps concatenating such measurements.

diameter: NX_FLOAT {units=NX_LENGTH}
    The diameter of a cylindrical detector

acquisition_mode: NX_CHAR
    The acquisition mode of the detector.
    Any of these values:
    • gated
    • triggered
    • summed
    • event
    • histogrammed

angular_calibration_applied: NX_BOOLEAN
    True when the angular calibration has been applied in the electronics, false otherwise.

angular_calibration[i, j]: NX_FLOAT
    Angular calibration data.

flatfield_applied: NX_BOOLEAN
    True when the flat field correction has been applied in the electronics, false otherwise.

flatfield[i, j]: NX_FLOAT
    Flat field correction data.

flatfield_error[i, j]: NX_FLOAT
    Errors of the flat field correction data.

pixel_mask_applied: NX_BOOLEAN
    True when the pixel mask correction has been applied in the electronics, false otherwise.

pixel_mask[i, j]: NX_INT
The 32-bit pixel mask for the detector. Contains a bit field for each pixel to signal dead, blind or high or otherwise unwanted or undesirable pixels. They have the following meaning:

- bit 0: gap (pixel with no sensor)
- bit 1: dead
- bit 2: under responding
- bit 3: over responding
- bit 4: noisy
- bit 5: undefined
- bit 6: pixel is part of a cluster of problematic pixels (bit set in addition to others)
- bit 7: undefined
- bit 8: user defined mask (e.g. around beamstop)
- bits 9-30: undefined
- bit 31: virtual pixel (corner pixel with interpolated value)

The normal data analysis software would not take pixels into account when a bit in (mask & 0x00FF) is set. Tag bit in the upper two bytes would indicate special pixel properties that normally would not be a sole reason to reject the intensity value (unless lower bits are set as well of course).

count_rate_correction_applied: NX_BOOLEAN

True when a count-rate correction has already been applied in the electronics, false otherwise.

bit_depth_readout: NX_INT

How many bits the electronics reads per pixel. With CCD’s and single photon counting detectors, this must not align with traditional integer sizes. This can be 4, 8, 12, 14, 16, ...

detector_readout_time: NX_FLOAT {units=NX_TIME}

Time it takes to read the detector (typically milliseconds). This is important to know for time resolved experiments.

trigger_delay_time: NX_FLOAT {units=NX_TIME}

Time it takes to start exposure after a trigger signal has been received. This is important to know for time resolved experiments.

trigger_dead_time: NX_FLOAT {units=NX_TIME}

Time during which no new trigger signal can be accepted. Typically this is the trigger_delay_time + exposure_time + readout_time. This is important to know for time resolved experiments.

frame_time[NP]: NX_FLOAT {units=NX_TIME}

This is time for each frame. This is exposure_time + readout_time.

gain_setting: NX_CHAR

The gain setting of the detector. This influences background etc.

Any of these values: high | standard | fast | auto

saturation_value: NX_INT
The value at which the detector goes into saturation. Especially common to CCD detectors, the data is known to be invalid above this value.

**number_of_cycles**: *NX_INT*

CCD images are sometimes constructed by summing together multiple short exposures in the electronics. This reduces background etc. This is the number of short exposures used to sum images for an image.

**sensor_material**: *NX_CHAR*

At times, radiation is not directly sensed by the detector. Rather, the detector might sense the output from some converter like a scintillator. This is the name of this converter material.

**sensor_thickness**: *NX_FLOAT {units=NX_LENGTH}*

At times, radiation is not directly sensed by the detector. Rather, the detector might sense the output from some converter like a scintillator. This is the thickness of this converter material.

**threshold_energy**: *NX_FLOAT {units=NX_ENERGY}*

Single photon counter detectors can be adjusted for a certain energy range in which they work optimally. This is the energy setting for this.

**(geometry)**: *NXgeometry*

Position and orientation of detector

**efficiency**: *NXdata*

Spectral efficiency of detector with respect to e.g. wavelength

**efficiency[i, j, k]**: *NX_FLOAT {units=NX_DIMENSIONLESS}*

efficiency of the detector

**wavelength[i, j, k]**: *NX_FLOAT {units=NX_WAVELENGTH}*

TODO: need documentation

**start_time[np]**: *NX_FLOAT {units=NX_TIME}*

start time for each frame, with the `start` attribute as absolute reference

@`start`: *NX_DATE_TIME*

**stop_time[np]**: *NX_FLOAT {units=NX_TIME}*

stop time for each frame, with the `start` attribute as absolute reference

@`start`: *NX_DATE_TIME*

**real_time[i, j, k]**: *NX_NUMBER {units=NX_TIME}*

real-time of the exposure (use this if exposure time varies for each array element, otherwise use `count_time` field)

**calibration_method**: *NXnote*

summary of conversion of array data to pixels (e.g. polynomial approximations) and location of details of the calibrations

**data_file**: *NXnote*

**(characterization)**: *NXcharacterization*

deprecated, use NXcollection instead

for more details, see https://github.com/nexusformat/definitions/issues/177
(collection): NXcollection

Use this group to provide other data related to this NXdetector group.

(detector_module): NXdetector_module

For use in special cases where the data in NXdetector is represented in several parts, each with a separate geometry.

Use one or more instances of the NXdetector_module group to declare regions of interest or some other subdivision of a detector.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXdetector.nxdl.xml

3.3.13 NXdetector_group

Status:

base class, extends NXobject, version 1.0

Description:

This class is used to allow a logical grouping of detector elements (e.g. which tube, bank or group of banks) to be recorded in the file. As well as allowing you to e.g just select the “left” or “east” detectors, it may also be useful for determining which elements belong to the same PSD tube and hence have e.g. the same dead time.

For example, if we had “bank1” composed of “tube1”, “tube2” and “tube3” then group_names would be the string “bank1,bank1/tube1,bank1/tube2,bank1/tube3” group_index would be {1,2,3,4} group_parent would be {-1,1,1,1}

The mapping array is interpreted as group 1 is a top level group containing groups 2, 3 and 4

A group_index array in NXdetector gives the base group for a detector element.

Symbols:

No symbol table

Groups cited: none

Structure:

  group_names: NX_CHAR
    Comma separated list of names

  group_index[i]: NX_INT
    Unique ID for group. A group_index array in NXdetector gives the base group for a detector element.

  group_parent[ref(group_index)]: NX_INT
    Index of group parent in the hierarchy: -1 means no parent (i.e. a top level) group

  group_type[ref(group_index)]: NX_INT
    Code number for group type, e.g. bank=1, tube=2 etc.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXdetector_group.nxdl.xml
3.3.14 NXdetector_module

Status:

base class, extends NXobject, version 1.0

Description:

This is the description of a detector module. Many detectors consist of multiple smaller modules. Sometimes it is important to know the exact position of such modules. This is the purpose of this group. It is a child group to NXdetector.

Symbols:

No symbol table

Groups cited: none

Structure:

- **data_origin**: NX_INT
  A two value field which gives the index of the start of the modules data in the main area detector image in the underlying NXdetector module.

- **data_size**: NX_INT
  Two values for the size of the module in pixels in each direction.

- **module_offset**: NX_NUMBER \{units=NX_LENGTH\}
  Offset of the module in regards to the origin of the detector in an arbitrary direction.

  - **transformation_type**: NX_CHAR
    Obligatory value: translation

  - **vector**: NX_NUMBER
    Three values that define the axis for this transformation

  - **offset**: NX_NUMBER
    A fixed offset applied before the transformation

  - **offset_units**: NX_CHAR
    Units of the offset.

  - **depends_on**: NX_CHAR
    Points to the path of the next element in the geometry chain.

- **fast_pixel_direction**: NX_NUMBER \{units=NX_LENGTH\}
  Values along the direction of fastest varying pixel direction. The direction itself is given through the vector attribute

  - **transformation_type**: NX_CHAR
    Obligatory value: translation

  - **vector**: NX_NUMBER
    Three values that define the axis for this transformation

  - **offset**: NX_NUMBER
    A fixed offset applied before the transformation
slow_pixel_direction: NX_NUMBER {units=NX_LENGTH}

Values along the direction of slow varying pixel direction. The direction itself is given through the vector attribute

@transformation_type: NX_CHAR

Obligatory value: translation

@vector: NX_NUMBER

Three values that define the axis for this transformation

@offset: NX_NUMBER

A fixed offset applied before the transformation

@offset_units: NX_CHAR

Units of the offset.

@depends_on: NX_CHAR

Points to the path of the next element in the geometry chain.

fast_pixel_size: NX_NUMBER {units=NX_LENGTH}

Values along the direction of fastest varying pixel direction. The direction itself is given through the vector attribute

@transformation_type: NX_CHAR

Obligatory value: translation

@vector: NX_NUMBER

Three values that define the axis for this transformation

@offset: NX_NUMBER

A fixed offset applied before the transformation

@offset_units: NX_CHAR

Units of the offset.

@depends_on: NX_CHAR

Points to the path of the next element in the geometry chain.

slow_pixel_size: NX_NUMBER {units=NX_LENGTH}

Values along the direction of slow varying pixel direction. The direction itself is given through the vector attribute

@transformation_type: NX_CHAR

Obligatory value: translation

@vector: NX_NUMBER

Three values that define the axis for this transformation
@offset: NX_NUMBER
   A fixed offset applied before the transformation

@offset_units: NX_CHAR
   Units of the offset.

@depends_on: NX_CHAR
   Points to the path of the next element in the geometry chain.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXdetector_module.nxdl.xml

### 3.3.15 NXdisk_chopper

**Status:**
- base class, extends NXobject, version 1.0

**Description:**
- TODO: need documentation

**Symbols:**
- No symbol table

**Groups cited:** NXgeometry

**Structure:**
- **type:** NX_CHAR
  - Type of the disk-chopper: only one from the enumerated list (match text exactly)
  - Any of these values:
    - Chopper type single
    - contra_rotating_pair
    - synchro_pair

- **rotation_speed:** NX_FLOAT {units=NX_FREQUENCY}
  - chopper rotation speed

- **slits:** NX_INT
  - Number of slits

- **slit_angle:** NX_FLOAT {units=NX_ANGLE}
  - angular opening

- **pair_separation:** NX_FLOAT {units=NX_LENGTH}
  - disc spacing in direction of beam

- **radius:** NX_FLOAT {units=NX_LENGTH}
  - radius to centre of slit

- **slit_height:** NX_FLOAT {units=NX_LENGTH}
  - total slit height

- **phase:** NX_FLOAT {units=NX_ANGLE}
chopper phase angle

**ratio:** `NX_INT`

pulse reduction factor of this chopper in relation to other choppers/fastest pulse in the instrument

**distance:** `NX_FLOAT {units=NX_LENGTH}`

Effective distance to the origin

**wavelength_range[2]:** `NX_FLOAT {units=NX_WAVELENGTH}`

low and high values of wavelength range transmitted

(geometrical: `NXgeometry`

**Source:** Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXdisk_chopper.nxdl.xml

### 3.3.16 NXentry

**Status:**

base class, extends `NXobject`, version 1.0

**Description:**

(required) Template of the top-level NeXus group which contains all the data and associated information that comprise a single measurement. It is mandatory that there is at least one group of this type in the NeXus file.

**Symbols:**

No symbol table

**Groups cited:** `NXcharacterization, NXcollection, NXdata, NXinstrument, NXmonitor, NXnote, NXparameters, NXprocess, NXsample, NXsubentry, NXuser`

**Structure:**

@IDF_Version: `NX_CHAR`

ISIS Muon IDF_Version

title: `NX_CHAR`

Extended title for entry

experiment_identifier: `NX_CHAR`

Unique identifier for the experiment, defined by the facility, possibly linked to the proposals

experiment_description: `NX_CHAR`

Brief summary of the experiment, including key objectives.

collection_identifier: `NX_CHAR`

User or Data Acquisition defined group of NeXus files or NXentry

collection_description: `NX_CHAR`

Brief summary of the collection, including grouping criteria.

type_identifier: `NX_CHAR`

unique identifier for the measurement, defined by the facility.
definition: *NX_CHAR*
(alternate use: see same field in *NXsubentry* for preferred)
Official NeXus NXDL schema to which this file conforms.
This field is provided so that *NXentry* can be the overlay position in a NeXus data file for an
application definition and its set of groups, fields, and attributes.

*It is advised to use* *NXsubentry*, *instead, as the overlay position.*

@version: *NX_CHAR*
NXDL version number

@URL: *NX_CHAR*
URL of NXDL file

definition_local: *NX_CHAR*
(deprecated use: see same field in *NXsubentry* for preferred) Local NXDL schema extended
from the file specified in the *definition* field. This contains any locally-defined, additional
fields in the file.

@version: *NX_CHAR*
NXDL version number

@URL: *NX_CHAR*
URL of NXDL file

start_time: *NX_DATE_TIME*
Starting time of measurement

duration: *NX_INT* *{units=NX_TIME}*
Duration of measurement

collection_time: *NX_FLOAT* *{units=NX_TIME}*
Time transpired actually collecting data i.e. taking out time when collection was suspended
due to e.g. temperature out of range

run_cycle: *NX_CHAR*
Such as “2007-3”. Some user facilities organize their beam time into run cycles.

program_name: *NX_CHAR*
Name of program used to generate this file

@version: *NX_CHAR*
Program version number

@configuration: *NX_CHAR*
configuration of the program

revision: *NX_CHAR*
Revision id of the file due to re-calibration, reprocessing, new analysis, new instrument definition format, ...

@comment: NX_CHAR

pre_sample_flightpath: NX_FLOAT {units=NX_LENGTH}

This is the flightpath before the sample position. This can be determined by a chopper, by the moderator or the source itself. In other words: it the distance to the component which gives the T0 signal to the detector electronics. If another component in the NXinstrument hierarchy provides this information, this should be a link.

(data): NXdata

The required data group

experiment_documentation: NXnote

Description of the full experiment (document in pdf, latex, ...)

notes: NXnote

Notes describing entry

thumbnail: NXnote

A small image that is representative of the entry. An example of this is a 640x480 jpeg image automatically produced by a low resolution plot of the NXdata.

@mimeType: NX_CHAR

The value should be an image/*

Obligatory value: image/*

(characterization): NXcharacterization

(user): NXuser

(sample): NXsample

(instrument): NXinstrument

(collection): NXcollection

(monitor): NXmonitor

(parameters): NXparameters

(process): NXprocess

(subentry): NXsubentry

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXentry.nxdl.xml

3.3.17 NXenvironment

Status:

base class, extends NXobject, version 1.0

Description:

This class describes an external condition applied to the sample

Symbols:

No symbol table
Groups cited:  *NXgeometry*, *NXnote*, *NXsensor*

Structure:

**name**: *NX_CHAR*

Apparatus identification code/model number; e.g. OC100 011

**short_name**: *NX_CHAR*

Alternative short name, perhaps for dashboard display like a present Seblock name

**type**: *NX_CHAR*

Type of apparatus. This could be the SE codes in scheduling database; e.g. OC/100

**description**: *NX_CHAR*

Description of the apparatus; e.g. 100mm bore orange cryostat with Roots pump

**program**: *NX_CHAR*

Program controlling the apparatus; e.g. LabView VI name

**position**: *NXgeometry*

The position and orientation of the apparatus

**(note)**: *NXnote*

Additional information, LabView logs, digital photographs, etc

**(sensor)**: *NXsensor*

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXenvironment.nxdl.xml

### 3.3.18 NXevent_data

**Status**:

base class, extends *NXObject*, version 1.0

**Description**:

Time-of-flight events

**Symbols**:

No symbol table

**Groups cited**: none

**Structure**:

**time_of_flight[i]**: *NX_INT* {units=*NX_TIME_OF_FLIGHT*}

A list of time of flight for each event as it comes in. This list is for all pulses with information to attach to a particular pulse located in events_per_pulse.

**pixel_number[i]**: *NX_INT* {units=*NX_DIMENSIONLESS*}

There will be extra information in the NXdetector to convert pixel_number to detector_number. This list is for all pulses with information to attach to a particular pulse located in events_per_pulse.

**pulse_time[j]**: *NX_INT* {units=*NX_TIME*}
The time that each pulse started with respect to the offset

@offset: NX_DATE_TIME

ISO8601

events_per_pulse[j]: NX_INT \{units=NX_DIMENSIONLESS\}

This connects the index “i” to the index “j”. The jth element is the number of events in “i” that occurred during the jth pulse.

pulse_height[i, k]: NX_FLOAT \{units=NX_DIMENSIONLESS\}

If voltages from the ends of the detector are read out this is where they go. This list is for all events with information to attach to a particular pulse height. The information to attach to a particular pulse is located in events_per_pulse.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXevent_data.nxdl.xml

3.3.19 NXfermi_chopper

Status:

base class, extends NXobject, version 1.0

Description:

Description of a Fermi chopper, possibly with curved slits.

Symbols:

No symbol table

Groups cited: NXgeometry

Structure:

type: NX_CHAR

Fermi chopper type

rotation_speed: NX_FLOAT \{units=NX_FREQUENCY\}

chopper rotation speed

radius: NX_FLOAT \{units=NX_LENGTH\}

radius of chopper

slit: NX_FLOAT \{units=NX_LENGTH\}

width of an individual slit

r_slit: NX_FLOAT \{units=NX_LENGTH\}

radius of curvature of slits

number: NX_INT \{units=NX_UNITLESS\}

number of slits

height: NX_FLOAT \{units=NX_LENGTH\}

input beam height

width: NX_FLOAT \{units=NX_LENGTH\}

input beam width
3.3.20 NXfilter

Status:

base class, extends NXobject, version 1.0

Description:

Template for specifying the state of band pass filters. If uncertain whether to use NXfilter (band-pass filter) orNXattenuator (reduces beam intensity), then use NXattenuator.

Symbols:

No symbol table

Groups cited: NXdata, NXgeometry, NXlog, NXsensor

Structure:

description: NX_CHAR

Composition of the filter. Chemical formula can be specified separately.

This field was changed (2010-11-17) from an enumeration to a string since common usage showed a wider variety of use than a simple list. These are the items in the list at the time of the change: Beryllium | Pyrolytic Graphite | Graphite | Sapphire | Silicon | Supermirror.

status: NX_CHAR

position with respect to in or out of the beam (choice of only “in” or “out”)

Any of these values:

• in: in the beam
• out: out of the beam

temperature: NX_FLOAT {units=NX_TEMPERATURE}

average/nominal filter temperature

thickness: NX_FLOAT {units=NX_LENGTH}
Thickness of the filter

density: **NX_NUMBER** {units=**NX_MASS_DENSITY**}

mass density of the filter

chemical_formula: **NX_CHAR**

The chemical formula specified using CIF conventions. Abbreviated version of CIF standard:

- Only recognized element symbols may be used.
- Each element symbol is followed by a ‘count’ number. A count of ‘1’ may be omitted.
- A space or parenthesis must separate each cluster of (element symbol + count).
- Where a group of elements is enclosed in parentheses, the multiplier for the group must follow the closing parentheses. That is, all element and group multipliers are assumed to be printed as subscripted numbers.
- Unless the elements are ordered in a manner that corresponds to their chemical structure, the order of the elements within any group or moiety depends on whether or not carbon is present.
- If carbon is present, the order should be:
  - C, then H, then the other elements in alphabetical order of their symbol.
  - If carbon is not present, the elements are listed purely in alphabetic order of their symbol.
- This is the Hill system used by Chemical Abstracts.

unit_cell_a: **NX_FLOAT** {units=**NX_LENGTH**}

Unit cell lattice parameter: length of side a

unit_cell_b: **NX_FLOAT** {units=**NX_LENGTH**}

Unit cell lattice parameter: length of side b

unit_cell_c: **NX_FLOAT** {units=**NX_LENGTH**}

Unit cell lattice parameter: length of side c

unit_cell_alpha: **NX_FLOAT** {units=**NX_ANGLE**}

Unit cell lattice parameter: angle alpha

unit_cell_beta: **NX_FLOAT** {units=**NX_ANGLE**}

Unit cell lattice parameter: angle beta

unit_cell_gamma: **NX_FLOAT** {units=**NX_ANGLE**}

Unit cell lattice parameter: angle gamma

unit_cell_volume[n_comp]: **NX_FLOAT** {units=**NX_VOLUME**}

Unit cell

orientation_matrix[n_comp, 3, 3]: **NX_FLOAT**

Orientation matrix of single crystal filter using Busing-Levy convention

m_value: **NX_FLOAT** {units=**NX_DIMENSIONLESS**}

m value of supermirror filter

substrate_material: **NX_CHAR**
substrate material of supermirror filter

**substrate_thickness**: `NX_FLOAT {units=NX_LENGTH}`

substrate thickness of supermirror filter

**coating_material**: `NX_CHAR`

coating material of supermirror filter

**substrate_roughness**: `NX_FLOAT {units=NX_LENGTH}`

substrate roughness (RMS) of supermirror filter

**coating_roughness[nsurf]**: `NX_FLOAT {units=NX_LENGTH}`

coating roughness (RMS) of supermirror filter

**(geometry)**: `NXgeometry`

Geometry of the filter

**transmission**: `NXdata`

Wavelength transmission profile of filter

**temperature_log**: `NXlog`

Linked temperature_log for the filter

**sensor_type**: `NXsensor`

Sensor(s) used to monitor the filter temperature

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXfilter.nxdl.xml

### 3.3.21 NXflipper

**Status:**

base class, extends `NXObject`, version 1.0

**Description:**

Template of a beamline spin flipper.

**Symbols:**

No symbol table

**Groups cited:** none

**Structure:**

- **type**: `NX_CHAR`
  
  Any of these values: `coil` | `current-sheet`

- **flip_turns**: `NX_FLOAT {units=NX_PER_LENGTH}`
  
  Linear density of turns (such as number of turns/cm) in flipping field coils

- **comp_turns**: `NX_FLOAT {units=NX_PER_LENGTH}`
  
  Linear density of turns (such as number of turns/cm) in compensating field coils

- **guide_turns**: `NX_FLOAT {units=NX_PER_LENGTH}`
  
  Linear density of turns (such as number of turns/cm) in guide field coils
flip_current: \textit{NX_FLOAT} \{units=\textit{NX_CURRENT}\}

Flipping field coil current in “on” state

comp_current: \textit{NX_FLOAT} \{units=\textit{NX_CURRENT}\}

Compensating field coil current in “on” state

guide_current: \textit{NX_FLOAT} \{units=\textit{NX_CURRENT}\}

Guide field coil current in “on” state

thickness: \textit{NX_FLOAT} \{units=\textit{NX_LENGTH}\}

thickness along path of neutron travel

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXflipper.nxdl.xml

3.3.22 NXgeometry

Status:

base class, extends \textit{NXObject}, version 1.0

Description:

This is the description for a general position of a component. It is recommended to name an instance of NXgeometry as “geometry” to aid in the use of the definition in simulation codes such as McStas. Also, in HDF, linked items must share the same name. However, it might not be possible or practical in all situations.

Symbols:

No symbol table

Groups cited: \textit{NXorientation}, \textit{NXshape}, \textit{NXtranslation}

Structure:

\textbf{description: \textit{NX_CHAR}}

Optional description/label. Probably only present if we are an additional reference point for components rather than the location of a real component.

\textbf{component_index: \textit{NX_INT}}

Position of the component along the beam path. The sample is at 0, components upstream have negative component_index, components downstream have positive component_index.

\textbf{(shape): \textit{NXshape}}

shape/size information of component

\textbf{(translation): \textit{NXtranslation}}

translation of component

\textbf{(orientation): \textit{NXorientation}}

orientation of component

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXgeometry.nxdl.xml
3.3.23 NXguide

Status:

base class, extends NXobject, version 1.0

Description:

NXguide is used by neutron instruments to describe a guide consists of several mirrors building a shape through which neutrons can be guided or directed. The simplest such form is box shaped although elliptical guides are gaining in popularity. The individual parts of a guide usually have common characteristics but there are cases where they are different. For example, a neutron guide might consist of 2 or 4 coated walls or a supermirror bender with multiple, coated vanes.

To describe polarizing supermirrors such as used in neutron reflection, it may be necessary to revise this definition of NXguide to include NXpolarizer and/or NXmirror.

When even greater complexity exists in the definition of what constitutes a guide, it has been suggested that NXguide be redefined as a NXcollection of NXmirrors each having their own NXgeometries describing their location(s).

For the more general case when describing mirrors, consider using NXmirror.

NOTE: The NeXus International Advisory Committee welcomes comments for revision and improvement of this definition of NXguide.

Symbols:

nsurf: number of reflecting surfaces
nwl: number of wavelengths

Groups cited: NXdata, NXgeometry

Structure:

description: NX_CHAR
   A description of this particular instance of NXguide.
incident_angle: NX_FLOAT {units=NX_ANGLE}
   TODO: documentation needed
bend_angle_x: NX_FLOAT {units=NX_ANGLE}
   TODO: documentation needed
bend_angle_y: NX_FLOAT {units=NX_ANGLE}
   TODO: documentation needed
interior_atmosphere: NX_CHAR
   Any of these values: vacuum|helium|argon
external_material: NX_CHAR
   external material outside substrate
m_value[nsurf]: NX_FLOAT
   The m value for a supermirror, which defines the supermirror regime in multiples of the critical angle of Nickel.
substrate_material[nsurf]: NX_FLOAT
   TODO: documentation needed
substrate_thickness[nsurf]: `NX_FLOAT {units=NX_LENGTH}`

TODO: documentation needed

coating_material[nsurf]: `NX_FLOAT`

TODO: documentation needed

substrate_roughness[nsurf]: `NX_FLOAT {units=NX_LENGTH}`

TODO: documentation needed

coating_roughness[nsurf]: `NX_FLOAT {units=NX_LENGTH}`

TODO: documentation needed

number_sections: `NX_INT {units=NX_UNITLESS}`

number of substrate sections (also called nsurf as an index in the NXguide specification)

(geometry): `NXgeometry`

TODO: Explain what this NXgeometry group means. What is intended here?

reflectivity: `NXdata`

Reflectivity as function of reflecting surface and wavelength

data[nsurf, nwl]: `NX_NUMBER`

reflectivity of each surface as a function of wavelength

@signal: `NX_POSINT`

Use signal=1 to indicate that this is the plottable data for NeXus.

Obligatory value: 1

@axes: `NX_CHAR`

Use axes="surface:wavelength" to indicate the dimension scales to be used when plotting this data.

Obligatory value: surface:wavelength

surface[nsurf]: `NX_NUMBER {units=NX_ANY}`

List of surfaces. Probably best to use index numbers but the specification is very loose.

wavelength[nwl]: `NX_NUMBER {units=NX_WAVELENGTH}`

wavelengths at which reflectivity was measured

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXguide.nxdl.xml

3.3.24 NXinsertion_device

Status:

base class, extends `NXobject`, version 1.0

Description:

Description of an insertion device, as in a synchrotron.

Symbols:

No symbol table
Groups cited:  *NXdata, NXgeometry*

Structure:

- **type**: *NX_CHAR*
  - Any of these values: undulator, wiggler
- **gap**: *NX_FLOAT* \{units=*NX_LENGTH*\}
  - separation between opposing pairs of magnetic poles
- **taper**: *NX_FLOAT* \{units=*NX_ANGLE*\}
  - angular of gap difference between upstream and downstream ends of the insertion device
- **phase**: *NX_FLOAT* \{units=*NX_ANGLE*\}
- **poles**: *NX_INT* \{units=*NX_UNITLESS*\}
  - number of poles
- **magnetic_wavelength**: *NX_FLOAT* \{units=*NX_WAVELENGTH*\}
- **k**: *NX_FLOAT* \{units=*NX_DIMENSIONLESS*\}
  - beam displacement parameter
- **length**: *NX_FLOAT* \{units=*NX_LENGTH*\}
  - length of insertion device
- **power**: *NX_FLOAT* \{units=*NX_POWER*\}
  - total power delivered by insertion device
- **energy**: *NX_FLOAT* \{units=*NX_ENERGY*\}
  - energy of peak intensity in output spectrum
- **bandwidth**: *NX_FLOAT* \{units=*NX_ENERGY*\}
  - bandwidth of peak energy
- **harmonic**: *NX_INT* \{units=*NX_UNITLESS*\}
  - harmonic number of peak
- **spectrum**: *NXdata*
  - spectrum of insertion device
- **(geometry)**: *NXgeometry*
  - “Engineering” position of insertion device

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXinsertion_device.nxdl.xml

### 3.3.25 NXinstrument

Status:

- base class, extends *NXobject*, version 1.0

Description:
Template of instrument descriptions comprising various beamline components. Each component will also be a NeXus group defined by its distance from the sample. Negative distances represent beamline components that are before the sample while positive distances represent components that are after the sample. This device allows the unique identification of beamline components in a way that is valid for both reactor and pulsed instrumentation.

Symbols:
No symbol table

Groups cited: NXaperture, NXattenuator, NXbeam_stop, NXbeam, NXbending_magnet, NXcapillary, NXcollection, NXcollimator, NXcrystal, NXdetector_group, NXdetector, NXdisk_chopper, NXevent_data, NFermi_chopper, NXfilter, NXflipper, NXguide, NXinsertion_device, NXmirror, N Xm moderator, NXmonochromator, NXpolarizer, NXpositioner, NXsource, NXvelocity_selector, NXraylens

Structure:

name: NX_CHAR
  Name of instrument
  @short_name: NX_CHAR
    short name for instrument, perhaps the acronym

(aperture): NXaperture
(attenuator): NXattenuator
(beam): NXbeam
(beam_stop): NXbeam_stop
(bending_magnet): NXbending_magnet
(collimator): NXcollimator
(collection): NXcollection
(capillary): NXcapillary
(crystal): NXcrystal
(detector): NXdetector
(detector_group): NXdetector_group
(disk_chopper): NXdisk_chopper
(event_data): NXevent_data
(fermi_chopper): NFermi_chopper
(filter): NXfilter
(flipper): NXflipper
(guide): NXguide
(insertion_device): NXinsertion_device
(mirror): NXmirror
(moderator): N Xm moderator
(monochromator): NXmonochromator
(polarizer): NXpolarizer
3.3.26 **NXlog**

**Status:**

base class, extends *NXobject*, version 1.0

**Description:**

Definition of information that is recorded against time, such as information monitored during the run. It contains the logged values and the times at which they were measured as elapsed time since a starting time recorded in ISO8601 format. This method of storing logged data helps to distinguish instances in which a variable is a dimension scale of the data, in which case it is stored in an *NXdata* group, and instances in which it is logged during the run, when it should be stored in an *NXlog* group. Note: When using multiple *NXlog* groups, it is suggested to place them inside a *NXcollection* group. In such cases, when *NXlog* is used in another class, *NXcollection/NXlog* is then constructed.

**Symbols:**

No symbol table

**Groups cited:** none

**Structure:**

- **time**: *NX_FLOAT* \{units=NX\_TIME\}
  - Time of logged entry. The times are relative to the “start” attribute and in the units specified in the “units” attribute.
  - **@start**: *NX\_DATE\_TIME*

- **value**: *NX\_NUMBER* \{units=NX\_ANY\}
  - Array of logged value, such as temperature

- **raw\_value**: *NX\_NUMBER* \{units=NX\_ANY\}
  - Array of raw information, such as thermocouple voltage

- **description**: *NX\_CHAR*
  - Description of logged value

- **average\_value**: *NX\_FLOAT* \{units=NX\_ANY\}

- **average\_value\_error**: *NX\_FLOAT* \{units=NX\_ANY\}
  - estimated uncertainty (often used: standard deviation) of average\_value

- **minimum\_value**: *NX\_FLOAT* \{units=NX\_ANY\}

- **maximum\_value**: *NX\_FLOAT* \{units=NX\_ANY\}

- **duration**: *NX\_FLOAT* \{units=NX\_ANY\}
  - Total time log was taken
3.3.27 NXmirror

Status:
  base class, extends NXobject, version 1.0

Description:
  Template of a beamline mirror or supermirror.

Symbols:
  No symbol table

Groups cited: NXdata, NXgeometry, NXshape

Structure:

  type: NX_CHAR
    Any of these values:
      • single: mirror with a single material as a reflecting surface
      • multi: mirror with stacked, multiple layers as a reflecting surface

  description: NX_CHAR
    description of this mirror

  incident_angle: NX_FLOAT {units=NX_ANGLE}

  bend_angle_x: NX_FLOAT {units=NX_ANGLE}

  bend_angle_y: NX_FLOAT {units=NX_ANGLE}

  interior_atmosphere: NX_CHAR
    Any of these values: vacuum|helium|argon

  external_material: NX_CHAR
    external material outside substrate

  m_value: NX_FLOAT {units=NX_UNITLESS}
    The m value for a supermirror, which defines the supermirror regime in multiples of the critical angle of Nickel.

  substrate_material: NX_CHAR

  substrate_density: NX_FLOAT {units=NX_MASS_DENSITY}

  substrate_thickness: NX_FLOAT {units=NX_LENGTH}

  coating_material: NX_CHAR

  substrate_roughness: NX_FLOAT {units=NX_LENGTH}

  coating_roughness: NX_FLOAT {units=NX_LENGTH}

  even_layer_material: NX_CHAR

  even_layer_density: NX_FLOAT {units=NX_MASS_DENSITY}

  odd_layer_material: NX_CHAR
**odd_layer_density:** `NX_FLOAT` [units=NX_MASS_DENSITY]

**layer_thickness:** `NX_FLOAT` [units=NX_LENGTH]

An array describing the thickness of each layer

(geography): `NXgeometry`

**reflectivity:** `NXdata`

Reflectivity as function of wavelength

**shape:** `NXshape`

A NXshape group describing the shape of the mirror

**figure_data:** `NXdata`

Numerical description of the surface figure of the mirror.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXmirror.nxdl.xml

### 3.3.28 NXmoderator

**Status:**

base class, extends `NXobject`, version 1.0

**Description:**

This is the description for a general moderator

**Symbols:**

No symbol table

**Groups cited:** `NXdata, NXgeometry, NXlog`

**Structure:**

**distance:** `NX_FLOAT` [units=NX_LENGTH]

Effective distance as seen by measuring radiation

**type:** `NX_CHAR`

Any of these values:

- H2O
- D2O
- Liquid H2
- Liquid CH4
- Liquid D2
- Solid D2
- C
- Solid CH4
- Solid H2

**poison_depth:** `NX_FLOAT` [units=NX_LENGTH]

**coupled:** `NX_BOOLEAN`
whether the moderator is coupled

coupling_material:  **NX_CHAR**

The material used for coupling. Usually Cd.

poison_material:  **NX_CHAR**

Any of these values: Gd | Cd

temperature:  **NX_FLOAT**  \{units=NX_TEMPERATURE\}

average/nominal moderator temperature

(geomeiry):  **NXgeometry**

“Engineering” position of moderator

temperature_log:  **NXlog**

log file of moderator temperature

pulse_shape:  **NXdata**

moderator pulse shape

Source:  Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXmoderator.nxdl.xml

### 3.3.29 NXmonitor

**Status:**

base class, extends  **NXobject**, version 1.0

**Description:**

Template of monitor data. It is similar to the NXdata groups containing monitor data and its associated dimension scale, e.g. time_of_flight or wavelength in pulsed neutron instruments. However, it may also include integrals, or scalar monitor counts, which are often used in both in both pulsed and steady-state instrumentation.

**Symbols:**

No symbol table

**Groups cited:**  **NXgeometry, NXlog**

**Structure:**

mode:  **NX_CHAR**

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: monitor|timer

start_time:  **NX_DATE_TIME**

Starting time of measurement

distance:  **NX_FLOAT**  \{units=NX_LENGTH\}

preset:  **NX_NUMBER**  \{units=NX_ANY\}

presest value for time or monitor
Distance of monitor from sample

**range[2]:** `NX_FLOAT {units=NX_ANY}`

Range (X-axis, Time-of-flight, etc.) over which the integral was calculated

**integral:** `NX_NUMBER {units=NX_ANY}`

Total integral monitor counts

**type:** `NX_CHAR`

Any of these values: Fission Chamber|Scintillator

**time_of_flight[ref(efficiency)]:** `NX_FLOAT {units=NX_TIME_OF_FLIGHT}`

Time-of-flight

**efficiency[ref(i)]:** `NX_NUMBER {units=NX_DIMENSIONLESS}`

Monitor efficiency

**data[n]:** `NX_NUMBER {units=NX_ANY}`

Monitor data

The signal and axes attributes take the same definitions as in `NXdata`:

**signal** signal=1 means this is the plottable data

**axes** axes="names" where names are defined as a colon-delimited string within this attribute in the C-order of the data array

**@signal:** `NX_POSINT`

as defined for NXdata

**@axes:** `NX_CHAR`

as defined for NXdata

**sampled_fraction:** `NX_FLOAT {units=NX_DIMENSIONLESS}`

Proportion of incident beam sampled by the monitor (0<\(x\)<1)

**count_time:** `NX_FLOAT {units=NX_TIME}`

Elapsed actual counting time, can be an array of size \(n_p\) when scanning. This is not the difference of the calendar time but the time the instrument was really counting, without pauses or times lost due beam unavailability

**integral_log:** `NXlog`

Time variation of monitor counts

**(geometry):** `NXgeometry`

Geometry of the monitor

**Source:** Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXmonitor.nxdl.xml

### 3.3.30 NXmonochromator

**Status:**

base class, extends `NXobject`, version 1.0

**Description:**
This is a base class for everything which selects a wavelength or energy, be it a monochromator crystal, a velocity selector, an undulator or whatever.

The expected units are:

- wavelength: angstrom
- energy: eV

Symbols:

No symbol table

Groups cited: NXcrystal, NXdata, NXgeometry, NXvelocity_selector

Structure:

- **wavelength**: `NX_FLOAT {units=NX_WAVELENGTH}`
  - wavelength selected
- **wavelength_error**: `NX_FLOAT {units=NX_WAVELENGTH}`
  - wavelength standard deviation
- **energy**: `NX_FLOAT {units=NX_ENERGY}`
  - energy selected
- **energy_error**: `NX_FLOAT {units=NX_ENERGY}`
  - energy standard deviation
- **distribution**: `NXdata`
- **geometry**: `NXgeometry`
- **(crystal)**: `NXcrystal`
  - Use as many crystals as necessary to describe
- **(velocity selector)**: `NXvelocity_selector`

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXmonochromator.nxdl.xml

### 3.3.31 NXnote

**Status:**

- base class, extends `NXobject`, version 1.0

**Description:**

- This class can be used to store additional information in a NeXus file e.g. pictures, movies, audio, additional text logs

Symbols:

- No symbol table

Groups cited: none

Structure:

- **author**: `NX_CHAR`
  - Author or creator of note
- **date**: `NX_DATE_TIME`
Date note created/added
type: \textit{NX\_CHAR}
Mime content type of note data field e.g. image/jpeg, text/plain, text/html

\textbf{Mime content type of note data field e.g. image/jpeg, text/plain, text/html}

file\_name: \textit{NX\_CHAR}
Name of original file name if note was read from an external source
description: \textit{NX\_CHAR}
Title of an image or other details of the note
data: \textit{NX\_BINARY}
Binary note data - if text, line terminator is [CR][LF].

Source: Automatically generated from \url{https://github.com/nexusformat/definitions/blob/master/base_classes/NXnote.nxdl.xml}

\subsection{3.3.32 NXobject}

\textbf{Status:}
base class, extends none, version 1.0

\textbf{Description:}
This is the base object of NeXus

\textbf{Symbols:}
No symbol table

\textbf{Groups cited: } none

\textbf{Structure:}
Source: Automatically generated from \url{https://github.com/nexusformat/definitions/blob/master/base_classes/NXobject.nxdl.xml}

\subsection{3.3.33 NXorientation}

\textbf{Status:}
base class, extends \textit{NXobject}, version 1.0

\textbf{Description:}
This is the description for a general orientation of a component - it is used by the NXgeometry class

\textbf{Symbols:}
No symbol table

\textbf{Groups cited: } \textit{NXgeometry}

\textbf{Structure:}
\begin{verbatim}
  value[numobj, 6]: \textit{NX\_FLOAT} [units=\textit{NX\_UNITLESS}]
\end{verbatim}
The orientation information is stored as direction cosines. The direction cosines will be between the local coordinate directions and the reference directions (to origin or relative NXgeometry). Calling the local unit vectors (x’,y’,z’) and the reference unit vectors (x,y,z) the six numbers will be [x’ dot x, x’ dot y, x’ dot z, y’ dot x, y’ dot y, y’ dot z] where “dot” is the
scalar dot product (cosine of the angle between the unit vectors). The unit vectors in both the local and reference coordinates are right-handed and orthonormal.

The pair of groups NXtranslation and NXorientation together describe the position of a component.

(geometry): **NXgeometry**

Link to another object if we are using relative positioning, else absent

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXorientation.nxdl.xml

### 3.3.34 NXparameters

**Status:**

base class, extends **NXobject**, version 1.0

**Description:**

Container for parameters, usually used in processing or analysis.

**Symbols:**

No symbol table

**Groups cited:** none

**Structure:**

**term**: **NX_CHAR**

A parameter (also known as a term) that is used in or results from processing.

@units: **NX_CHAR**

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXparameters.nxdl.xml

### 3.3.35 NXpinhole

**Status:**

base class, extends **NXobject**, version 1.0

**Description:**

Template of a simple pinhole. For more complex geometries NXaperture should be used.

**Symbols:**

No symbol table

**Groups cited:** none

**Structure:**

**depends_on**: **NX_CHAR**

Points to the path of the last element in the geometry chain that places this object in space. When followed through that chain is supposed to end at an element depending on "." i.e. the origin of the coordinate system. If desired the location of the slit can also be described relative to an NXbeam, which will allow a simple description of a non-centred pinhole.

**diameter**: **NX_NUMBER** {units=**NX_LENGTH**}
Size of the circular hole defining the transmitted beam size.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXpinhole.nxdl.xml

### 3.3.36 NXpolarizer

**Status:**
base class, extends *NXobject*, version 1.0

**Description:**
Template of a beamline spin polarizer. This is a draft and is subject to revision.

**Symbols:**
No symbol table

**Groups cited:** none

**Structure:**
- **type**: *NX_CHAR*
  one of these values: “crystal”, “supermirror”, “3He”
- **composition**: *NX_CHAR*
  description of the composition of the polarizing material
- **reflection[3]**: *NX_INT* \{units=NX_UNITSLESS\}
  \[hkl\] values of nominal reflection
- **efficiency**: *NX_FLOAT* \{units=NX_DIMENSIONLESS\}
  polarizing efficiency

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXpolarizer.nxdl.xml

### 3.3.37 NXpositioner

**Status:**
base class, extends *NXobject*, version 1.0

**Description:**
This group describes a generic positioner such as a motor or piezo-electric transducer. It is used to document the current information of a piece of beam line equipment. Note: When using multiple NXpositioner groups, it is suggested to place them inside a NXcollection group. In such cases, when NXpositioner is used in another class, NXcollection/NXpositioner is then constructed.

**Symbols:**
No symbol table

**Groups cited:** none

**Structure:**
- **name**: *NX_CHAR*
  symbolic or mnemonic name (one word)
description: \textit{NX\_CHAR}

description of positioner

description

value\[n\]: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

best known value of positioner - need \[n\] as may be scanned

raw\_value\[n\]: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

raw value of positioner - need \[n\] as may be scanned

target\_value\[n\]: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

targeted (commanded) value of positioner - need \[n\] as may be scanned

tolerance\[n\]: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

maximum allowable difference between target\_value and value

soft\_limit\_min: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

minimum allowed limit to set value

soft\_limit\_max: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

maximum allowed limit to set value

velocity: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

velocity of the positioner (distance moved per unit time)

acceleration\_time: \textit{NX\_NUMBER} \{units=\textit{NX\_ANY}\}

time to ramp the velocity up to full speed

controller\_record: \textit{NX\_CHAR}

Hardware device record, e.g. EPICS process variable, taco/tango ...

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXpositioner.nxdl.xml

3.3.38 **NXprocess**

Status:

base class, extends \textit{NXobject}, version 1.0

Description:

Document an event of data processing, reconstruction, or analysis for this data.

Symbols:

No symbol table

Groups cited: \textit{NXnote}

Structure:

\textbf{program}: \textit{NX\_CHAR}

Name of the program used

\textbf{version}: \textit{NX\_CHAR}

Version of the program used

\textbf{date}: \textit{NX\_DATE\_TIME}
Date and time of processing.

(note): **NXnote**

The note will contain information about how the data was processed or anything about the data provenance. The contents of the note can be anything that the processing code can understand, or simple text.

The name will be numbered to allow for ordering of steps.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXprocess.nxdl.xml

### 3.3.39 NXroot

**Status:**

base class, extends **NXobject**, version 1.0

**Description:**

Definition of the root NeXus group.

**Symbols:**

No symbol table

**Groups cited:** **NXentry**

**Structure:**

@**NX_class**: **NX_CHAR**

The root of any NeXus data file is an NXroot class (no other choice is allowed for a valid NeXus data file). This attribute cements that definition.

Obligatory value: **NXroot**

@**file_time**: **NX_CHAR**

Date and time file was originally created

@**file_name**: **NX_CHAR**

File name of original NeXus file

@**file_update_time**: **NX_CHAR**

Date and time of last file change at close

@**NeXus_version**: **NX_CHAR**

Version of NeXus API used in writing the file

@**HDF_version**: **NX_CHAR**

Version of NeXus API used in writing the file

@**HDF5_Version**: **NX_CHAR**

Version of NeXus API used in writing the file. Note this attribute is spelled with uppercase “V”, different than other version attributes.

@**XML_version**: **NX_CHAR**

Version of NeXus API used in writing the file

@**creator**: **NX_CHAR**
facility or program where file originated

(entry): NXentry

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXroot.nxdl.xml

### 3.3.40 NXsample

**Status:**

base class, extends NXobject, version 1.0

**Description:**

Template of the state of the sample. This could include scanned variables that are associated with one of the data dimensions, e.g. the magnetic field, or logged data, e.g. monitored temperature vs elapsed time.

**Symbols:**

symbolic array lengths to be coordinated between various fields

- **n_comp**: number of compositions
- **n_Temp**: number of temperatures
- **n_eField**: number of values in applied electric field
- **n_mField**: number of values in applied magnetic field
- **n_pField**: number of values in applied pressure field
- **n_sField**: number of values in applied stress field

**Groups cited:** NXbeam, NXdata, NXenvironment, NXgeometry, NXlog, NXpositioner

**Structure:**

- **name**: NX_CHAR
  
  Descriptive name of sample

- **chemical_formula**: NX_CHAR

  The chemical formula specified using CIF conventions. Abbreviated version of CIF standard:

  - Only recognized element symbols may be used.
  - Each element symbol is followed by a ‘count’ number. A count of ‘1’ may be omitted.
  - A space or parenthesis must separate each cluster of (element symbol + count).
  - Where a group of elements is enclosed in parentheses, the multiplier for the group must follow the closing parentheses. That is, all element and group multipliers are assumed to be printed as subscripted numbers.
  - Unless the elements are ordered in a manner that corresponds to their chemical structure, the order of the elements within any group or moiety depends on whether or not carbon is present.
  - If carbon is present, the order should be:
    - C, then H, then the other elements in alphabetical order of their symbol.
    - If carbon is not present, the elements are listed purely in alphabetic order of their symbol.
• This is the Hill system used by Chemical Abstracts.

**temperature**[n\_Temp]: *NX\_FLOAT* \{units=NX\_TEMPERATURE\}

Sample temperature. This could be a scanned variable

**electric\_field**[n\_eField]: *NX\_FLOAT* \{units=NX\_VOLTAGE\}

Applied electric field

@**direction**: *NX\_CHAR*

Any of these values: \( x | y | z \)

**magnetic\_field**[n\_mField]: *NX\_FLOAT* \{units=NX\_ANY\}

Applied magnetic field

@**direction**: *NX\_CHAR*

Any of these values: \( x | y | z \)

**stress\_field**[n\_sField]: *NX\_FLOAT* \{units=NX\_ANY\}

Applied external stress field

@**direction**: *NX\_CHAR*

Any of these values: \( x | y | z \)

**pressure**[n\_pField]: *NX\_FLOAT* \{units=NX\_PRESSURE\}

Applied pressure

**changer\_position**: *NX\_INT* \{units=NX\_UNITLESS\}

Sample changer position

**unit\_cell**[n\_comp, 6]: *NX\_FLOAT* \{units=NX\_LENGTH\}

Unit cell parameters (lengths and angles)

**unit\_cell\_volume**[n\_comp]: *NX\_FLOAT* \{units=NX\_VOLUME\}

Volume of the unit cell

**sample\_orientation**[3]: *NX\_FLOAT* \{units=NX\_ANGLE\}

This will follow the Busing and Levy convention from Acta.Crysta v22, p457 (1967)

**orientation\_matrix**[n\_comp, 3, 3]: *NX\_FLOAT*

Orientation matrix of single crystal sample. This will follow the Busing and Levy convention from Acta.Crysta v22, p457 (1967)

**mass**[n\_comp]: *NX\_FLOAT* \{units=NX\_MASS\}

Mass of sample

**density**[n\_comp]: *NX\_FLOAT* \{units=NX\_MASS\_DENSITY\}

Density of sample

**relative\_molecular\_mass**[n\_comp]: *NX\_FLOAT* \{units=NX\_MASS\}

Relative Molecular Mass of sample

**type**: *NX\_CHAR*

Any of these values:
• sample
• sample+can
• can
• calibration sample
• normalisation sample
• simulated data
• none
• sample environment

situation: NX_CHAR

The atmosphere will be one of the components, which is where its details will be stored; the relevant components will be indicated by the entry in the sample_component member.

Any of these values:
• air
• vacuum
• inert atmosphere
• oxidising atmosphere
• reducing atmosphere
• sealed can
• other

description: NX_CHAR

Description of the sample

preparation_date: NX_DATE_TIME

Date of preparation of the sample

component[n_comp]: NX_CHAR

Details of the component of the sample and/or can

sample_component[n_comp]: NX_CHAR

Type of component

Any of these values: sample|can|atmosphere|kit

concentration[n_comp]: NX_FLOAT \{units=\text{NX\_MASS\_DENSITY}\}

Concentration of each component

volume_fraction[n_comp]: NX_FLOAT

Volume fraction of each component

scattering_length_density[n_comp]: NX_FLOAT \{units=\text{NX\_SCATTERING\_LENGTH\_DENSITY}\}

Scattering length density of each component

unit_cell_class[n_comp]: NX_CHAR
In case it is all we know and we want to record/document it

Any of these values:

- cubic
- tetragonal
- orthorhombic
- monoclinic
- triclinic

unit_cell_group[n_comp]: \textit{NX\_CHAR}

Crystallographic point or space group

\textbf{path-length}: \textit{NX\_FLOAT} \{units=\textit{NX\_LENGTH}\}

Path length through sample/can for simple case when it does not vary with scattering direction

\textbf{path-length-window}: \textit{NX\_FLOAT} \{units=\textit{NX\_LENGTH}\}

Thickness of a beam entry/exit window on the can (mm) - assumed same for entry and exit

\textbf{thickness}: \textit{NX\_FLOAT} \{units=\textit{NX\_LENGTH}\}

Sample thickness

\textbf{external\_DAC}: \textit{NX\_FLOAT} \{units=\textit{NX\_ANY}\}

Value sent to user’s sample setup

\textbf{short-title}: \textit{NX\_CHAR}

20 character fixed length sample description for legends

\textbf{rotation-angle}: \textit{NX\_FLOAT} \{units=\textit{NX\_ANGLE}\}

Optional rotation angle for the case when the powder diagram has been obtained through an omega-2theta scan like from a traditional single detector powder diffractometer

\textbf{x-translation}: \textit{NX\_FLOAT} \{units=\textit{NX\_LENGTH}\}

Translation of the sample along the X-direction of the laboratory coordinate system

\textbf{distance}: \textit{NX\_FLOAT} \{units=\textit{NX\_LENGTH}\}

Translation of the sample along the Z-direction of the laboratory coordinate system

\textbf{geometry}: \textit{NXgeometry}

The position and orientation of the center of mass of the sample

\textbf{(beam)}: \textit{NXbeam}

Details of beam incident on sample - used to calculate sample/beam interaction point

\textbf{transmission}: \textit{NXdata}

As a function of Wavelength

\textbf{temperature-log}: \textit{NXlog}

temperature_log.value is a link to e.g. temperature_env.sensor1.value_log.value

\textbf{temperature-env}: \textit{NXenvironment}

Additional sample temperature environment information
magnetic_field_log: NXlog
magnetic_field_log.value is a link to e.g. magnetic_field_env.sensor1.value_log.value

magnetic_field_env: NXenvironment
Additional sample magnetic environment information

external_ADC: NXlog
logged value (or logic state) read from user’s setup

(positioner): NXpositioner
Any positioner (motor, PZT, ...) used to locate the sample

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXsample.nxdl.xml

3.3.41 NXsensor

Status:
base class, extends NXobject, version 1.0

Description:
This class describes a sensor used to monitor an external condition - the condition itself is described in NXenvironment

Symbols:
No symbol table

Groups cited: NXgeometry, NXlog, NXorientation

Structure:
model: NX_CHAR
Sensor identification code/model number

name: NX_CHAR
Name for the sensor

short_name: NX_CHAR
Short name of sensor used e.g. on monitor display program

attached_to: NX_CHAR
where sensor is attached to ("sample" | "can")

measurement: NX_CHAR
name for measured signal
Any of these values:
• temperature
• pH
• magnetic_field
• electric_field
• conductivity
• resistance
• voltage
• pressure
• flow
• stress
• strain
• shear
• surface_pressure

type: **NX_CHAR**

The type of hardware used for the measurement. Examples (suggestions but not restrictions):

- **Temperature**  J | K | T | E | R | S | Pt100 | Rh/Fe
- **pH**  Hg/Hg2Cl2 | Ag/AgCl | ISFET
- **Ion selective electrode** specify species; e.g. Ca2+
- **Magnetic field**  Hall
- **Surface pressure**  wilhelmy plate

**run_control**: **NX_BOOLEAN**

Is data collection controlled or synchronised to this quantity: 1=no, 0=to “value”, 1=to “value_deriv1”, etc.

**high_trip_value**: **NX_FLOAT** {units=**NX_ANY**}

Upper control bound of sensor reading if using run_control

**low_trip_value**: **NX_FLOAT** {units=**NX_ANY**}

Lower control bound of sensor reading if using run_control

**value[n]**: **NX_FLOAT** {units=**NX_ANY**}

nominal setpoint or average value - need [n] as may be a vector

**value_deriv1[ref(value)]**: **NX_FLOAT** {units=**NX_ANY**}

Nominal/average first derivative of value e.g. strain rate - same dimensions as “value” (may be a vector)

**value_deriv2[ref(value)]**: **NX_FLOAT** {units=**NX_ANY**}

Nominal/average second derivative of value - same dimensions as “value” (may be a vector)

**external_field_brief**: **NX_CHAR**

Any of these values:

• along beam
• across beam
• transverse
• solenoidal
• flow shear gradient
• flow vorticity
geometry: \textit{NXgeometry}

Defines the axes for logged vector quantities if they are not the global instrument axes

value\_log: \textit{NXlog}

Time history of sensor readings

value\_deriv1\_log: \textit{NXlog}

Time history of first derivative of sensor readings

value\_deriv2\_log: \textit{NXlog}

Time history of second derivative of sensor readings

external\_field\_full: \textit{NXorientation}

For complex external fields not satisfied by External\_field\_brief

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXsensor.nxdl.xml

3.3.42 \textbf{NXshape}

Status:

base class, extends \textit{NXobject}, version 1.0

Description:

This is the description of the general shape and size of a component, which may be made up of \textit{numobj} separate elements - it is used by the \textit{NXgeometry} class

Symbols:

No symbol table

Groups cited: none

Structure:

\textbf{shape}: \textit{NX\_CHAR}

general shape of a component

Any of these values:

- \texttt{nxflat}
- \texttt{nxcylinder}
- \texttt{nxbox}
- \texttt{nxsphere}
- \texttt{nxcone}
- \texttt{nxelliptical}
- \texttt{nxtoroidal}
- \texttt{nxparabolic}
- \texttt{nxpolynomial}

\textbf{size[\textit{numobj, nshapepar}]}: \textit{NX\_FLOAT} \{units=\textit{NX\_LENGTH}\}
physical extent of the object along its local axes (after NXorientation) with the center of mass at the local origin (after NXtranslation). The meaning and location of these axes will vary according to the value of the “shape” variable. nshapepar defines how many parameters:

- For “nxcylinder” type the parameters are (diameter, height) and a three value orientation vector of the cylinder.
- For the “nxbox” type the parameters are (length, width, height).
- For the “nxsphere” type the parameters are (diameter).
- For nxcone cone half aperture
- For nxelliptical, semi-major axis, semi-minor-axis, angle of major axis and pole
- For nxtoroidal, major radius, minor radius
- For nxparabolic, parabolic parameter a
- For nxpolynomial, an array of polynom coefficients, the dimension of the array encodes the degree of the polynom

**direction**: NX_CHAR

Any of these values: concave | convex

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXshape.nxdl.xml

### 3.3.43 NXslit

**Status**: base class, extends NXobject, version 1.0

**Description**: Template of a simple slit. For more complex geometries NXaperture should be used.

**Symbols**: No symbol table

**Groups cited**: none

**Structure**:

- **depends_on**: NX_CHAR
  
  Points to the path of the last element in the geometry chain that places this object in space. When followed through that chain is supposed to end at an element depending on ”.” i.e. the origin of the coordinate system. If desired the location of the slit can also be described relative to an NXbeam, which will allow a simple description of a non-centred slit.

- **x_gap**: NX_NUMBER {units=NX_LENGTH}
  
  Size of the gap opening in the first dimension of the local coordinate system.

- **y_gap**: NX_NUMBER {units=NX_LENGTH}
  
  Size of the gap opening in the second dimension of the local coordinate system.

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXslit.nxdl.xml
### 3.3.44 NXsource

**Status:**

base class, extends `NXobject`, version 1.0

**Description:**

Template of the neutron or x-ray source, insertion devices and/or moderators.

**Symbols:**

No symbol table

**Groups cited:** `NXdata, NXgeometry, NXnote`

**Structure:**

- **distance:** `NX_FLOAT {units=NX_LENGTH}`
  
  Effective distance from sample. Distance as seen by radiation from sample. This number should be negative to signify that it is upstream of the sample.

- **name:** `NX_CHAR`
  
  Name of source

- **@short_name:** `NX_CHAR`
  
  Short name for source, perhaps the acronym

- **type:** `NX_CHAR`
  
  Type of radiation source (pick one from the enumerated list and spell exactly)

  Any of these values:

  - Spallation Neutron Source
  - Pulsed Reactor Neutron Source
  - Reactor Neutron Source
  - Synchrotron X-ray Source
  - Pulsed Muon Source
  - Rotating Anode X-ray
  - Fixed Tube X-ray
  - UV Laser
  - Free-Electron Laser
  - Optical Laser
  - Ion Source
  - UV Plasma Source

- **probe:** `NX_CHAR`
  
  Type of radiation probe (pick one from the enumerated list and spell exactly)

  Any of these values:

  - neutron
  - x-ray
• muon
• electron
• ultraviolet
• visible light
• positron
• proton

**power:** *NX_FLOAT* {units=*NX_POWER*}

Source power

**emittance_x:** *NX_FLOAT* {units=*NX_EMITTANCE*}

Source emittance (nm-rad) in X (horizontal) direction.

**emittance_y:** *NX_FLOAT* {units=*NX_EMITTANCE*}

Source emittance (nm-rad) in Y (horizontal) direction.

**sigma_x:** *NX_FLOAT* {units=*NX_LENGTH*}

Particle beam size in x

**sigma_y:** *NX_FLOAT* {units=*NX_LENGTH*}

Particle beam size in y

**flux:** *NX_FLOAT* {units=*NX_FLUX*}

Source intensity/area (example: s-1 cm-2)

**energy:** *NX_FLOAT* {units=*NX_ENERGY*}

Source energy. For storage rings, this would be the particle beam energy. For X-ray tubes, this would be the excitation voltage.

**current:** *NX_FLOAT* {units=*NX_CURRENT*}

Accelerator, X-ray tube, or storage ring current

**voltage:** *NX_FLOAT* {units=*NX_VOLTAGE*}

Accelerator voltage

**frequency:** *NX_FLOAT* {units=*NX_FREQUENCY*}

Frequency of pulsed source

**period:** *NX_FLOAT* {units=*NX_PERIOD*}

Period of pulsed source

**target_material:** *NX_CHAR*

Pulsed source target material

Any of these values:

• Ta
• W
• depleted_U
• enriched_U
number_of_bunches: **NX_INT**
For storage rings, the number of bunches in use.

bunch_length: **NX_FLOAT** \{units=**NX_TIME**\}
For storage rings, temporal length of the bunch

bunch_distance: **NX_FLOAT** \{units=**NX_TIME**\}
For storage rings, time between bunches

pulse_width: **NX_FLOAT** \{units=**NX_TIME**\}
temporal width of source pulse

mode: **NX_CHAR**
source operating mode

Any of these values:
- Single Bunch: for storage rings
- Multi Bunch: for storage rings

top_up: **NX_BOOLEAN**
Is the synchrotron operating in top_up mode?

last_fill: **NX_NUMBER** \{units=**NX_CURRENT**\}
For storage rings, the current at the end of the most recent injection.

@time: **NX_DATE_TIME**
date and time of the most recent injection.

notes: **NXnote**
any source/facility related messages/events that occurred during the experiment

bunch_pattern: **NXdata**
For storage rings, description of the bunch pattern. This is useful to describe irregular bunch patterns.

title: **NX_CHAR**
name of the bunch pattern

pulse_shape: **NXdata**
source pulse shape

geometry: **NXgeometry**
“Engineering” location of source

distribution: **NXdata**
The wavelength or energy distribution of the source

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXsource.nxdl.xml
### 3.3.45 NXsubentry

**Status:**

base class, extends *NXobject*, version 1.0

**Description:**

NXsubentry is a base class virtually identical to NXentry and is used as the (overlay) location for application definitions. Use a separate NXsubentry for each application definition.

To use NXsubentry with a hypothetical application definition called NXmyappdef:

- Create a group with attribute `NX_class="NXsubentry"
- Within that group, create a field called `definition="NXmyappdef"`
- There are two optional attributes of definition: `version` and `URL`

The intended use is to define application definitions for a multi-technique NXentry. Previously, an application definition replaced NXentry with its own definition. With the increasing popularity of instruments combining multiple techniques for data collection (such as SAXS/WAXS instruments), it was recognized the application definitions must be entered in the NeXus data file tree as children of NXentry.

**Symbols:**

No symbol table

**Groups cited:** NXcharacterization, NXcollection, NXdata, NXinstrument, NXmonitor, NXnote, NXparameters, NXprocess, NXsample, NXuser

**Structure:**

```python
@IDF_Version: NX_CHAR
  ISIS Muon IDF_Version
title: NX_CHAR
  Extended title for entry
experiment_identifier: NX_CHAR
  Unique identifier for the experiment, defined by the facility, possibly linked to the proposals
experiment_description: NX_CHAR
  Brief summary of the experiment, including key objectives.
collection_identifier: NX_CHAR
  User or Data Acquisition defined group of NeXus files or NXentry
collection_description: NX_CHAR
  Brief summary of the collection, including grouping criteria.
entry_identifier: NX_CHAR
  unique identifier for the measurement, defined by the facility.
definition: NX_CHAR
  Official NeXus NXDL schema to which this file conforms
@version: NX_CHAR
  NXDL version number
```
@URL: NX_CHAR
URL of NXDL file

definition_local: NX_CHAR
Local NXDL schema extended from the file specified in the definition field. This contains any locally-defined, additional fields in the file.

@version: NX_CHAR
NXDL version number

@URL: NX_CHAR
URL of NXDL file

start_time: NX_DATE_TIME
Starting time of measurement

duration: NX_INT (units=NX_TIME)
Duration of measurement

start_time: NX_DATE_TIME
Ending time of measurement

end_time: NX_DATE_TIME

run_cycle: NX_CHAR
Such as “2007-3”. Some user facilities organize their beam time into run cycles.

program_name: NX_CHAR
Name of program used to generate this file

@version: NX_CHAR
Program version number

@configuration: NX_CHAR
configuration of the program

revision: NX_CHAR
Revision id of the file due to re-calibration, reprocessing, new analysis, new instrument definition format, ...

@comment: NX_CHAR

pre_sample_flightpath: NX_FLOAT (units=NX_LENGTH)
This is the flightpath before the sample position. This can be determined by a chopper, by the moderator or the source itself. In other words: it the distance to the component which gives the T0 signal to the detector electronics. If another component in the NXinstrument hierarchy provides this information, this should be a link.

experiment_documentation: NXnote
Description of the full experiment (document in pdf, latex, ...)

notes: NXnote

3.3. Base Class Definitions
Notes describing entry

**thumbnail:** NXnote

A small image that is representative of the entry. An example of this is a 640x480 jpeg image automatically produced by a low resolution plot of the NXdata.

**@mime_type:** NX_CHAR

The value should be an image/*

Obligatory value: image/*

(characterization): NXcharacterization

(user): NXuser

(sample): NXsample

(instrument): NXinstrument

(collection): NXcollection

(monitor): NXmonitor

(data): NXdata

(parameters): NXparameters

(process): NXprocess

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXsubentry.nxdl.xml

### 3.3.46 NXtranslation

**Status:**

base class, extends NXobject, version 1.0

**Description:**

This is the description for the general spatial location of a component - it is used by the NXgeometry class

**Symbols:**

No symbol table

**Groups cited:** NXgeometry

**Structure:**

```
 distances[numobj, 3]: NX_FLOAT {units=NX_LENGTH}
```

(x,y,z) This field describes the lateral movement of a component. The pair of groups NXtranslation and NXorientation together describe the position of a component. For absolute position, the origin is the scattering center (where a perfectly aligned sample would be) with the z-axis pointing downstream and the y-axis pointing gravitationally up. For a relative position the NXtranslation is taken into account before the NXorientation. The axes are right-handed and orthonormal.

**geometry:** NXgeometry

Link to other object if we are relative, else absent

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXtranslation.nxdl.xml
3.3.47 **NXuser**

**Status:**
base class, extends *NXobject*, version 1.0

**Description:**
Template of user’s contact information. The format allows more than one user with the same affiliation and contact information, but a second NXuser group should be used if they have different affiliations, etc.

**Symbols:**
No symbol table

**Groups cited:** none

**Structure:**
- **name**: `NX_CHAR`
  Name of user responsible for this entry
- **role**: `NX_CHAR`
  Role of user responsible for this entry. Suggested roles are “local_contact”, “principal_investigator”, and “proposer”
- **affiliation**: `NX_CHAR`
  Affiliation of user
- **address**: `NX_CHAR`
  Address of user
- **telephone_number**: `NX_CHAR`
  Telephone number of user
- **fax_number**: `NX_CHAR`
  Fax number of user
- **email**: `NX_CHAR`
  Email of user
- **facility_user_id**: `NX_CHAR`
  facility based unique identifier for this person e.g. their identification code on the facility address/contact database

**Source:** Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXuser.nxdl.xml

3.3.48 **NXvelocity_selector**

**Status:**
base class, extends *NXobject*, version 1.0

**Description:**
This is the description for a (typically neutron) velocity selector

**Symbols:**
No symbol table

Groups cited: NXgeometry

Structure:

- **type**: `NX_CHAR`
  - velocity selector type

- **rotation_speed**: `NX_FLOAT {units=NX_FREQUENCY}`
  - velocity selector rotation speed

- **radius**: `NX_FLOAT {units=NX_LENGTH}`
  - radius at beam centre

- **spwidth**: `NX_FLOAT {units=NX_LENGTH}`
  - spoke width at beam centre

- **length**: `NX_FLOAT {units=NX_LENGTH}`
  - rotor length

- **num**: `NX_INT {units=NX_UNITLESS}`
  - number of spokes/lamella

- **twist**: `NX_FLOAT {units=NX_ANGLE}`
  - twist angle along axis

- **table**: `NX_FLOAT {units=NX_ANGLE}`
  - offset vertical angle

- **height**: `NX_FLOAT {units=NX_LENGTH}`
  - input beam height

- **width**: `NX_FLOAT {units=NX_LENGTH}`
  - input beam width

- **wavelength**: `NX_FLOAT {units=NX_WAVELENGTH}`
  - wavelength

- **wavelength_spread**: `NX_FLOAT {units=NX_WAVELENGTH}`
  - deviation FWHM /Wavelength

geometry: NXgeometry

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/base_classes/NXvelocity_selector.nxdl.xml

### 3.3.49 NXxraylens

**Status:**

- base class, extends NXobject, version 1.0

**Description:**

This is a dictionary of field names to use for describing a X-ray lens as used at synchrotron beam lines. Based on information provided by Gerd Wellenreuther.
Symbols:

No symbol table

Groups cited: NXnote

Structure:

lens_geometry: NX_CHAR
  Geometry of the lens
  Any of these values:
  • paraboloid
  • spherical
  • elliptical
  • hyperbolical

symmetric: NX_BOOLEAN
  Is the device symmetric?

cylindrical: NX_BOOLEAN
  Is the device cylindrical?

focus_type: NX_CHAR
  The type of focus of the lens
  Any of these values: line|point

lens_thickness: NX_FLOAT {units=NX_LENGTH}
  Thickness of the lens

lens_length: NX_FLOAT {units=NX_LENGTH}
  Length of the lens

curvature: NX_FLOAT {units=NX_LENGTH}
  Radius of the curvature as measured in the middle of the lens

aperture: NX_FLOAT {units=NX_LENGTH}
  Diameter or radius of the lens.

number_of_lenses: NX_INT
  Number of lenses that make up the compound lens.

lens_material: NX_CHAR
  Material used to make the lens.

gas: NX_CHAR
  Gas used to fill the lens

gas_pressure: NX_FLOAT {units=NX_PRESSURE}
  Gas pressure in the lens

cylinder_orientation: NXnote
  Orientation of the cylinder axis.

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3.4 Application Definitions

A description of each NeXus application definition is given. NeXus application definitions define the \textit{minimum} set of terms that \textit{must} be used in an instance of that class. Consider the application definitions as a \textit{contract} between a data provider (such as the beam line control system) and a data consumer (such as a data analysis program for a scientific technique) that describes the information is certain to be available in a data file.

3.4.1 NXarchive

Status:

application definition, extends \textit{NXobject}, version 1.0b

Description:

This is a definition for data to be archived by ICAT (http://www.icatproject.org/).

Symbols:

No symbol table

Groups cited: \textit{NXentry, NXinstrument, NXsample, NXsource, NXuser}

Structure:

\begin{verbatim}
entry: NXentry
  @index: NX_CHAR
  title: NX_CHAR
  experiment_identifier: NX_CHAR
    unique identifier for the experiment
  experiment_description: NX_CHAR
    Brief description of the experiment and its objectives
  collection_identifier: NX_CHAR
    ID of user or DAQ define group of data files
  collection_description: NX_CHAR
    Brief summary of the collection, including grouping criteria
  entry_identifier: NX_CHAR
    unique identifier for this measurement as provided by the facility
  start_time: NX_DATE_TIME
  end_time: NX_DATE_TIME
  duration: NX_FLOAT \{ units=NX_TIME \}
    TODO: needs documentation
  collection_time: NX_FLOAT \{ units=NX_TIME \}
    TODO: needs documentation
\end{verbatim}
run_cycle: NX_CHAR
   TODO: needs documentation

revision: NX_CHAR
   revision ID of this file, may be after recalibration, reprocessing etc.

definition: NX_CHAR
   Official NeXus NXDL schema to which this file conforms
   Obligatory value: NXarchive

program: NX_CHAR
   The program and version used for generating this file
   @version: NX_CHAR

release_date: NX_CHAR {units=NX_TIME}
   when this file is to be released into PD

user: NXuser
   name: NX_CHAR
   role: NX_CHAR
      role of the user
   facility_user_id: NX_CHAR
      ID of the user in the facility bureaucracy database

instrument: NXinstrument
   name: NX_CHAR
   description: NX_CHAR
      Brief description of the instrument
   (source): NXsource
      type: NX_CHAR
         Any of these values:
         • Spallation Neutron Source
         • Pulsed Reactor Neutron Source
         • Reactor Neutron Source
         • Synchrotron X-Ray Source
         • Pulsed Muon Source
         • Rotating Anode X-Ray
         • Fixed Tube X-Ray
   name: NX_CHAR
   probe: NX_CHAR
      Any of these values: neutron|x-ray|electron

sample: NXsample
name: NX_CHAR
Descriptive name of sample

sample_id: NX_CHAR
Unique database id of the sample

description: NX_CHAR

type: NX_CHAR
Any of these values:
• sample
• sample+can
• calibration sample
• normalisation sample
• simulated data
• none
• sample_environment

chemical_formula: NX_CHAR
Chemical formula formatted according to CIF conventions

preparation_date: NX_CHAR {units=NX_TIME}

situation: NX_CHAR
Description of the environment the sample is in: air, vacuum, oxidizing atmosphere, dehydrated, etc.

temperature: NX_FLOAT {units=NX_TEMPERATURE}
magnetic_field: NX_FLOAT {units=NX_CURRENT}
electric_field: NX_FLOAT {units=NX_VOLTAGE}
stress_field: NX_FLOAT {units=NX_UNITLESS}
pressure: NX_FLOAT {units=NX_PRESSURE}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXarchive.nxdl.xml

3.4.2 NXdirecttof

Status:
application definition, extends NXtofraw, version 1.0b

Description:
This is a application definition for raw data from a direct geometry TOF spectrometer

Symbols:
No symbol table

Groups cited: NXentry, NXfermi_chopper, NXinstrument

Structure:
entry: NXentry
    title: NX_CHAR
    start_time: NX_DATE_TIME
    definition: NX_CHAR
        Official NeXus NXDL schema to which this file conforms
        Obligatory value: NXdirecttof
    (instrument): NXinstrument
        fermi_chopper: NXfermi_chopper
            rotation_speed: NX_FLOAT {units=NX_FREQUENCY}
                chopper rotation speed
            energy: NX_FLOAT {units=NX_ENERGY}
                energy selected

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXdirecttof.nxdl.xml

3.4.3 NXfluo

Status:
    application definition, extends NXobject, version 1.0

Description:
    This is an application definition for raw data from an X-ray fluorescence experiment

Symbols:
    No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXmonochromator, NXsample, NXsource

Structure:

entry: NXentry
    title: NX_CHAR
    start_time: NX_DATE_TIME
    definition: NX_CHAR
        Official NeXus NXDL schema to which this file conforms.
        Obligatory value: NXfluo
    (instrument): NXinstrument
        (source): NXsource
            type: NX_CHAR
            name: NX_CHAR
            probe: NX_CHAR
                Obligatory value: x-ray
            monochromator: NXmonochromator
wavelength: \texttt{NX\_FLOAT}

fluorescence: \texttt{NXdetector}

data\[nenergy\]: \texttt{NX\_INT}

energy\[nenergy\]: \texttt{NX\_FLOAT}

(sam\(ple\)): \texttt{NXsample}

name: \texttt{NX\_CHAR}

Descriptive name of sample

(monitor): \texttt{NXmonitor}

mode: \texttt{NX\_CHAR}

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: \texttt{monitor} | \texttt{timer}

preset: \texttt{NX\_FLOAT}

preset value for time or monitor

data: \texttt{NX\_INT}

data: \texttt{NXdata}

e\textit{nergy} \rightarrow /\texttt{entry/instrument/fluorescence/energy}

data \rightarrow /\texttt{entry/instrument/fluorescence/data}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXfluo.nxdl.xml

3.4.4 \texttt{NXindirecttof}

Status:

application definition, extends \texttt{NXtofraw}, version 1.0b

Description:

This is an application definition for raw data from a direct geometry TOF spectrometer

Symbols:

No symbol table

Groups cited: \texttt{NXentry, NXinstrument, NXmonochromator}

Structure:

\texttt{entry: NXentry}

\hspace{1cm} title: \texttt{NX\_CHAR}

\hspace{1cm} start\_time: \texttt{NX\_DATE\_TIME}

\hspace{1cm} definition: \texttt{NX\_CHAR}

\hspace{1cm} Official NeXus NXDL schema to which this file conforms

\hspace{1cm} Obligatory value: \texttt{NXindirecttof}

\hspace{1cm} (instrument): \texttt{NXinstrument}
analyser: `NXmonochromator`

`energy[nDet]: NX_FLOAT {units=NX_ENERGY}`
analyzed energy

`polar_angle[nDet]: NX_FLOAT {units=NX_ANGLE}`
polar angle towards sample

`distance[nDet]: NX_FLOAT {units=NX_LENGTH}`
distance from sample

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXindirecttof.nxdl.xml

### 3.4.5 NXiqproc

Status:
application definition, extends `NXobject`, version 1.0b

Description:
Application definition for any $I(Q)$ data.

Symbols:
No symbol table

Groups cited: `NXdata`, `NXentry`, `NXinstrument`, `NXparameters`, `NXprocess`, `NXsample`, `NXsource`

Structure:

```
(entry): NXentry
  @entry: NX_CHAR
  title: NX_CHAR
  definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms
    Obligatory value: NXiqproc
  instrument: NXinstrument
    name: NX_CHAR
      Name of the instrument from which this data was reduced.
  (source): NXsource
    type: NX_CHAR
    name: NX_CHAR
    probe: NX_CHAR
      Any of these values: neutron\textbar{}x-ray\textbar{}electron
  (sample): NXsample
    name: NX_CHAR
      Descriptive name of sample
```

(reduction: NXprocess)
program: NX_CHAR
version: NX_CHAR
input: NXparameters
   Input parameters for the reduction program used
filenames: NX_CHAR
   Raw data files used to generate this I(Q)
output: NXparameters
   Eventual output parameters from the data reduction program used
(data): NXdata
data[NE, NQX, NQY]: NX_INT
   This is I(Q). The client has to analyse the dimensions of I(Q). Often, multiple I(Q) for various environment conditions are measured; that would be the first dimension. Q can be multidimensional, this accounts for the further dimensions in the data
variable[NE]: NX_CHAR
   @varied_variable: NX_CHAR
   The real name of the varied variable in the first dim of data, temperature, P, MF etc...
qx[NQX]: NX_CHAR
   Values for the first dimension of Q
qy[NQY]: NX_CHAR
   Values for the second dimension of Q

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXiqproc.nxdl.xml

3.4.6 NXlauetof

Status:
   application definition, extends NXobject, version 1.0b
Description:
   This is the application definition for a TOF laue diffractometer
Symbols:
   No symbol table
Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample
Structure:
   entry: NXentry
definition: NX_CHAR
   Official NeXus NXDL schema to which this file conforms
   Obligatory value: NXlauetof
instrument: NXinstrument

detector: NXdetector

This assumes a planar 2D detector. All angles and distances refer to the center of the detector.

**polar_angle**: NX_FLOAT \{units=NX_ANGLE\}

The polar_angle (two theta) where the detector is placed.

**azimuthal_angle**: NX_FLOAT \{units=NX_ANGLE\}

The azimuthal angle where the detector is placed.

data[number of x pixels, number of y pixels, nTOF]: NX_INT

@signal: NX_POSINT

Obligatory value: 1

**x_pixel_size**: NX_FLOAT \{units=NX_LENGTH\}

**y_pixel_size**: NX_FLOAT \{units=NX_LENGTH\}

distance: NX_FLOAT \{units=NX_LENGTH\}

time_of_flight[nTOF]: NX_FLOAT \{units=NX_TIME_OF_FLIGHT\}

sample: NXsample

**name**: NX_CHAR

Descriptive name of sample

**orientation_matrix[3, 3]**: NX_FLOAT

The orientation matrix according to Busing and Levy conventions. This is not strictly necessary as the UB can always be derived from the data. But let us bow to common usage which includes the UB nearly always.

**unit_cell[6]**: NX_FLOAT

The unit cell, a, b, c, alpha, beta, gamma. Again, not strictly necessary, but normally written.

control: NXmonitor

**mode**: NX_CHAR

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: monitor | timer

**preset**: NX_FLOAT

preset value for time or monitor

data[nTOF]: NX_INT

use these attributes primary=1 signal=1

time_of_flight[nTOF]: NX_FLOAT \{units=NX_TIME_OF_FLIGHT\}

name: NXdata

data –> /NXentry/NXinstrument/NXdetector/data

time_of_flight –> /NXentry/NXinstrument/NXdetector/time_of_flight
3.4.7 NXmonopd

Status:
application definition, extendsNXobject, version 1.0b

Description:
Monochromatic Neutron and X-Ray Powder Diffraction. Instrument definition for a powder diffractometer at a monochromatic neutron or X-ray beam. This is both suited for a powder diffractometer with a single detector or a powder diffractometer with a position sensitive detector.

Symbols:
No symbol table

Groups cited: NXcrystal, NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample, NXsource

Structure:
entry: NXentry
  title: NX_CHAR
  start_time: NX_DATE_TIME
  definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms
    Obligatory value: NXmonopd
  (instrument): NXinstrument
    (source): NXsource
      type: NX_CHAR
      name: NX_CHAR
      probe: NX_CHAR
        Any of these values: neutron|x-ray|electron
  (crystal): NXcrystal
    wavelength[i]: NX_FLOAT {units=NX_WAVELENGTH}
      Optimum diffracted wavelength
  (detector): NXdetector
    polar_angle[ndet]: NX_FLOAT
      where ndet = number of detectors
    data[ndet]: NX_INT
      detector signal (usually counts) are already corrected for detector efficiency
  (sample): NXsample
    name: NX_CHAR
    Descriptive name of sample
rotation_angle: \texttt{NX\_FLOAT} \{units=\texttt{NX\_ANGLE}\}

Optional rotation angle for the case when the powder diagram has been obtained through an omega-2theta scan like from a traditional single detector powder diffractometer

(monitor): \texttt{NX\_monitor}

mode: \texttt{NX\_CHAR}

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: monitor | timer

preset: \texttt{NX\_FLOAT}

preset value for time or monitor

integral: \texttt{NX\_FLOAT} \{units=\texttt{NX\_ANY}\}

Total integral monitor counts

(data): \texttt{NX\_data}

polar_angle \rightarrow /NXentry/NXinstrument/NXdetector/polar_angle

Link to polar angle in /NXentry/NXinstrument/NXdetector

data \rightarrow /NXentry/NXinstrument/NXdetector/data

Link to data in /NXentry/NXinstrument/NXdetector

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXmonopd.nxdl.xml

3.4.8 NXrefscan

Status:

application definition, extends \texttt{NX\_object}, version 1.0b

Description:

This is an application definition for a monochromatic scanning reflectometer. It does not have the information to calculate the resolution since it does not have any apertures.

Symbols:

No symbol table

Groups cited: \texttt{NX\_data}, \texttt{NX\_detector}, \texttt{NX\_entry}, \texttt{NX\_instrument}, \texttt{NX\_monitor}, \texttt{NX\_monochromator}, \texttt{NX\_sample}, \texttt{NX\_source}

Structure:

entry: \texttt{NX\_entry}

title: \texttt{NX\_CHAR}

start_time: \texttt{NX\_DATE\_TIME}

data_time: \texttt{NX\_DATE\_TIME}

definition: \texttt{NX\_CHAR}

Official NeXus NDXL schema to which this file conforms

Obligatory value: \texttt{NXrefscan}
instrument: NXinstrument
  (source): NXsource
    type: NX_CHAR
    name: NX_CHAR
    probe: NX_CHAR
      Any of these values: neutron|x-ray|electron
monochromator: NXmonochromator
  wavelength: NX_FLOAT {units=NX_WAVELENGTH}
(detector): NXdetector
  data[NP]: NX_INT
  polar_angle[NP]: NX_FLOAT {units=NX_ANGLE}
sample: NXsample
  name: NX_CHAR
    Descriptive name of sample
  rotation_angle[NP]: NX_FLOAT {units=NX_ANGLE}
control: NXmonitor
  mode: NX_CHAR
    Count to a preset value based on either clock time (timer) or received monitor counts (monitor).
    Any of these values: monitor|timer
  preset: NX_FLOAT
    preset value for time or monitor
  data[NP]: NX_FLOAT {units=NX_ANY}
    Monitor counts for each step
data: NXdata
  data -> /NXentry/NXinstrument/NXdetector/data
  rotation_angle -> /NXentry/NXsample/rotation_angle
  polar_angle -> /NXentry/NXinstrument/NXdetector/polar_angle

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXrefscan.nxdl.xml

3.4.9 NXreftof

Status:
  application definition, extends NXobject, version 1.0b

Description:
  This is an application definition for raw data from a TOF reflectometer.

Symbols:
No symbol table

Groups cited: NXdata, NXdetector, NXdisk_chopper, NXentry, NXinstrument, NXmonitor, NXsample

Structure:

entry: NXentry
  title: NX_CHAR
  start_time: NX_DATE_TIME
  end_time: NX_DATE_TIME
  definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms
    Obligatory value: NXreftof

instrument: NXinstrument
  name: NX_CHAR
  chopper: NXdisk_chopper
    distance: NX_FLOAT {units=NX_LENGTH}
    Distance between chopper and sample

detector: NXdetector
  data[xsize, ysize, nTOF]: NX_INT
  time_of_flight[nTOF]: NX_FLOAT {units=NX_TIME_OF_FLIGHT}
    Array of time values for each bin in a time-of-flight measurement
  distance: NX_FLOAT {units=NX_LENGTH}
  polar_angle: NX_FLOAT {units=NX_ANGLE}
  x_pixel_size: NX_FLOAT {units=NX_LENGTH}
  y_pixel_size: NX_FLOAT {units=NX_LENGTH}

sample: NXsample
  name: NX_CHAR
  Descriptive name of sample
  rotation_angle: NX_FLOAT {units=NX_ANGLE}

control: NXmonitor
  mode: NX_CHAR
    Count to a preset value based on either clock time (timer) or received monitor counts (monitor).
    Any of these values: monitor | timer
  preset: NX_FLOAT {units=NX_ANY}
    preset value for time or monitor
  integral: NX_INT
    Total integral monitor counts
**time_of_flight**: `NX_FLOAT` [units=NX_TIME_OF_FLIGHT]

Time channels

data: `NX_INT`

Monitor counts in each time channel

data: `NXdata`

data -> /NXentry/NXinstrument/NXdetector/data
time_binning -> /NXentry/NXinstrument/NXdetector/time_binning

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXreftof.nxdl.xml

### 3.4.10 NXsas

**Status:**

application definition, extends `NXobject`, version 1.0b

**Description:**

This is an application definition for raw data (not processed or reduced data) from a 2-D small angle scattering instrument collected with a monochromatic beam and an area detector. It is meant to be suitable both for neutron SANS and X-ray SAXS data.

It covers all raw data from all SAS techniques: SAS, WSAS, grazing incidence, GISAS

**Symbols:**

No symbol table

**Groups cited:** `NXcollimator, NXdata, NXdetector, NXentry, NXgeometry, NXinstrument, NXmonitor, NXmonochromator, NXsample, NXshape, NXsource`

**Structure:**

(entry): `NXentry`

@entry: `NX_CHAR`

NeXus convention is to use entry1, entry2, ... for analysis software to locate each entry
title: `NX_CHAR`

start_time: `NX_DATE_TIME`

end_time: `NX_DATE_TIME`

definition: `NX_CHAR`

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXsas

instrument: `NXinstrument`

name: `NX_CHAR`

Name of the instrument actually used to perform the experiment

source: `NXsource`

type: `NX_CHAR`
type of radiation source

**name**: *NX_CHAR*

Name of the radiation source

**probe**: *NX_CHAR*

Any of these values: **neutron**, x-ray

**monochromator**: *NXmonochromator*

**wavelength**: *NX_FLOAT {units=NX_WAVELENGTH}*

The wavelength of the radiation

**wavelength_spread**: *NX_FLOAT*

\(\Delta \lambda / \lambda\): Important for resolution calculations

**collimator**: *NXcollimator*

**geometry**: *NXgeometry*

**shape**: *NXshape*

**shape**: *NX_CHAR*

Any of these values: nxylinder, nxbox

**size**: *NX_FLOAT {units=NX_LENGTH}*

The collimation length

**detector**: *NXdetector*

**data[nXPxel, nYPixel]**: *NX_NUMBER*

This is area detector data, of number of x-pixel versus number of y-pixels. Since the beam center is to be determined as a step of data reduction, it is not necessary to document or assume the position of the beam center in acquired data.

**distance**: *NX_FLOAT {units=NX_LENGTH}*

The distance between detector and sample

**x_pixel_size**: *NX_FLOAT {units=NX_LENGTH}*

Physical size of a pixel in x-direction

**y_pixel_size**: *NX_FLOAT {units=NX_LENGTH}*

Size of a pixel in y direction

**polar_angle**: *NX_FLOAT {units=NX_ANGLE}*

**azimuthal_angle**: *NX_FLOAT {units=NX_ANGLE}*

**rotation_angle**: *NX_FLOAT {units=NX_ANGLE}*

**aequatorial_angle**: *NX_FLOAT {units=NX_ANGLE}*

**beam_center_x**: *NX_FLOAT {units=NX_LENGTH}*

This is the x position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

**beam_center_y**: *NX_FLOAT {units=NX_LENGTH}*


This is the y position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

**sample:** `NXsample`

  **name:** `NX_CHAR`
  
  Descriptive name of sample

  **aequatorial_angle:** `NX_FLOAT` `[units=NX_ANGLE]`

**control:** `NXmonitor`

  **mode:** `NX_CHAR`
  
  Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

  *Any of these values:* monitor | timer

  **preset:** `NX_FLOAT`
  
  preset value for time or monitor

  **integral:** `NX_FLOAT` `[units=NX_ANY]`
  
  Total integral monitor counts

**data:** `NXdata`

  **data** -> /NXentry/NXinstrument/NXdetector/data

---

**3.4.11 NXsastof**

**Status:**

  application definition, extends `NXobject`, version 1.0b

**Description:**

  This is an application definition for small angle scattering using a 2D detector in TOF mode. It strives to cover all the SAS techniques in the file again

**Symbols:**

  No symbol table

**Groups cited:** `NXcollimator, NXdata, NXdetector, NXentry, NXgeometry, NXinstrument, NXmonitor, NXsample, NXshape, NXsource`

**Structure:**

  (entry): `NXentry`

    @entry: `NX_CHAR`

    NeXus convention is to use “entry1”, “entry2”, ... for analysis software to locate each entry

    **title:** `NX_CHAR`

    **start_time:** `NX_DATE_TIME`

    **definition:** `NX_CHAR`
Official NeXus NXDL schema to which this file conforms

Obligatory value: NXsastof

**instrument**: NXinstrument

**name**: NX_CHAR

Name of the instrument actually used to perform the experiment

**source**: NXsource

**type**: NX_CHAR

type of radiation source

**name**: NX_CHAR

Name of the radiation source

**probe**: NX_CHAR

Any of these values: neutron|x-ray

**collimator**: NXcollimator

**geometry**: NXgeometry

**shape**: NXshape

**shape**: NX_CHAR

Any of these values: nxcylinder|nxbox

**size**: NX_FLOAT \{units=NX_LENGTH\}

The collimation length

**detector**: NXdetector

**data**: NX_NUMBER

This is area detector data, of number of x-pixel versus number of y-pixels. Since the beam center is to be determined as a step of data reduction, it is not necessary to document or assume the position of the beam center in acquired data.

**time_of_flight**: NX_FLOAT \{units=NX_TIME_OF_FLIGHT\}

**distance**: NX_FLOAT \{units=NX_LENGTH\}

The distance between detector and sample

**x_pixel_size**: NX_FLOAT \{units=NX_LENGTH\}

Physical size of a pixel in x-direction

**y_pixel_size**: NX_FLOAT \{units=NX_LENGTH\}

Size of a pixel in y direction

**polar_angle**: NX_FLOAT \{units=NX_ANGLE\}

**azimuthal_angle**: NX_FLOAT \{units=NX_ANGLE\}

**rotation_angle**: NX_FLOAT \{units=NX_ANGLE\}

**aequatorial_angle**: NX_FLOAT \{units=NX_ANGLE\}

**beam_center_x**: NX_FLOAT \{units=NX_LENGTH\}
This is the x position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

**beam_center_y**: `NX_FLOAT {units=NX_LENGTH}`

This is the y position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

**sample**: `NXSample`

**name**: `NX_CHAR`

Descriptive name of sample

**aequatorial_angle**: `NX_FLOAT {units=NX_ANGLE}`

**control**: `NXMonitor`

**mode**: `NX_CHAR`

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: `monitor | timer`

**preset**: `NX_FLOAT`

preset value for time or monitor

**data[nTOF]**: `NX_INT`

**time_of_flight[nTOF]**: `NX_FLOAT {units=NX_TIME_OF_FLIGHT}`

**data**: `NXdata`

data -> /NXentry/NXinstrument/NXdetector/data

time_of_flight -> /NXentry/NXinstrument/NXdetector/time_of_flight

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXSastof.nxdl.xml

### 3.4.12 NXscan

**Status**:

Application definition, extends `NXObject`, version 1.0b

**Description**:

Application definition for a generic scan instrument. This definition is more an example then a stringent definition as the content of a given NeXus scan file needs to differ for different types of scans. This example definition shows a scan like done on a rotation camera: the sample is rotated and a detector image, the rotation angle and a monitor value is stored at each step in the scan. In the following, the symbol `NP` is used to represent the number of scan points. These are the rules for storing scan data in NeXus files which are implemented in this example:

- Each value varied throughout a scan is stored as an array of length `NP` at its respective location within the NeXus hierarchy.

- For area detectors, `NP` is the first dimension, example for a detector of 256x256: data[256,256]

- The NXdata group contains links to all variables varied in the scan and the data. This to give an equivalent to the more familiar classical tabular representation of scans.
These rules exist for a reason: HDF allows the first dimension of a data set to be unlimited. This means the data can be appended too. Thus a NeXus file built according to the rules given above can be used in the following way:

- At the start of a scan, write all the static information.
- At each scan point, append new data from varied variables and the detector to the file.

Symbols:

No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample

Structure:

(entry): NXentry
  title: NX_CHAR
  start_time: NX_DATE_TIME
  end_time: NX_DATE_TIME
  definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms
    Obligatory value: NXscan

(instrument): NXinstrument

(detector): NXdetector
  data[NP, xdim, ydim]: NX_INT

(sample): NXsample
  rotation_angle[NP]: NX_FLOAT

(monitor): NXmonitor
  data[NP]: NX_INT

(data): NXdata
  data -> /NXentry/NXinstrument/NXdetector/data
  rotation_angle -> /NXentry/NXsample/rotation_angle

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXscan.nxdl.xml

3.4.13 NXspe

Status:
  application definition, extends NXobject, version 1.0

Description:
  NXSPE Inelastic Format. Application definition for NXSPE file format.

Symbols:
  No symbol table

Groups cited: NXcollection, NXdata, NXentry, NXfermi_chopper, NXinstrument, NXsample

Structure:
(entry): NXentry

  program_name: NX_CHAR

  definition: NX_CHAR

    Official NeXus NXDL schema to which this file conforms.

    Any of these values: NXSPE | NXspe

    @version: NX_CHAR

NXSPE_info: Nxcollection

  fixed_energy: NX_FLOAT {units=NX_ENERGY}

    The fixed energy used for this file.

  ki_over_kf_scaling: NX_BOOLEAN

    Indicates whether ki/kf scaling has been applied or not.

  psi: NX_FLOAT {units=NX_ANGLE}

    Orientation angle as expected in DCS-MSlice

data: NXdata

  azimuthal: NX_FLOAT {units=NX_ANGLE}

  azimuthal_width: NX_FLOAT {units=NX_ANGLE}

  polar: NX_FLOAT {units=NX_ANGLE}

  polar_width: NX_FLOAT {units=NX_ANGLE}

  distance: NX_FLOAT {units=NX_LENGTH}

  data: NX_NUMBER

  error: NX_NUMBER

  energy: NX_FLOAT {units=NX_ENERGY}

(instrument): NXinstrument

  name: NX_CHAR

  (fermi_chopper): NXfermi_chopper

    energy: NX_NUMBER {units=NX_ENERGY}

(sample): NXsample

  rotation_angle: NX_NUMBER {units=NX_ANGLE}

  seblock: NX_CHAR

  temperature: NX_NUMBER {units=NX_TEMPERATURE}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXspe.nxdl.xml

3.4.14 NXsqom

Status:

  application definition, extends NXobject, version 1.0b

Description:
This is the application definition for S(Q,OM) processed data. As this kind of data is in general not on a rectangular grid after data reduction, it is stored as Q,E positions plus their intensity, table like. It is the task of a possible visualisation program to regrid this data in a sensible way.

Symbols:

No symbol table

Groups cited: NXdata, NXentry, NXinstrument, NXparameters, NXprocess, NXsample, NXsource

Structure:

(entry): NXentry

@entry: NX_CHAR
title: NX_CHAR
definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms
    Obligatory value: NXsqom

instrument: NXinstrument

name: NX_CHAR
    Name of the instrument from which this data was reduced.

(source): NXsource

type: NX_CHAR
name: NX_CHAR
probe: NX_CHAR
    Any of these values: neutron|x-ray|electron

(sample): NXsample

name: NX_CHAR
    Descriptive name of sample

reduction: NXprocess

program: NX_CHAR
version: NX_CHAR
input: NXparameters
    Input parameters for the reduction program used

filenames: NX_CHAR
    Raw data files used to generate this I(Q)

output: NXparameters
    Eventual output parameters from the data reduction program used

(data): NXdata

data[NP]: NX_INT
    This is the intensity for each point in QE

qx[NP]: NX_CHAR {units=NX_WAVENUMBER}
Positions for the first dimension of Q

\( q_y[NP] \):  \textit{NX_CHAR} \{\text{units=}NX\_WAVENUMBER\}

Positions for the the second dimension of Q

\( q_z[NP] \):  \textit{NX_CHAR} \{\text{units=}NX\_WAVENUMBER\}

Positions for the the third dimension of Q

\( e[n] \):  \textit{NX\_FLOAT} \{\text{units=}NX\_ENERGY\}

Values for the energy transfer for each point

\textbf{Source:} Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXsqom.nxdl.xml

### 3.4.15 \textit{NXtas}

\textbf{Status:}

application definition, extends \textit{NXobject}, version 1.0b

\textbf{Description:}

This is an application definition for a triple axis spectrometer. It is for the trademark scan of the TAS, the Q-E scan. For your alignment scans use the rules in NXscan.

\textbf{Symbols:}

No symbol table

\textbf{Groups cited:} \textit{NXcrystal, NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample, NXsource}

\textbf{Structure:}

\textit{entry}: \textit{NXentry}

\textit{title}: \textit{NX_CHAR}

\textit{start\_time}: \textit{NX\_DATE\_TIME}

\textit{definition}: \textit{NX\_CHAR}

Official NeXus NXDL schema to which this file conforms

Obligatory value: \textit{NXtas}

\textbf{(instrument)}: \textit{NXinstrument}

\textbf{(source)}: \textit{NXsource}

\textit{name}: \textit{NX\_CHAR}

\textit{probe}: \textit{NX\_CHAR}

Any of these values: neutron, x-ray

\textit{monochromator}: \textit{NXcrystal}

\textit{ei}[NP]: \textit{NX\_FLOAT} \{\text{units=}NX\_ENERGY\}

\textit{rotation\_angle}[NP]: \textit{NX\_FLOAT} \{\text{units=}NX\_ANGLE\}

\textit{analyser}: \textit{NXcrystal}
ef[np]: NX_FLOAT {units=NX_ENERGY}
rotation_angle[np]: NX_FLOAT {units=NX_ANGLE}
polar_angle[np]: NX_FLOAT {units=NX_ANGLE}
(detector): NXdetector
data[np]: NX_INT
polar_angle[np]: NX_FLOAT {units=NX_ANGLE}
(same): NXsample
name: NX_CHAR
Descriptive name of sample
qh[np]: NX_FLOAT {units=NX_DIMENSIONLESS}
qk[np]: NX_FLOAT {units=NX_DIMENSIONLESS}
ql[np]: NX_FLOAT {units=NX_DIMENSIONLESS}
en[np]: NX_FLOAT {units=NX_ENERGY}
rotation_angle[np]: NX_FLOAT {units=NX_ANGLE}
polar_angle[np]: NX_FLOAT {units=NX_ANGLE}
sgu[np]: NX_FLOAT {units=NX_ANGLE}
sgl[np]: NX_FLOAT {units=NX_ANGLE}
unit_cell[6]: NX_FLOAT {units=NX_LENGTH}
orientation_matrix[9]: NX_FLOAT {units=NX_DIMENSIONLESS}
(monitor): NXmonitor
mode: NX_CHAR
Count to a preset value based on either clock time (timer) or received monitor
counts (monitor).
Any of these values: monitor|timer
preset: NX_FLOAT
preset value for time or monitor
data[np]: NX_FLOAT {units=NX_ANY}
Total integral monitor counts
(data): NXdata
One of the ei,ef,qh,qk,ql,en should get a primary=1 attribute to denote the main scan
axis
ei -> /NXentry/NXinstrument/monochromator:NXcrystal/ei
ef -> /NXentry/NXinstrument/analyzer:NXcrystal/ef
en -> /NXentry/NXsample/en
qh -> /NXentry/NXsample/qh
qk -> /NXentry/NXsample/qk
ql -> /NXentry/NXsample/ql
3.4.16 NXtofnpd

Status:
application definition, extends NXobject, version 1.0b

Description:
This is a application definition for raw data from a TOF neutron powder diffractometer

Symbols:
No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample, NXuser

Structure:
entry: NXentry
   title: NX_CHAR
   start_time: NX_DATE_TIME
   definition: NX_CHAR
      Official NeXus NXDL schema to which this file conforms
      Obligatory value: NXtofnpd
   pre_sample_flightpath: NX_FLOAT {units=NX_LENGTH}
      This is the flight path before the sample position. This can be determined by a chopper, by the moderator or the source itself. In other words: it the distance to the component which gives the T0 signal to the detector electronics. If another component in the NXinstrument hierarchy provides this information, this should be a link.

user: NXuser
   name: NX_CHAR

(instrument): NXinstrument
   detector: NXdetector
      data[ndet, ntimechan]: NX_INT
      detector_number[ndet]: NX_INT
      distance[ndet]: NX_FLOAT {units=NX_LENGTH}
         distance to sample for each detector
      time_of_flight[ntimechan]: NX_FLOAT {units=NX_TIME_OF_FLIGHT}
      polar_angle[ndet]: NX_FLOAT {units=NX_ANGLE}
         polar angle for each detector element
      azimuthal_angle[ndet]: NX_FLOAT {units=NX_ANGLE}
         azimuthal angle for each detector element

(sample): NXsample
name: $NX_CHAR$

Descriptive name of sample

(monitor): $NXmonitor$

mode: $NX_CHAR$

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: monitor \| timer

preset: $NX_FLOAT$

preset value for time or monitor
distance: $NX_FLOAT$ \{units=$NX_LENGTH$\}
data[ntimechan]: $NX_INT$
time_of_flight[ntimechan]: $NX_FLOAT$ \{units=$NX_TIME_OF_FLIGHT$\}
data: $NXdata$

data $\rightarrow$ /NXentry/NXinstrument/NXdetector/data
detector_number $\rightarrow$ /NXentry/NXinstrument/NXdetector/detector_number
time_of_flight $\rightarrow$ /NXentry/NXinstrument/NXdetector/time_of_flight

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXtofnpd.nxdl.xml

### 3.4.17 NXtofraw

Status:

application definition, extends $NXobject$, version 1.0b

Description:

This is an application definition for raw data from a generic TOF instrument

Symbols:

No symbol table

Groups cited: $NXdata$, $NXdetector$, $NXentry$, $NXinstrument$, $NXmonitor$, $NXsample$, $NXuser$

Structure:

entry: $NXentry$

title: $NX_CHAR$
start_time: $NX_DATE_TIME$
definition: $NX_CHAR$

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXtofraw
duration: $NX_FLOAT$

run_number: $NX_INT$

pre_sample_flightpath: $NX_FLOAT$ \{units=$NX_LENGTH$\}
This is the flight path before the sample position. This can be determined by a chopper, by the moderator, or the source itself. In other words: it is the distance to the component which gives the T0 signal to the detector electronics. If another component in the NXInstrument hierarchy provides this information, this should be a link.

**user:** *NXuser*

**name:** *NX_CHAR*

**instrument:** *NXinstrument*

**detector:** *NXdetector*

- **data[ndet, ntimechan]:** *NX_INT*
- **detector_number[ndet]:** *NX_INT*
- **distance[ndet]:** *NX_FLOAT* {units=*NX_LENGTH*}
  - distance to sample for each detector
- **time_of_flight[ntimechan]:** *NX_FLOAT* {units=*NX_TIME_OF_FLIGHT*}
- **polar_angle[ndet]:** *NX_FLOAT* {units=*NX_ANGLE*}
  - polar angle for each detector element
- **azimuthal_angle[ndet]:** *NX_FLOAT* {units=*NX_ANGLE*}
  - polar angle for each detector element

**(sample):** *NXsample*

- **name:** *NX_CHAR*
  - Descriptive name of sample
- **nature:** *NX_CHAR*
  - Any of these values: powder\|liquid\|single crystal

**(monitor):** *NXmonitor*

- **mode:** *NX_CHAR*
  - Count to a preset value based on either clock time (timer) or received monitor counts (monitor).
  - Any of these values: monitor\|timer
- **preset:** *NX_FLOAT*
  - preset value for time or monitor
- **distance:** *NX_FLOAT* {units=*NX_LENGTH*}
- **data[ntimechan]:** *NX_INT*
- **time_of_flight[ntimechan]:** *NX_FLOAT* {units=*NX_TIME_OF_FLIGHT*}
- **integral_counts:** *NX_INT* {units=*NX_UNITLESS*}

**data:** *NXdata*

- **data** → /NXentry/NXinstrument/NXdetector/data
- **detector_number** → /NXentry/NXinstrument/NXdetector/detector_number
- **time_of_flight** → /NXentry/NXinstrument/NXdetector/time_of_flight
3.4.18 NXtofsingle

Status:
application definition, extends NXobject, version 1.0b

Description:
This is a application definition for raw data from a generic TOF instrument

Symbols:
No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample, NXuser

Structure:
entry: NXentry
  title: NX_CHAR
  start_time: NX_DATE_TIME
  definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms
    Obligatory value: NXtofsingle
  duration: NX_FLOAT
  pre_sample_flightpath: NX_FLOAT {units=NX_LENGTH}
    This is the flight path before the sample position. This can be determined by a chopper, by the moderator or the source itself. In other words: it the distance to the component which gives the T0 signal to the detector electronics. If another component in the NXinstrument hierarchy provides this information, this should be a link.

user: NXuser
  name: NX_CHAR

(instrument): NXinstrument
  detector: NXdetector
    data[xsize, ysize, ntimechan]: NX_INT
    distance[1]: NX_FLOAT {units=NX_LENGTH}
      Distance to sample for the center of the detector
    time_of_flight[ntimechan]: NX_FLOAT {units=NX_TIME_OF_FLIGHT}
    polar_angle[ndet]: NX_FLOAT {units=NX_ANGLE}
      polar angle for each detector element
    azimuthal_angle[ndet]: NX_FLOAT {units=NX_ANGLE}
      azimuthal angle for each detector element

(sample): NXsample
  name: NX_CHAR
Descriptive name of sample

**nature**: *NX_CHAR*

Any of these values: powder | liquid | single crystal

**(monitor)**: *NXmonitor*

**mode**: *NX_CHAR*

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: monitor | timer

**preset**: *NX_FLOAT*

preset value for time or monitor

**distance**: *NX_FLOAT* {units=NX_LENGTH}

**data[ntimechan]**: *NX_INT*

**time_of_flight[ntimechan]**: *NX_FLOAT* {units=NX_TIME_OF_FLIGHT}

**data**: *NXdata*

data → /NXentry/NXinstrument/NXdetector/data
detector_number → /NXentry/NXinstrument/NXdetector/detector_number
time_of_flight → /NXentry/NXinstrument/NXdetector/time_of_flight

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXtofsingle.nxdl.xml

### 3.4.19 NXtomo

**Status**:

application definition, extends *NXobject*, version 2.0

**Description**:

This is the application definition for x-ray or neutron tomography raw data. In tomography a number of dark field images are measured, some bright field images and, of course the sample. In order to distinguish between them images carry a image_key.

**Symbols**:

These symbols will be used below to coordinate datasets with the same shape.

- **nFrames**: number of frames
- **xsize**: number of pixels in X direction
- **ysize**: number of pixels in Y direction

**Groups cited**: *NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample, NXsource*

**Structure**:

- **entry**: *NXentry*
  - **title**: *NX_CHAR*
  - **start_time**: *NX_DATE_TIME*
  - **end_time**: *NX_DATE_TIME*
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

**definition:** *NX_CHAR*

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXtomo

**instrument:** *NXinstrument*

(source): *NXsource*

*type:* *NX_CHAR*

*name:* *NX_CHAR*

*probe:* *NX_CHAR*

Any of these values: neutron|x-ray|electron

**detector:** *NXdetector*

data[nFrames, xsize, ysize]: *NX_INT*

image_key[nFrames]: *NX_INT*

In order to distinguish between sample projections, dark and flat images, a magic number is recorded per frame. The key is as follows:

- projection = 0
- flat field = 1
- dark field = 2
- invalid = 3

*x_pixel_size*: *NX_FLOAT* {units=*NX_LENGTH*}

*y_pixel_size*: *NX_FLOAT* {units=*NX_LENGTH*}

*distance*: *NX_FLOAT* {units=*NX_LENGTH*}

Distance between detector and sample

*x_rotation_axis_pixel_position*: *NX_FLOAT*

*y_rotation_axis_pixel_position*: *NX_FLOAT*

**sample:** *NXsample*

*name:* *NX_CHAR*

Descriptive name of sample

*rotation_angle*[nFrames]: *NX_FLOAT* {units=*NX_ANGLE*}

In practice this axis is always aligned along one pixel direction on the detector and usually vertical. There are experiments with horizontal rotation axes, so this would need to be indicated somehow. For now the best way for that is an open question.

*x_translation*[nFrames]: *NX_FLOAT* {units=*NX_LENGTH*}

*y_translation*[nFrames]: *NX_FLOAT* {units=*NX_LENGTH*}

*z_translation*[nFrames]: *NX_FLOAT* {units=*NX_LENGTH*}

**control:** *NXmonitor*

*data*[nFrames]: *NX_FLOAT* {units=*NX_ANY*}
Total integral monitor counts for each measured frame. Allows a to correction for fluctuations in the beam between frames.

data: NXdata
    data → /NXentry/NXinstrument/detector:NXdetector/data
    rotation_angle → /NXentry/NXsample/rotation_angle
    image_key → /NXentry/NXinstrument/detector:NXdetector/image_key

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXtomo.nxdl.xml

3.4.20 NXtomophase

Status:
application definition, extends NXobject, version 1.0b

Description:
This is the application definition for x-ray or neutron tomography raw data with phase contrast variation at each point. In tomography first some dark field images are measured, some bright field images and, of course the sample. In order to properly sort the order of the images taken, a sequence number is stored with each image.

Symbols:
These symbols will be used below to coordinate datasets with the same shape.

nBrightFrames: number of bright frames
nDarkFrames: number of dark frames
nSampleFrames: number of image (sample) frames
nPhase: number of phase settings
xsize: number of pixels in X direction
ysize: number of pixels in Y direction

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXsample, NXsource

Structure:
entry: NXentry
    title: NX_CHAR
    start_time: NX_DATE_TIME
    end_time: NX_DATE_TIME
    definition: NX_CHAR
        Official NeXus NXDL schema to which this file conforms
        Obligatory value: NXtomophase
    instrument: NXinstrument
        (source): NXsource
**NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1**

- **type**: `NX_CHAR`
- **name**: `NX_CHAR`
- **probe**: `NX_CHAR`
  
  Any of these values: `neutron`, `x-ray`, `electron`

- **bright_field**: `NXdetector`

  - **data**[nBrightFrames, xsize, ysize]: `NX_INT`
  - **sequence_number**[nBrightFrames]: `NX_INT`

- **dark_field**: `NXdetector`

  - **data**[nDarkFrames, xsize, ysize]: `NX_INT`
  - **sequence_number**[nDarkFrames]: `NX_INT`

- **sample**: `NXsample`

  - **data**[nSampleFrames, nPhase, xsize, ysize]: `NX_INT`
  - **sequence_number**[nSampleFrames, nPhase]: `NX_INT`
  - **x_pixel_size**: `NX_FLOAT` {units=`NX_LENGTH`}
  - **y_pixel_size**: `NX_FLOAT` {units=`NX_LENGTH`}
  - **distance**: `NX_FLOAT` {units=`NX_LENGTH`}

  Distance between detector and sample

- **sample**: `NXsample`

  - **name**: `NX_CHAR`
  
  Descriptive name of sample

  - **rotation_angle**[nSampleFrames]: `NX_FLOAT` {units=`NX_ANGLE`}
  - **x_translation**[nSampleFrames]: `NX_FLOAT` {units=`NX_LENGTH`}
  - **y_translation**[nSampleFrames]: `NX_FLOAT` {units=`NX_LENGTH`}
  - **z_translation**[nSampleFrames]: `NX_FLOAT` {units=`NX_LENGTH`}

- **control**: `NXmonitor`

  - **integral**[nDarkFrames + nBrightFrames + nSampleFrame]: `NX_FLOAT` {units=`NX_ANY`}

  Total integral monitor counts for each measured frame. Allows a correction for fluctuations in the beam between frames.

- **data**: `NXdata`

  - **data** → `/NXentry/NXinstrument/sample:NXdetector/data`

  - **rotation_angle** → `/NXentry/NXsample/rotation_angle`

**Source**: Automatically generated from `https://github.com/nexusformat/definitions/blob/master/applications/NXtomophase.nxdl.xml`
3.4.21 NXtomoproc

Status:

application definition, extends *NXobject*, version 1.0b

Description:

This is an application definition for the final result of a tomography experiment: a 3D construction of some volume of physical properties.

Symbols:

These symbols will be used below to coordinate datasets with the same shape.

nx: number of voxels in X direction
ny: number of voxels in Y direction
nz: number of voxels in Z direction

Groups cited: *NXdata, NXentry, NXinstrument, NXparameters, NXprocess, NXsample, NXsource*

Structure:

entry: *NXentry*

  title: *NX_CHAR*

  definition: *NX_CHAR*

    Official NeXus NXDL schema to which this file conforms

    Obligatory value: NXtomoproc

  (instrument): *NXinstrument*

    type: *NX_CHAR*

    name: *NX_CHAR*

    probe: *NX_CHAR*

      Any of these values: neutron|x-ray|electron

  (source): *NXsource*

    type: *NX_CHAR*

    name: *NX_CHAR*

  (sample): *NXsample*

    name: *NX_CHAR*

      Descriptive name of sample

  reconstruction: *NXprocess*

    program: *NX_CHAR*

      Name of the program used for reconstruction

    version: *NX_CHAR*

      Version of the program used

    date: *NX_DATE_TIME*

      Date and time of reconstruction processing.

  parameters: *NXparameters*

    raw_file: *NX_CHAR*
Original raw data file this data was derived from

data: \textit{NXdata}

\texttt{data[nx, nx, nz]}: \textit{NX\_INT}

This is the reconstructed volume. This can be different things. Please indicate in the unit attribute what physical quantity this really is.

@transform: \textit{NX\_CHAR}

@offset: \textit{NX\_CHAR}

@scaling: \textit{NX\_CHAR}

\texttt{x[nx]}: \textit{NX\_FLOAT} \{units=\textit{NX\_ANY}\}

This is an array holding the values to use for the x-axis of data. The units must be appropriate for the measurement.

\texttt{y[ny]}: \textit{NX\_FLOAT} \{units=\textit{NX\_ANY}\}

This is an array holding the values to use for the y-axis of data. The units must be appropriate for the measurement.

\texttt{z[nz]}: \textit{NX\_FLOAT} \{units=\textit{NX\_ANY}\}

This is an array holding the values to use for the z-axis of data. The units must be appropriate for the measurement.

\textbf{Source}: Automatically generated from \url{https://github.com/nexusformat/definitions/blob/master/applications/NXtomoproc.nxdl.xml}

### 3.4.22 NXxas

\textbf{Status}:

application definition, extends \textit{NXobject}, version 1.0

\textbf{Description}:

This is an application definition for raw data from an X-ray absorption spectroscopy experiment. This is essentially a scan on energy versus incoming/absorbed beam.

\textbf{Symbols}:

No symbol table

\textbf{Groups cited}: \textit{NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXmonochromator, NXsample, NXsource}

\textbf{Structure}:

\begin{itemize}
  \item (entry): \textit{NXentry}
  \begin{itemize}
    \item @entry: \textit{NX\_CHAR}
      
      NeXus convention is to use “entry1”, “entry2”, ... for analysis software to locate each entry.
    \item title: \textit{NX\_CHAR}
    \item start\_time: \textit{NX\_DATE\_TIME}
    \item definition: \textit{NX\_CHAR}
      
      Official NeXus NXDL schema to which this file conforms
  \end{itemize}
  \end{itemize}

\textbf{3.4. Application Definitions}
(instrument): \textit{NXinstrument}

(source): \textit{NXsource}

type: \textit{NX\_CHAR}

name: \textit{NX\_CHAR}

probe: \textit{NX\_CHAR}

Obligatory value: \textit{x-ray}

monochromator: \textit{NXmonochromator}

energy[np]: \textit{NX\_FLOAT}

incoming\_beam: \textit{NXdetector}

\text{data[np]:} \textit{NX\_INT}

absorbed\_beam: \textit{NXdetector}

\text{data[np]:} \textit{NX\_INT}

mark this field with attribute \texttt{signal}=1

(sample): \textit{NXsample}

name: \textit{NX\_CHAR}

Descriptive name of sample

(monitor): \textit{NXmonitor}

mode: \textit{NX\_CHAR}

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: \texttt{monitor}\mid\texttt{timer}

preset: \textit{NX\_FLOAT}

preset value for time or monitor

\text{data[np]:} \textit{NX\_INT}

(data): \textit{NXdata}

energy \rightarrow /entry/instrument/monochromator/energy

absorbed\_beam \rightarrow /entry/instrument/absorbed\_beam/data

Source: Automatically generated from \url{https://github.com/nexusformat/definitions/blob/master/applications/NXxas.nxdl.xml}

3.4.23 \textbf{NXxasproc}

Status:

application definition, extends \textit{NXobject}, version 1.0

Description:

This is an application definition for processed data from XAS. This is energy versus \( \frac{I(\text{incoming})}{I(\text{absorbed})} \).

Symbols:

No symbol table
Groups cited: NXdata, NXentry, NXparameters, NXprocess, NXsample

Structure:

(entry): NXentry

@entry: NX_CHAR

NeXus convention is to use “entry1”, “entry2”, ... for analysis software to locate each entry.

title: NX_CHAR

definition: NX_CHAR

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXasproc

(sample): NXsample

name: NX_CHAR

Descriptive name of sample

XAS_data_reduction: NXprocess

program: NX_CHAR

Name of the program used for reconstruction

version: NX_CHAR

Version of the program used

date: NX_DATE_TIME

Date and time of reconstruction processing.

parameters: NXparameters

raw_file: NX_CHAR

Original raw data file this data was derived from

(data): NXdata

energy[np]: NX_CHAR

data[np]: NX_FLOAT

This is corrected and calibrated I(incoming)/I(absorbed). So it is the absorption. Expect attribute signal=1

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXasproc.nxdl.xml

3.4.24 NXxbase

Status:

application definition, extends NXobject, version 1.0b

Description:

This definition covers the common parts of all monochromatic single crystal raw data application definitions.

Symbols:
No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXmonitor, NXmonochromator, NXsample, NXsource

Structure:

**entry**: NXentry

**title**: NX_CHAR

**start_time**: NX_DATE_TIME

**definition**: NX_CHAR

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXxbase

**instrument**: NXinstrument

**source**: NXsource

**type**: NX_CHAR

**name**: NX_CHAR

**probe**: NX_CHAR

Any of these values: neutron|x-ray|electron

**monochromator**: NXmonochromator

**wavelength**: NX_FLOAT {units=NX_WAVELENGTH}

**detector**: NXdetector

The name of the group is detector if there is only one detector, if there are several, names have to be detector1, detector2, ...detector{n}

**data[np, number of x pixels, number of y pixels]**: NX_INT

The area detector data, the first dimension is always the number of scan points, the second and third are the number of pixels in x and y. The origin is always assumed to be in the center of the detector. maxOccurs is limited to the number of detectors on your instrument.

**@signal**: NX_POSINT

Obligatory value: 1

**x_pixel_size**: NX_FLOAT {units=NX_LENGTH}

**y_pixel_size**: NX_FLOAT {units=NX_LENGTH}

**distance**: NX_FLOAT {units=NX_LENGTH}

**frame_start_number**: NX_INT

This is the start number of the first frame of a scan. In PX one often scans a couple of frames on a give sample, then does something else, then returns to the same sample and scans some more frames. Each time with a new data file. This number helps concatenating such measurements.

**sample**: NXsample

**name**: NX_CHAR

Descriptive name of sample
orientation_matrix[3, 3]: *NX_FLOAT*

The orientation matrix according to Busing and Levy conventions. This is not strictly necessary as the UB can always be derived from the data. But let us bow to common usage which includes the UB nearly always.

unit_cell[6]: *NX_FLOAT*

The unit cell, a, b, c, alpha, beta, gamma. Again, not strictly necessary, but normally written.

temperature[NP]: *NX_FLOAT*

The sample temperature or whatever sensor represents this value best

x_translation: *NX_FLOAT* {units=*NX_LENGTH*}

Translation of the sample along the X-direction of the laboratory coordinate system

y_translation: *NX_FLOAT* {units=*NX_LENGTH*}

Translation of the sample along the Y-direction of the laboratory coordinate system

distance: *NX_FLOAT* {units=*NX_LENGTH*}

Translation of the sample along the Z-direction of the laboratory coordinate system

control: *NXmonitor*

mode: *NX_CHAR*

Count to a preset value based on either clock time (timer) or received monitor counts (monitor).

Any of these values: monitor | timer

preset: *NX_FLOAT*

preset value for time or monitor

integral: *NX_FLOAT* {units=*NX_ANY*}

Total integral monitor counts

(data): *NXdata*

The name of this group id data if there is only one detector; if there are several the names will be data1, data2, data3 and will point to the corresponding detector groups in the instrument hierarchy.

data → /NXentry/NXinstrument/NXdetector/data

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXxbase.nxdl.xml

### 3.4.25 NXxeuler

**Status:**

application definition, extends *NXbase*, version 1.0b

**Description:**
This is the application definition for raw data from a four-circle diffractometer with an eulerian cradle. It extends \textit{NXxbase}, so the full definition is the content of \textit{NXxbase} plus the data defined here. All four angles are logged in order to support arbitrary scans in reciprocal space.

\textbf{Symbols:} \\
No symbol table

\textbf{Groups cited:} \textit{NXdata, NXdetector, NXentry, NXinstrument, NXsample}

\textbf{Structure:} \\
\texttt{entry: NXEntry} \\
\texttt{definition: NX_CHAR} \\
Official NeXus NXDL schema to which this file conforms \\
Obligatory value: \texttt{NXxeuler}

\texttt{instrument: NXinstrument} \\
\texttt{detector: NXdetector} \\
\texttt{polar_angle[np]: NX_FLOAT \{units=NX_ANGLE\}} \\
The polar_angle (two theta) where the detector is placed at each scan point.

\texttt{sample: NXsample} \\
\texttt{rotation_angle[np]: NX_FLOAT \{units=NX_ANGLE\}} \\
This is an array holding the sample rotation angle at each scan point

\texttt{chi[np]: NX_FLOAT \{units=NX_ANGLE\}} \\
This is an array holding the chi angle of the eulerian cradle at each scan point

\texttt{phi[np]: NX_FLOAT \{units=NX_ANGLE\}} \\
This is an array holding the phi rotation of the eulerian cradle at each scan point

\texttt{name: NXdata} \\
polar_angle \rightarrow /NXentry/NXinstrument/NXdetector/polar_angle \\
rotation_angle \rightarrow /NXentry/NXsample/rotation_angle \\
chi \rightarrow /NXentry/NXsample/chi \\
phi \rightarrow /NXentry/NXsample/phi

\textbf{Source:} Automatically generated from \url{https://github.com/nexusformat/definitions/blob/master/applications/NXxeuler.nxdl.xml}

\subsection*{3.4.26 \textit{NXxkappa}}

\textbf{Status:} \\
\texttt{application definition, extends NXxbase, version 1.0b}

\textbf{Description:} \\
This is the application definition for raw data from a kappa geometry (CAD4) single crystal diffractometer. It extends \textit{NXxbase}, so the full definition is the content of \textit{NXxbase} plus the data defined here.

\textbf{Symbols:}
No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXsample

Structure:

entry: NXentry

definition: NX_CHAR

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXxkappa

instrument: NXinstrument
detector: NXdetector

polar_angle[np]: NX_FLOAT {units=NX_ANGLE}

The polar_angle (two theta) at each scan point

sample: NXsample

rotation_angle[np]: NX_FLOAT {units=NX_ANGLE}

This is an array holding the sample rotation angle at each scan point

kappa[np]: NX_FLOAT {units=NX_ANGLE}

This is an array holding the kappa angle at each scan point

phi[np]: NX_FLOAT {units=NX_ANGLE}

This is an array holding the phi angle at each scan point

alpha: NX_FLOAT {units=NX_ANGLE}

This holds the inclination angle of the kappa arm.

name: NXdata

polar_angle -> /NXentry/NXinstrument/NXdetector/polar_angle
rotation_angle -> /NXentry/NXsample/rotation_angle
kappa -> /NXentry/NXsample/kappa
phi -> /NXentry/NXsample/phi

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXxkappa.nxdl.xml

3.4.27 NXxlaue

Status:

application definition, extends NXxrot, version 1.0b

Description:

This is the application definition for raw data from a single crystal laue camera. It extends NXxrot.

Symbols:

No symbol table

Groups cited: NXdata, NXentry, NXinstrument, NXsource

Structure:
entry: NXEntry

definition: NX_CHAR

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXxlaue

instrument: NXinstrument

source: NXsource

distribution: NXdata

This is the wavelength distribution of the beam

data[ne]: NX_CHAR

expect signal=1 axes="energy"

wavelength[ne]: NX_CHAR {units=NX_WAVELENGTH}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXxlaue.nxdl.xml

3.4.28 NXxlaueplate

Status:

application definition, extends NXxlaue, version 1.0b

Description:

This is the application definition for raw data from a single crystal Laue camera with an image plate as a detector. It extends NXxlaue.

Symbols:

No symbol table

Groups cited: NXdetector, NXentry, NXinstrument

Structure:

entry: NXEntry

definition: NX_CHAR

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXxlaueplate

instrument: NXinstrument

detector: NXdetector

diameter: NX_FLOAT {units=NX_LENGTH}

The diameter of a cylindrical detector

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXxlaueplate.nxdl.xml
3.4.29 NXxnb

Status:
application definition, extends NXxbase, version 1.0b

Description:
This is the application definition for raw data from a single crystal diffractometer measuring in normal beam mode. It extends NXxbase, so the full definition is the content of NXxbase plus the data defined here. All angles are logged in order to support arbitrary scans in reciprocal space.

Symbols:
No symbol table

Groups cited: NXdata, NXdetector, NXentry, NXinstrument, NXsample

Structure:

entry: NXentry

definition: NX_CHAR

Official NeXus NXDL schema to which this file conforms

Obligatory value: NXxnb

instrument: NXinstrument
detector: NXdetector

polar_angle[np]: NX_FLOAT {units=NX_ANGLE}
The polar_angle (gamma) of the detector for each scan point.
tilt_angle[np]: NX_FLOAT {units=NX_ANGLE}
The angle by which the detector has been tilted out of the scattering plane.
sample: NXsample

rotation_angle[np]: NX_FLOAT {units=NX_ANGLE}
This is an array holding the sample rotation angle at each scan point

name: NXdata

polar_angle -> /NXentry/NXinstrument/NXdetector/polar_angle
tilt -> /NXentry/NXinstrument/NXdetector/tilt
rotation_angle -> /NXentry/NXsample/rotation_angle

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXxnb.nxdl.xml

3.4.30 NXxrot

Status:
application definition, extends NXxbase, version 1.0b

Description:
This is the application definition for raw data from a rotation camera. It extends NXxbase, so the full definition is the content of NXxbase plus the data defined here.
Symbols:

No symbol table

Groups cited:  *NXattenuator, NXdata, NXdetector, NXentry, NXinstrument, NXsample*

Structure:

```
entry: NXentry
definition: NX_CHAR

Official NeXus NXDL schema to which this file conforms.

Obligatory value: NXxrot

instrument: NXinstrument
detector: NXdetector

polar_angle: NX_FLOAT {units=NX_ANGLE}
The polar_angle (two theta) where the detector is placed.

beam_center_x: NX_FLOAT {units=NX_LENGTH}
This is the x position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

beam_center_y: NX_FLOAT {units=NX_LENGTH}
This is the y position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

attenuator: NXattenuator

attenuator_transmission: NX_FLOAT {units=NX_ANY}

sample: NXsample

rotation_angle[np]: NX_FLOAT {units=NX_ANGLE}
This is an array holding the sample rotation start angle at each scan point

rotation_angle_step[np]: NX_FLOAT {units=NX_ANGLE}
This is an array holding the step made for sample rotation angle at each scan point

name: NXdata

rotation_angle -> /NXentry/NXsample/rotation_angle
```

Source:  Automatically generated from https://github.com/nexusformat/definitions/blob/master/applications/NXxrot.nxdl.xml

### 3.5 Contributed Definitions

A description of each NeXus contributed definition is given. NXDL files in the NeXus contributed definitions include propositions from the community for NeXus base classes or application definitions, as well as other NXDL files for long-term archival by NeXus. Consider the contributed definitions as either in incubation or a special case not for general use. The NIAC: The NeXus International Advisory Committee is charged to review any new contributed definitions and provide feedback to the authors before ratification and acceptance as either a base class or application definition.
### 3.5.1 NXarpes

**Status:**
contributed definition, extends \textit{NXobject}, version 1.0

**Description:**
This is an application definition for angular resolved photo electron spectroscopy. It has been drawn up with hemispherical electron analysers in mind.

**Symbols:**
No symbol table

**Groups cited:** \textit{NXdata, NXdetector, NXentry, NXinstrument, NXmonochromator, NXsample, NXsource}

**Structure:**

\begin{verbatim}
(entry): NXentry
  @entry: NX_CHAR
    NeXus convention is to use “entry1”, “entry2”, ... for analysis software to locate each entry.
  title: NX_CHAR
  start_time: NX_DATE_TIME
  definition: NX_CHAR
    Official NeXus NXDL schema to which this file conforms.
    Obligatory value: NXarpes

(instrument): NXinstrument
  (source): NXsource
    type: NX_CHAR
    name: NX_CHAR
    probe: NX_CHAR
      Obligatory value: x-ray
    monochromator: NXmonochromator
    energy: NX_NUMBER \{units=NXENERGY\}

analyzer: NXdetector
  data: NX_NUMBER
  lens_mode: NX_CHAR
    setting for the electron analyzer lens
  acquisition_mode: NX_CHAR
    Any of these values: swept | fixed
  entrance_slit_shape: NX_CHAR
    Any of these values: curved | straight
  entrance_slit_setting: NX_NUMBER \{units=NXANY\}
\end{verbatim}
dial setting of the entrance slit

**entrance_slit_size:** `NX_CHAR {units=NX_LENGTH}`

size of the entrance slit

**pass_energy:** `NX_CHAR {units=NX_ENERGY}`

energy of the electrons on the mean path of the analyser

**time_per_channel:** `NX_CHAR {units=NX_TIME}`

todo: define more clearly

**angles:** `NX_NUMBER {units=NX_ANGLE}`

Angular axis of the analyser data which dimension the axis applies to is defined using the normal NXdata methods.

**energies:** `NX_NUMBER {units=NX_ENERGY}`

Energy axis of the analyser data which dimension the axis applies to is defined using the normal NXdata methods.

**sensor_size[]:** `NX_INT`

number of raw active elements in fast and slow pixel dimension direction

**region_origin[]:** `NX_INT`

origin of rectangular region selected for readout

**region_size[]:** `NX_INT`

size of rectangular region selected for readout

**(sample):** `NXsample`

**name:** `NX_CHAR`

Descriptive name of sample

**temperature:** `NX_NUMBER {units=NX_TEMPERATURE}`

**(data):** `NXdata`

**Source:** Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXarpes.nxdl.xml

### 3.5.2 NXbeamline

**Status:**

contributed definition, extends `NXobject`, version 1.0

**Description:**

container for elements describing beamline.

**Symbols:**

No symbol table

**Groups cited:** `NXaperture`, `NXbending_magnet`, `NXcollection`, `NXelectrostatic_kicker`, `NXmagnetic_kicker`, `NXquadrupole_magnet`, `NXseparator`, `NXsolenoid_magnet`, `NXspin_rotator`

**Structure:**
beamline: \texttt{NX_CHAR}

name of beamline.

diagnostics: \texttt{NXcollection}

(bending_magnet): \texttt{NXbending_magnet}

(quadrupole_magnet): \texttt{NXquadrupole_magnet}

(solenoid_magnet): \texttt{NXsolenoid_magnet}

(separator): \texttt{NXseparator}

(spin_rotator): \texttt{NXspin_rotator}

(electrostatic_kicker): \texttt{NXelectrostatic_kicker}

(magnetic_kicker): \texttt{NXmagnetic_kicker}

(aperture): \texttt{NXaperture}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXbeamline.nxdl.xml

3.5.3 \textbf{NXcanSAS}

Status:

commited definition, extends \texttt{NXobject}, version 1.0

Description:

Implementation of the canSAS standard to store reduced multi-dimensional small-angle scattering data.

for more details, see:

- http://www.cansas.org/
- http://www.cansas.org/formats/canSAS2012/1.0/doc

Symbols:

No symbol table

Groups cited: \texttt{NXcollection, NXcollimator, NXdata, NXdetector, NXinstrument, NXorientation, NXparameters, NX-process, NXsample, NXsource, NXsubentry, NXtranslation}

Structure:

(subentry): \texttt{NXsubentry}

Place the canSAS SASentry group as a child of a NeXus \texttt{NXentry} group.

Note: It is required for all numerical objects to provide a \textit{units} attribute that describes the engineering units. Use the Unidata UDunits \textsuperscript{3} specification as this is compatible with various community standards.

\texttt{@canSAS_class: NX_CHAR}

Official canSAS group: SASentry

Obligatory value: SASentry

\texttt{@version: NX_CHAR}

\textsuperscript{3} The UDunits specification also includes instructions for derived units.
Describes the version of the canSAS standard used to write this data. This must be a
text (not numerical) representation. Such as:

@version="1.0"

Obligatory value: 1.0

definition: NX_CHAR

Official NeXus NXDL schema to which this subentry conforms.

Obligatory value: NXcanSAS

title: NX_CHAR

Title of this SASentry.

run: NX_CHAR

Run identification for this SASentry. For many facilities, this is an integer. Use
multiple instances of run as needed.

@name: NX_CHAR

Optional string attribute to identify this particular run. Could use this to
associate (correlate) multiple SASdata elements with run elements.

(data): NXdata

A SASData group contains reduced a single small-angle scattering data set that can
be represented as $I(\vec{Q})$ or $I(|\vec{Q}|)$.

![Diagram of SASdata](image)

Figure 3.10: The SASdata element

$Q$ can be either a vector ($\vec{Q}$) or a vector magnitude ($|\vec{Q}|$)
The name of each SASdata must be unique within a SASentry group. Such as sasdata01.

A SASdata group has several attributes:

- \texttt{I\_axes}
- \texttt{Q\_indices}
- \texttt{Mask\_indices}

To indicate the dependency relationships of other varied parameters, use attributes similar to \texttt{@Mask\_indices} (such as \texttt{@Temperature\_indices} or \texttt{@Pressure\_indices}).

\texttt{@canSAS\_class: NX\_CHAR}

Official canSAS group: NXcanSAS (contributed definition); SASdata

Obligatory value: \texttt{SASdata}

\texttt{@I\_axes: NX\_CHAR}

Comma-separated list that describes the names of the data objects that correspond to the indices of the \texttt{I} data object. Such as:

\texttt{@I\_axes=\textquotesingle\textquotesingle;Temperature,Time,Pressure,Q,Q\textquotesingle\textquotesingle;}

Since there are five items in the list, the intensity field \texttt{I} must be a five-dimensional array (rank=5).

\texttt{@Q\_indices: NX\_CHAR}

Array that describes which indices (of the \texttt{I} data object) are used to reference the \texttt{Q} data object. The items in this array use zero-based indexing. Such as:

\texttt{@Q\_indices=1,3,4}

which indicates that \texttt{Q} requires three indices from the \texttt{I} data object: one for time and two for \texttt{Q} position. Thus, in this example, the \texttt{Q} data is time-dependent: $\vec{Q}(t)$.

\texttt{@Mask\_indices: NX\_CHAR}

Array that describes which indices (of the \texttt{I} data object) are used to reference the Mask data object. The items in this array use zero-based indexing. Such as:

\texttt{@Mask\_indices=3,4}

which indicates that \texttt{Q} requires two indices from the \texttt{I} data object for \texttt{Q} position.

\texttt{Q: NX\_NUMBER \{units=NX\_PER\_LENGTH\}}

Array of \texttt{Q} data to accompany \texttt{I}.

\texttt{Q} may be represented either as the three-dimensional scattering vector $\vec{Q}$ or by the magnitude of the scattering vector, $|\vec{Q}|$.

\begin{equation}
|\vec{Q}| = (4\pi/\lambda)\sin(\theta)
\end{equation}

3.5. Contributed Definitions
When we write $Q$, we may refer to either or both of $|\vec{Q}|$ or $\vec{Q}$, depending on the context.

@uncertainty: **NX_CHAR**

Typically the estimated standard deviation. More general, this is the estimate of the uncertainty of each "$\text{math:Q}$.

Can use this to describe the slit-length at each datum. Use a subgroup to describe any supplementary uncertainty data such as two-dimensional.

(optional for numerical arrays) Name of the data object (in this SAS data group) that provides the uncertainty to be used for data analysis.

### I: **NX_NUMBER**

Array of intensity ($I$) data.

The intensity may be represented in one of these forms:

- **absolute units**: $d\Sigma/d\Omega(Q)$ differential cross-section per unit volume per unit solid angle (typical units: $1/\text{cm}/\text{sr}$)
- **absolute units**: $d\sigma/d\Omega(Q)$ differential cross-section per unit atom per unit solid angle (typical units: $\text{cm}^2$)
- **arbitrary units**: $I(Q)$ usually a ratio of two detectors but units are meaningless (typical units: a.u.)

This presents a few problems for analysis software to sort out when reading the data. Fortunately, it is possible to analyze the units to determine which type of intensity is being reported and make choices at the time the file is read. But this is an area for consideration and possible improvement.

One problem arises with software that automatically converts data into some canonical units used by that software. The software should not convert units between these different types of intensity indiscriminately.

A second problem is that when arbitrary units are used, then the set of possible analytical results is restricted. With such units, no meaningful volume fraction or number density can be determined directly from $I(Q)$.
In some cases, it is possible to apply a factor to convert the arbitrary units to an absolute scale. This should be considered as a possibility of the analysis process.

@uncertainty: **NX_CHAR**

Typically the estimated standard deviation. \(1/\sqrt{I}\) for Poisson statistics. More generally, though, this is the estimate of the uncertainty of each \(I\).

(optional for numerical arrays) Name of the data object (in this SAS data group) that provides the uncertainty to be used for data analysis.

\(I_{dev}\) is the canonical name from the 1D standard. The multi-D standard allows for this name to be described in this attribute.

Qmean: **NX_CHAR** \{units=NX_PER_LENGTH\}

Mean value of \(Q\) for this data point. Useful when describing data that has been binned from higher-resolution data. It is unexpected for \(Q\) and Qmean to have different units.

ShadowFactor: **NX_CHAR** \{units=NX_DIMENSIONLESS\}

A numerical factor applied to pixels affected by the beam stop penumbra. Used in data files from NIST/NCNR instruments.


(parameters): **NXparameters**

Supplementary Uncertainty Data

@canSAS_class: **NX_CHAR**

ad hoc canSAS group: NXcanSAS (contributed definition); SASuncertainties

Obligatory value: SASuncertainties

(instrument): **NXinstrument**

This the SAS instrument.

@canSAS_class: **NX_CHAR**

Official canSAS group: NXcanSAS (contributed definition); SASinstrument

Obligatory value: SASinstrument

(collimator): **NXcollimator**

Description of a collimating element in the instrument.

@canSAS_class: **NX_CHAR**

Official canSAS group: NXcanSAS (contributed definition); SAScollimation

Obligatory value: SAScollimation

length: **NX_NUMBER** \{units=NX_LENGTH\}

Amount/length of collimation inserted (as on a SANS instrument)

aperture: **NXparameters**
Figure 3.12: The SASinstrument element

Figure 3.13: The SAScollimation element
Description of a slit or aperture.

@canSAS_class: NX_CHAR

ad hoc canSAS_class.

Obligatory value: container

@name: NX_CHAR

Optional name attribute for this aperture

@type: NX_CHAR

Optional text attribute to describe the type of aperture (pinhole, 4-blade slit, Soller slit, ...).

distance: NX_NUMBER {units=NX_LENGTH}

Distance from this collimation element to the sample.

size: NXparameters

Opening dimensions of this aperture.

@canSAS_class: NX_CHAR

ad hoc canSAS_class.

Obligatory value: container

x: NX_NUMBER {units=NX_LENGTH}

Dimension of the collimation in x.

y: NX_NUMBER {units=NX_LENGTH}

Dimension of the collimation in y.

z: NX_NUMBER {units=NX_LENGTH}

Dimension of the collimation in z.

While z is allowed by the canSAS standard, it may not make sense to use it in some situations. Use of z may be ignored by processing software.

(detector): NXdetector

Description of a detector in the instrument.

@canSAS_class: NX_CHAR

Official canSAS group: NXcanSAS (contributed definition); SASdetector

Obligatory value: SASdetector

name: NX_CHAR

Identifies the name of this detector

SDD: NX_NUMBER {units=NX_LENGTH}

Distance between sample and detector.

slit_length: NX_NUMBER {units=NX_PER_LENGTH}

Slit length of the instrument for this detector, expressed in the same units as Q.
Figure 3.14: The SASdetector element
offset: \textit{NXparameters}

Offset of the detector position.

\texttt{@canSAS\_class: NX\_CHAR}

\textit{ad hoc} canSAS\_class.

\texttt{Obligatory value: container}

\texttt{x: NX\_NUMBER \{units=NX\_LENGTH\}}

\texttt{y: NX\_NUMBER \{units=NX\_LENGTH\}}

\texttt{z: NX\_NUMBER \{units=NX\_LENGTH\}}

While \texttt{z} is allowed by the canSAS standard, it may not make sense to use it in some situations. Use of \texttt{z} may be ignored by processing software.

orientation: \textit{NXparameters}

The orientation element describes simple rotations about a single axis (rather than a full set of rotations as in a crystallographic context).

\texttt{@canSAS\_class: NX\_CHAR}

\textit{ad hoc} canSAS\_class.

\texttt{Obligatory value: container}

\texttt{roll: NX\_NUMBER \{units=NX\_LENGTH\}}

Roll is rotation about the \texttt{z} axis.

\texttt{pitch: NX\_NUMBER \{units=NX\_LENGTH\}}

Pitch is rotation about the \texttt{x} axis.

\texttt{yaw: NX\_NUMBER \{units=NX\_LENGTH\}}

Yaw is rotation about the \texttt{y} axis.

beam\_center: \textit{NXparameters}

Position of the beam center on the detector.

\texttt{@canSAS\_class: NX\_CHAR}

\textit{ad hoc} canSAS\_class.

\texttt{Obligatory value: container}

\texttt{x: NX\_NUMBER \{units=NX\_LENGTH\}}

\texttt{y: NX\_NUMBER \{units=NX\_LENGTH\}}

\texttt{z: NX\_NUMBER \{units=NX\_LENGTH\}}

While \texttt{z} is allowed by the canSAS standard, it may not make sense to use it in some situations. Use of \texttt{z} may be ignored by processing software.

pixel\_size: \textit{NXparameters}

Size of the pixels on this detector.

\texttt{@canSAS\_class: NX\_CHAR}
ad hoc canSAS_class.

Obligatory value: container

x: \texttt{NX\_NUMBER \{units=NX\_LENGTH\}}
y: \texttt{NX\_NUMBER \{units=NX\_LENGTH\}}
z: \texttt{NX\_NUMBER \{units=NX\_LENGTH\}}

While z is allowed by the canSAS standard, it may not make sense to use it in some situations. Use of z may be ignored by processing software.

(source): \texttt{NXsource}

Description of the radiation source.

![Diagram](image)

Figure 3.15: The SASsource element

@canSAS_class: \texttt{NX\_CHAR}

Official canSAS group: NXcanSAS (contributed definition); SAS-source

Obligatory value: SASsource

radiation: \texttt{NX\_CHAR}
Name of the radiation used. Note that this is not the name of the facility!

Any of these values:

- Spallation Neutron Source
- Pulsed Reactor Neutron Source
- Reactor Neutron Source
- Synchrotron X-ray Source
- Pulsed Muon Source
- Rotating Anode X-ray
- Fixed Tube X-ray
- UV Laser
- Free-Electron Laser
- Optical Laser
- Ion Source
- UV Plasma Source
- neutron
- x-ray
- muon
- electron
- ultraviolet
- visible light
- positron
- proton

**beam_shape**: **NX_CHAR**

Text description of the shape of the beam (incident on the sample).

**wavelength**: **NX_NUMBER** \{units=NX_WAVELENGTH\}

wavelength ($\lambda$) of radiation incident on the sample

**wavelength_min**: **NX_NUMBER** \{units=NX_WAVELENGTH\}

Some facilities specify wavelength using a range. This is the lowest wavelength in such a range.

**wavelength_max**: **NX_NUMBER** \{units=NX_WAVELENGTH\}

Some facilities specify wavelength using a range. This is the highest wavelength in such a range.

**wavelength_spread**: **NX_NUMBER** \{units=NX_WAVELENGTH\}

Some facilities specify wavelength using a range. This is the width of such a range.

**beam_size**: **NXparameters**
Physical dimension of the beam (incident on the sample). If beam is round, just use $x$.

@canSAS_class: NX_CHAR

Ad hoc canSAS_class.

Obligatory value: container

x: NX_NUMBER \{units=NX_LENGTH\}
y: NX_NUMBER \{units=NX_LENGTH\}
z: NX_NUMBER \{units=NX_LENGTH\}

While $z$ is allowed by the canSAS standard, it may not make sense to use it in some situations. Use of $z$ may be ignored by processing software.

(sample): NXsample

Description of the sample.

@canSAS_class: NX_CHAR

Official canSAS group: NXcanSAS (contributed definition); SASsample

Obligatory value: SASsample

ID: NX_CHAR

Text string that identifies this sample.

thickness: NX_FLOAT \{units=NX_LENGTH\}

Thickness of this sample

transmission: NX_NUMBER \{units=NX_DIMENSIONLESS\}

Transmission ($I/I_0$) of this sample. Note that there is no units attribute as this number is dimensionless.

temperature: NX_NUMBER \{units=NX_TEMPERATURE\}

Temperature of this sample.

details: NX_CHAR

Any additional sample details.

position: NXtranslation

Location of the sample in $x$, $y$, and $z$.

orientation: NXorientation

Orientation (rotation) of the sample.

(process): NXprocess

Description of a processing or analysis step.

Add additional fields as needed to describe value(s) of any variable, parameter, or term related to the SASprocess step. Be sure to include units attributes for all numerical fields.

@canSAS_class: NX_CHAR
Figure 3.16: The SASsample element
Figure 3.17: The SASprocess element
Official canSAS group: NXcanSAS (contributed definition); SASprocess

Obligatory value: SASprocess

name: _NX_CHAR_

Optional name for this data processing or analysis step

date: _NX_DATE_TIME_

Optional date for this data processing or analysis step. ¹

description: _NX_CHAR_

Optional description for this data processing or analysis step

term: _NX_CHAR_

Specifies the value of a single variable, parameter, or term (while defined here as a string, it could be a number) related to the SASprocess step.

The name term is not required, it could take any name.

(collection): _NXcollection_

Describes anything about SASprocess that is not already described.

Any content not defined in the canSAS standard can be placed at this point.

@canSAS_class: _NX_CHAR_

Official canSAS group: NXcanSAS (contributed definition); SASprocessnote

Obligatory value: SASprocessnote

(collection): _NXcollection_

Free form description of anything not covered by other elements.

@canSAS_class: _NX_CHAR_

Official canSAS group: NXcanSAS (contributed definition); SASnote

Obligatory value: SASnote

(data): _NXdata_

This describes certain data obtained from a variable-wavelength source such as pulsed-neutron source.

@canSAS_class: _NX_CHAR_

Official canSAS group: NXcanSAS (contributed definition); SAStransmission_spectrum

Obligatory value: SAStransmission_spectrum

@name: _NX_CHAR_

Identify what type of spectrum is being described. It is expected that this value will take either of these two values:

¹ ISO-8601 standard time representation. Use a format for the date which is machine-readable such as ISO-8601 (e.g., yyyy-mm-ddThh:mm:ss) or modified ISO-8601 (e.g., yyyy-mm-dd hh:mm:ss). See: http://www.w3.org/TR/NOTE-datetime or http://en.wikipedia.org/wiki/ISO_8601 for more details.
### SASTransmission_spectrum Element

<table>
<thead>
<tr>
<th>value</th>
<th>meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td>sample</td>
<td>measurement with the sample and container</td>
</tr>
<tr>
<td>can</td>
<td>measurement with just the container</td>
</tr>
</tbody>
</table>

**@timestamp**: `NX_DATE_TIME`  
ISO-8601 time

**lambda**: `NX_NUMBER {units=NX_WAVELENGTH}`  
Wavelength of the radiation.

**T**: `NX_NUMBER {units=NX_DIMENSIONLESS}`  
Transmission value ($I/I_0$)

**@signal**: `NX_CHAR`  
Obligatory value:
- 1: the default data to plot in this group

**@axes**: `NX_CHAR`  
Obligatory value:
- T: the wavelengths field corresponding to this transmission

**@uncertainty**: `NX_CHAR`  
Estimate of the uncertainty of each “math:T.”

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXcanSAS.nxdl.xml

### 3.5.4 NXcite

**Status**:  
contributed definition, extends `NXobject`, version 1.0
**Description:**

Definition to include references for example for detectors, manuals, instruments, acquisition or analysis software used.

The idea would be to include this in the relevant NeXus object: NXdetector for detectors, NXinstrument for instruments, etc.

**Symbols:**

No symbol table

**Groups cited:** none

**Structure:**

- **description**: `NX_CHAR`
  
  This should describe the reason for including this reference. For example: The dataset in this group was normalised using the method which is described in detail in this reference.

- **doi**: `NX_CHAR`
  
  DOI referencing the document or data.

- **endnote**: `NX_CHAR`
  
  Bibliographic reference data in EndNote format.

- **bibtex**: `NX_CHAR`
  
  Bibliographic reference data in BibTeX format.

**Source:** Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXcite.nxdl.xml

### 3.5.5 **NXelectrostatic_kicker**

**Status:**

contributed definition, extends `NXobject`, version 1.0

**Description:**

definition for a electrostatic kicker.

**Symbols:**

No symbol table

**Groups cited:** `NXlog`

**Structure:**

- **description**: `NX_CHAR`
  
  extended description of the kicker.

- **beamline_distance**: `NX_FLOAT` (`units=NX_LENGTH`)
  
  define position of beamline element relative to production target

- ** timing**: `NX_FLOAT` (`units=NX_TIME`)
  
  kicker timing as defined by `description` attribute

  @ **description**: `NX_CHAR`

- **set_current**: `NX_FLOAT` (`units=NX_CURRENT`)

---

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current set on supply.

**set_voltage**: `NX_FLOAT` [units=`NX_VOLTAGE`]

Voltage set on supply.

**read_current**: `NXlog`

Current read from supply.

**value**: `NX_CHAR` [units=`NX_CURRENT`]

**read_voltage**: `NXlog`

Voltage read from supply.

**value**: `NX_CHAR` [units=`NX_VOLTAGE`]

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXelectrostatic_kicker.nxdl.xml

### 3.5.6 NXfresnel_zone_plate

**Status:**

contributed definition, extends `NXobject`, version 1.0

**Description:**

description for a fresnel zone plate

**Symbols:**

No symbol table

**Groups cited:** `NXgeometry`

**Structure:**

**focus_parameters**: `NX_FLOAT`

List of polynomial coefficients describing the focal length of the zone plate, in increasing order.

**outer_diameter**: `NX_FLOAT` [units=`NX_LENGTH`]

**outermost_zone_width**: `NX_FLOAT` [units=`NX_LENGTH`]

**central_stop_diameter**: `NX_FLOAT` [units=`NX_LENGTH`]

**fabrication**: `NX_CHAR`

How zone plate was manufactured

Any of these values: `etched` | `plated` | `zone doubled`

**zone_height**: `NX_FLOAT` [units=`NX_LENGTH`]

**zone_material**: `NX_CHAR`

Material of the zones themselves

**zone_support_material**: `NX_CHAR`

Material present between the zones. This is usually only present for the “zone doubled” fabrication process
3.5.7 NXmagnetic_kicker

Status:
contributed definition, extends NXobject, version 1.0

Description:
definition for a magnetic kicker.

Symbols:
No symbol table

Groups cited: NXlog

Structure:

description: NX_CHAR
extended description of the kicker.

beamline_distance: NX_FLOAT {units=NX_LENGTH}
define position of beamline element relative to production target

timing: NX_FLOAT {units=NX_TIME}
kicker timing as defined by description attribute

@description: NX_CHAR

set_current: NX_FLOAT {units=NX_CURRENT}
current set on supply.

set_voltage: NX_FLOAT {units=NX_VOLTAGE}
voltage set on supply.

read_current: NXlog
current read from supply.

value: NX_CHAR {units=NX_CURRENT}

read_voltage: NXlog
voltage read from supply.

**value:** `NX_CHAR` `{units=NX_VOLTAGE}`

**Source:** Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXmagnetic_kicker.nxdl.xml

### 3.5.8 NXmx

**Status:**

contributed definition, extends `NXobject`, version 1.2

**Description:**

functional application definition for macromolecular crystallography

**Symbols:**

These symbols will be used below to coordinate datasets with the same shape.

- **np**: number of scan points
- **i**: number of detector pixels in the slow direction
- **j**: number of detector pixels in the fast direction

**Groups cited:** `NXattenuator, NXbeam, NXcollection, NXdata, NXdetector_module, NXdetector, NXentry, NXinstrument, NXsample, NXtransformations`

**Structure:**

(entry): `NXentry`

- **title**: `NX_CHAR`
- **start_time**: `NX_DATE_TIME`
- **end_time**: `NX_DATE_TIME`
- **definition**: `NX_CHAR`

NeXus NXDL schema to which this file conforms

Obligatory value: `NXmx`

(instrument): `NXinstrument`

(attenuator): `NXattenuator`

- **attenuator_transmission**: `NX_NUMBER` `{units=NX_UNITLESS}`

(detector): `NXdetector`

- **depends_on**: `NX_CHAR`
- **data[np, i, j]**: `NX_NUMBER`
- **description**: `NX_CHAR`

  name/manufacturer/model/etc. information

- **time_per_channel**: `NX_CHAR` `{units=NX_TIME}`

  todo: define more clearly

- **distance**: `NX_FLOAT` `{units=NX_LENGTH}`
Distance from the sample to the beam center. This value is a guidance only, the proper geometry can be found following the depends_on axis chain.

**dead_time**: `NX_FLOAT` {units=NX\_TIME}
Detector dead time

**count_time**: `NX\_NUMBER` {units=NX\_TIME}
Elapsed actual counting time

**beam_center\_x**: `NX\_FLOAT` {units=NX\_LENGTH}
This is the x position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

**beam_center\_y**: `NX\_FLOAT` {units=NX\_LENGTH}
This is the y position where the direct beam would hit the detector. This is a length, not a pixel position, and can be outside of the actual detector.

**angular\_calibration\_applied**: `NX\_BOOLEAN`
True when the angular calibration has been applied in the electronics, false otherwise.

**angular\_calibration[i, j]**: `NX\_FLOAT`
Angular calibration data.

**flatfield\_applied**: `NX\_BOOLEAN`
True when the flat field correction has been applied in the electronics, false otherwise.

**flatfield[i, j]**: `NX\_FLOAT`
Flat field correction data.

**flatfield\_error[i, j]**: `NX\_FLOAT`
Errors of the flat field correction data.

**pixel\_mask\_applied**: `NX\_BOOLEAN`
True when the pixel mask correction has been applied in the electronics, false otherwise.

**pixel\_mask[i, j]**: `NX\_INT`
The 32-bit pixel mask for the detector. Contains a bit field for each pixel to signal dead, blind or high or otherwise unwanted or undesirable pixels. They have the following meaning:

- bit 0: gap (pixel with no sensor)
- bit 1: dead
- bit 2: under responding
- bit 3: over responding
- bit 4: noisy
- bit 5: undefined
• bit 6: pixel is part of a cluster of problematic pixels (bit set in addition to others)
• bit 7: -undefined-
• bit 8: user defined mask (e.g. around beamstop)
• bits 9-30: -undefined-
• bit 31: virtual pixel (corner pixel with interpolated value)

Normal data analysis software would not take pixels into account when a bit in (mask &amp; 0x0000FFFF) is set. Tag bit in the upper two bytes would indicate special pixel properties that normally would not be a sole reason to reject the intensity value (unless lower bits are set.

countrate_correction__applied: NX_BOOLEAN

True when a count-rate correction has already been applied in the data recorded here, false otherwise.

bit_depth_readout: NX_INT

How many bits the electronics record per pixel.

detector_readout_time: NX_FLOAT {units=NX_TIME}

Time it takes to read the detector (typically milliseconds). This is important to know for time resolved experiments.

frame_time: NX_FLOAT {units=NX_TIME}

This is time for each frame. This is exposure_time + readout time.

gain_setting: NX_CHAR

The gain setting of the detector. This influences background.

saturation_value: NX_INT

The value at which the detector goes into saturation. Data above this value is known to be invalid.

sensor_material: NX_CHAR

At times, radiation is not directly sensed by the detector. Rather, the detector might sense the output from some converter like a scintillator. This is the name of this converter material.

sensor_thickness: NX_FLOAT {units=NX_LENGTH}

At times, radiation is not directly sensed by the detector. Rather, the detector might sense the output from some converter like a scintillator. This is the thickness of this converter material.

threshold_energy: NX_FLOAT {units=NX_ENERGY}

Single photon counter detectors can be adjusted for a certain energy range in which they work optimally. This is the energy setting for this.

type: NX_CHAR

Description of type such as scintillator, ccd, pixel, image plate, CMOS, ...

(transformations): NXtransformations
Suggested location for axes (transformations) to do with the detector

(collection): NXcollection

Suggested container for detailed non-standard detector information like corrections applied automatically or performance settings.

(detector_module): NXdetector_module

This is the description of a detector module. Many detectors consist of multiple smaller modules. Sometimes it is important to know the exact position of such modules. This is the purpose of this group. It is a child group to NXdetector.

data_origin: NX_INT

A two value field which gives the index of the start of the modules data in the main area detector image in the underlying NXdetector module.

data_size: NX_INT

Two values for the size of the module in pixels in each direction.

module_offset: NX_NUMBER {units=NX_LENGTH}

Offset of the module in regards to the origin of the detector in an arbitrary direction.

@transformation_type: NX_CHAR

Obligatory value: translation

@vector: NX_CHAR

@offset: NX_CHAR

@depends_on: NX_CHAR

fast_pixel_direction: NX_NUMBER {units=NX_LENGTH}

Values along the direction of fastest varying pixel direction. The direction itself is given through the vector attribute.

@transformation_type: NX_CHAR

Obligatory value: translation

@vector: NX_CHAR

@offset: NX_CHAR

@depends_on: NX_CHAR

slow_pixel_direction: NX_NUMBER {units=NX_LENGTH}

Values along the direction of slow varying pixel direction. The direction itself is given through the vector attribute.

@transformation_type: NX_CHAR

Obligatory value: translation

@vector: NX_CHAR

@offset: NX_CHAR

@depends_on: NX_CHAR
fast_pixel_size: NX_NUMBER [units=NX_LENGTH]
Values along the direction of fastest varying pixel direction. The
direction itself is given through the vector attribute
   @transformation_type: NX_CHAR
       Obligatory value: translation
   @vector: NX_CHAR
   @offset: NX_CHAR
   @depends_on: NX_CHAR
slow_pixel_size: NX_NUMBER [units=NX_LENGTH]
Values along the direction of slow varying pixel direction. The
direction itself is given through the vector attribute
   @transformation_type: NX_CHAR
       Obligatory value: translation
   @vector: NX_CHAR
   @offset: NX_CHAR
   @depends_on: NX_CHAR

(sample): NXsample
   name: NX_CHAR
       Descriptive name of sample
   depends_on: NX_CHAR
       This should be an absolute requirement to have for any scan experiment. The
       reason it is optional is mainly to accommodate XFEL single shot exposures.
   temperature: NX_CHAR [units=NX_TEMPERATURE]
(beam): NXbeam
   incident_wavelength: NX_NUMBER [units=NX_WAVELENGTH]
   flux: NX_FLOAT [units=NX_FLUX]
       flux incident on beam plane area
   incident_polarisation_stokes[np, 4]: NX_CHAR
   incident_wavelength_spectrum: NXdata
(transformations): NXtransformations
       Suggested location for sample goniometer or other axes (transformations)
(data): NXdata

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXmx.nxdl.xml

3.5.9 NXquadrupole_magnet

Status:
   contributed definition, extends NXobject, version 1.0
Description:

definition for a quadrupole magnet.

Symbols:

No symbol table

Groups cited: NXlog

Structure:

- description: NX_CHAR
  extended description of the magnet.

- beamline_distance: NX_FLOAT {units=NX_LENGTH}
  define position of beamline element relative to production target

- set_current: NX_FLOAT {units=NX_CURRENT}
  current set on supply.

- read_current: NXlog
  current read from supply.
  value: NX_CHAR {units=NX_CURRENT}

- read_voltage: NXlog
  voltage read from supply.
  value: NX_CHAR {units=NX_VOLTAGE}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXquadruple_magnet.nxdl.xml

3.5.10 NXseparator

Status:

contributed definition, extends NXobject, version 1.0

Description:

definition for an electrostatic separator.

Symbols:

No symbol table

Groups cited: NXlog

Structure:

- description: NX_CHAR
  extended description of the separator.

- beamline_distance: NX_FLOAT {units=NX_LENGTH}
  define position of beamline element relative to production target

- set_Bfield_current: NX_FLOAT {units=NX_CURRENT}
  current set on magnet supply.

- set_Efield_voltage: NX_FLOAT {units=NX_VOLTAGE}
current set on HT supply.

**read_Bfield_current**: `NXlog`

- **value**: `NX_CHAR {units=NX_CURRENT}`

**read_Bfield_voltage**: `NXlog`

- **value**: `NX_CHAR {units=NX_VOLTAGE}`

**read_Efield_current**: `NXlog`

- **value**: `NX_CHAR {units=NX_CURRENT}`

**read_Efield_voltage**: `NXlog`

- **value**: `NX_CHAR {units=NX_VOLTAGE}`

**Source**: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXseparator.nxdl.xml

### 3.5.11 NXsnsevent

**Status:**

- contributed definition, extends `NXobject`, version 1.0

**Description:**

This is a definition for event data from Spallation Neutron Source (SNS) at ORNL.

**Symbols:**

- No symbol table

**Groups cited:** `NXaperture`, `NXattenuator`, `NXcollection`, `NXcrystal`, `NXdata`, `NXdetector`, `NXdisk_chopper`, `NXentry`, `NXevent_data`, `NXgeometry`, `NXinstrument`, `NXlog`, `NXmoderator`, `NXmonitor`, `NXnote`, `NXorientation`, `NXpolarizer`, `NXpositioner`, `NXsample`, `NXshape`, `NXsource`, `NXtranslation`, `NXuser`

**Structure:**

- `(entry)`: `NXentry`
  
  - **collection_identifier**: `NX_CHAR`
  
  - **collection_title**: `NX_CHAR`
  
  - **definition**: `NX_CHAR`
    
    Official NXDL schema after this file goes to applications.

    Obligatory value: `NXsnsevent`

  - **duration**: `NX_FLOAT {units=NX_TIME}`
  
  - **end_time**: `NX_DATE_TIME`
  
  - **entry_identifier**: `NX_CHAR`
  
  - **experiment_identifier**: `NX_CHAR`
  
  - **notes**: `NX_CHAR`
proton_charge:  \texttt{NX\_FLOAT} \{units=\texttt{NX\_CHARGE}\}

raw_frames:  \texttt{NX\_INT}

run_number:  \texttt{NX\_CHAR}

start_time:  \texttt{NX\_DATE\_TIME}

title:  \texttt{NX\_CHAR}

total_counts:  \texttt{NX\_UINT} \{units=\texttt{NX\_UNITLESS}\}

total_uncounted_counts:  \texttt{NX\_UINT} \{units=\texttt{NX\_UNITLESS}\}

DASlogs:  \texttt{NX\_collection}

Details of all logs, both from cvinfo file and from HistoTool (frequency and proton_charge).

(log):  \texttt{NX\_log}

average_value:  \texttt{NX\_FLOAT}

average_value_error:  \texttt{NX\_FLOAT}

description:  \texttt{NX\_CHAR}

duration:  \texttt{NX\_FLOAT}

maximum_value:  \texttt{NX\_FLOAT}

minimum_value:  \texttt{NX\_FLOAT}

time[nvalue]:  \texttt{NX\_FLOAT}

value[nvalue]:  \texttt{NX\_FLOAT}

(positioner):  \texttt{NX\_positioner}

Motor logs from cvinfo file.

average_value:  \texttt{NX\_FLOAT}

average_value_error:  \texttt{NX\_FLOAT}

description:  \texttt{NX\_CHAR}

duration:  \texttt{NX\_FLOAT}

maximum_value:  \texttt{NX\_FLOAT}

minimum_value:  \texttt{NX\_FLOAT}

time[numvalue]:  \texttt{NX\_FLOAT}

value[numvalue]:  \texttt{NX\_FLOAT}

SNSHistoTool:  \texttt{NX\_note}

SNSbanking_file_name:  \texttt{NX\_CHAR}

SNSmapping_file_name:  \texttt{NX\_CHAR}

author:  \texttt{NX\_CHAR}

command1:  \texttt{NX\_CHAR}

Command string for event2nxl.
date: \textit{NX_CHAR}
description: \textit{NX_CHAR}
version: \textit{NX_CHAR}

(data): \textit{NXdata}

\texttt{data\_x\_y} \rightarrow /NXentry/NXinstrument/NXdetector/data\_x\_y
\texttt{x\_pixel\_offset} \rightarrow /NXentry/NXinstrument/NXdetector/x\_pixel\_offset
\texttt{y\_pixel\_offset} \rightarrow /NXentry/NXinstrument/NXdetector/y\_pixel\_offset

(event data): \textit{NXevent\_data}

\texttt{event\_index} \rightarrow /NXentry/NXinstrument/NXdetector/event\_index
\texttt{event\_pixel\_id} \rightarrow /NXentry/NXinstrument/NXdetector/event\_pixel\_id
\texttt{event\_time\_of\_flight} \rightarrow /NXentry/NXinstrument/NXdetector/event\_time\_of\_flight
\texttt{pulse\_time} \rightarrow /NXentry/NXinstrument/NXdetector/pulse\_time

instrument: \textit{NXinstrument}

\texttt{SNSdetector\_calibration\_id}: \textit{NX_CHAR}
Detector calibration id from DAS.

\texttt{SNSgeometry\_file\_name}: \textit{NX_CHAR}

\texttt{SNSTranslation\_service}: \textit{NX_CHAR}

beamline: \textit{NX_CHAR}

name: \textit{NX_CHAR}

SNS: \textit{NXsource}

\texttt{frequency}: \textit{NX_FLOAT} \{units=\textit{NX_FREQUENCY}\}

name: \textit{NX_CHAR}

probe: \textit{NX_CHAR}

type: \textit{NX_CHAR}

(detector): \textit{NXdetector}

\texttt{azimuthal\_angle[\text{numx, numy}]}: \textit{NX_FLOAT} \{units=\textit{NX\_ANGLE}\}

\texttt{data\_x\_y[\text{numx, numy}]}: \textit{NX_UINT}

\texttt{distance[\text{numx, numy}]}: \textit{NX_FLOAT} \{units=\textit{NX\_LENGTH}\}

\texttt{event\_index[\text{numpulses}]}: \textit{NX_UINT}

\texttt{event\_pixel\_id[\text{numevents}]}: \textit{NX_UINT}

\texttt{event\_time\_of\_flight[\text{numevents}]}: \textit{NX_FLOAT}

\texttt{pixel\_id[\text{numx, numy}]}: \textit{NX_UINT}

\texttt{polar\_angle[\text{numx, numy}]}: \textit{NX_FLOAT} \{units=\textit{NX\_ANGLE}\}

\texttt{pulse\_time[\text{numpulses}]}: \textit{NX_FLOAT} \{units=\textit{NX\_TIME}\}
total_counts: NX_UINT
x_pixel_offset[numx]: NX_FLOAT \{units=NX_LENGTH\}
y_pixel_offset[numy]:NX_FLOAT \{units=NX_LENGTH\}
origin: NXgeometry
  orientation: NXorientation
  value[6]: NX_FLOAT
    Six out of nine rotation parameters.
shape: NXshape
  description: NX_CHAR
  shape: NX_CHAR
  size[3]: NX_FLOAT \{units=NX_LENGTH\}
translation: NXtranslation
  distance[3]: NX_FLOAT \{units=NX_LENGTH\}
(disk_chopper): NXdisk_chopper
  distance: NX_FLOAT \{units=NX_LENGTH\}
moderator: NXmoderator
  coupling_material: NX_CHAR
  distance: NX_FLOAT \{units=NX_LENGTH\}
  temperature: NX_FLOAT \{units=NX_TEMPERATURE\}
  type: NX_CHAR
(aperture): NXaperture
  x_pixel_offset: NX_FLOAT \{units=NX_LENGTH\}
origin: NXgeometry
  orientation: NXorientation
  value[6]: NX_FLOAT
    Six out of nine rotation parameters.
shape: NXshape
  description: NX_CHAR
  shape: NX_CHAR
  size[3]: NX_FLOAT \{units=NX_LENGTH\}
translation: NXtranslation
  distance[3]: NX_FLOAT \{units=NX_LENGTH\}
(attenuator): NXattenuator
  distance: NX_FLOAT \{units=NX_LENGTH\}
(polarizer): NXpolarizer
(crystal): NXcrystal
type: \texttt{NX\_CHAR}

wavelength: \texttt{NX\_FLOAT} \{units=\texttt{NX\_WAVELENGTH}\}

origin: \texttt{NX\_geometry}

description: \texttt{NX\_CHAR}

orientation: \texttt{NX\_orientation}

value[6]: \texttt{NX\_FLOAT}

Six out of nine rotation parameters.

shape: \texttt{NX\_shape}

description: \texttt{NX\_CHAR}

shape: \texttt{NX\_CHAR}

size: \texttt{NX\_FLOAT} \{units=\texttt{NX\_LENGTH}\}

translation: \texttt{NX\_translation}

distance[3]: \texttt{NX\_FLOAT} \{units=\texttt{NX\_LENGTH}\}

(monitor): \texttt{NX\_monitor}

data[numtimechannels]: \texttt{NX\_UINT}

expect signal=1 axes="time\_of\_flight"

distance: \texttt{NX\_FLOAT} \{units=\texttt{NX\_LENGTH}\}

mode: \texttt{NX\_CHAR}

time\_of\_flight[numtimechannels + 1]: \texttt{NX\_FLOAT} \{units=\texttt{NX\_TIME}\}

sample: \texttt{NX\_sample}

changer\_position: \texttt{NX\_CHAR}

holder: \texttt{NX\_CHAR}

identifier: \texttt{NX\_CHAR}

name: \texttt{NX\_CHAR}

Descriptive name of sample

nature: \texttt{NX\_CHAR}

(user): \texttt{NX\_user}

facility\_user\_id: \texttt{NX\_CHAR}

name: \texttt{NX\_CHAR}

role: \texttt{NX\_CHAR}

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXsnsevent.nxdl.xml

3.5.12 \textbf{NXsnshisto}

Status:

contributed definition, extends \texttt{NXobject}, version 1.0

Description:
This is a definition for histogram data from Spallation Neutron Source (SNS) at ORNL.

Symbols:
No symbol table

Groups cited: NXaperture, NXattenuator, NXcollection, NXcrystal, NXdata, NXdetector, NXdisk_chopper, NXentry, NXfermi_chopper, NXgeometry, NXinstrument, NXlog, NXmoderator, NXmonitor, NXnote, NXorientation, NXpolarizer, NXpositioner, NXsample, NXshape, NXsource, NXtranslation, NXuser

Structure:
(entry): NXentry
  collection_identifier: NX_CHAR
  collection_title: NX_CHAR
  definition: NX_CHAR
    Official NXDL schema after this file goes to applications.
    Obligatory value: NXsnshisto
duration: NX_FLOAT {units=NX_TIME}
end_time: NX_DATE_TIME
entry_identifier: NX_CHAR
experiment_identifier: NX_CHAR
notes: NX_CHAR
proton_charge: NX_FLOAT {units=NX_CHARGE}
raw_frames: NX_INT
run_number: NX_CHAR
start_time: NX_DATE_TIME
title: NX_CHAR
total_counts: NX_UINT {units=NX_UNITLESS}
total_uncounted_counts: NX_UINT {units=NX_UNITLESS}
DASlogs: NXcollection
  Details of all logs, both from cvinfo file and from HistoTool (frequency and proton_charge).
(log): NXlog
  average_value: NX_FLOAT
  average_value_error: NX_FLOAT
description: NX_CHAR
duration: NX_FLOAT
maximum_value: NX_FLOAT
minimum_value: NX_FLOAT
time[nvalue]: NX_FLOAT
value[nvalue]: NX_FLOAT
(positioner): NXpositioner
    Motor logs from cvinfo file.
    average_value: NX_FLOAT
    average_value_error: NX_FLOAT
    description: NX_CHAR
    duration: NX_FLOAT
    maximum_value: NX_FLOAT
    minimum_value: NX_FLOAT
    time[numvalue]: NX_FLOAT
    value[numvalue]: NX_FLOAT

SNSHistoTool: NXnote
    SNSbanking_file_name: NX_CHAR
    SNSmapping_file_name: NX_CHAR
    author: NX_CHAR
    command1: NX_CHAR
        Command string for event2histo_nxl.
    date: NX_CHAR
    description: NX_CHAR
    version: NX_CHAR

(data): NXdata
    data -> /NXentry/NXinstrument/NXdetector/data
    data_x_time_of_flight -> /NXentry/NXinstrument/NXdetector/data_x_time_of_flight
    data_x_y -> /NXentry/NXinstrument/NXdetector/data_x_y
    data_y_time_of_flight -> /NXentry/NXinstrument/NXdetector/data_y_time_of_flight
    pixel_id -> /NXentry/NXinstrument/NXdetector/pixel_id
    time_of_flight -> /NXentry/NXinstrument/NXdetector/time_of_flight
    total_counts -> /NXentry/NXinstrument/NXdetector/total_counts
    x_pixel_offset -> /NXentry/NXinstrument/NXdetector/x_pixel_offset
    y_pixel_offset -> /NXentry/NXinstrument/NXdetector/y_pixel_offset

instrument: NXinstrument
    SNSdetector_calibration_id: NX_CHAR
        Detector calibration id from DAS.
    SNSgeometry_file_name: NX_CHAR
    SNStranslation_service: NX_CHAR
    beamline: NX_CHAR
    name: NX_CHAR
SNS: **NXsource**

- frequency: **NX_FLOAT** \{units=**NX_FREQUENCY**\}
- name: **NX_CHAR**
- probe: **NX_CHAR**
- type: **NX_CHAR**

(detector): **NXdetector**

- azimuthal_angle[numx, numy]: **NX_FLOAT** \{units=**NX_ANGLE**\}
- data[numx, numy, numtof]: **NX_UINT**
- data_x_time_of_flight[numx, numtof]: **NX_UINT**
- data_y[numx, numy]: **NX_UINT**
- data_y_time_of_flight[numy, numtof]: **NX_UINT**
- distance[numx, numy]: **NX_FLOAT** \{units=**NX_LENGTH**\}
- pixel_id[numx, numy]: **NX_UINT**
- polar_angle[numx, numy]: **NX_FLOAT** \{units=**NX_ANGLE**\}
- time_of_flight[numtof + 1]: **NX_FLOAT** \{units=**NX_TIME_OF_FLIGHT**\}
- total_counts: **NX_UINT**
- x_pixel_offset[numx]: **NX_FLOAT** \{units=**NX_LENGTH**\}
- y_pixel_offset[numy]: **NX_FLOAT** \{units=**NX_LENGTH**\}
- origin: **NXgeometry**
  - orientation: **NXorientation**
    - value[6]: **NX_FLOAT**
      - Six out of nine rotation parameters.
  - shape: **NXshape**
    - description: **NX_CHAR**
    - shape: **NX_CHAR**
    - size[3]: **NX_FLOAT** \{units=**NX_LENGTH**\}
  - translation: **NXtranslation**
    - distance[3]: **NX_FLOAT** \{units=**NX_LENGTH**\}

(disk_chopper): **NXdisk_chopper**

- Original specification called for NXchopper, which is not a valid NeXus base class. Select either NXdisk_chopper or NXfermi_chopper, as appropriate.

(distance): **NX_FLOAT** \{units=**NX_LENGTH**\}

(fermi_chopper): **NXfermi_chopper**

- Original specification called for NXchopper, which is not a valid NeXus base class. Select either NXdisk_chopper or NXfermi_chopper, as appropriate.

(distance): **NX_FLOAT** \{units=**NX_LENGTH**\}

(moderator): **NXmoderator**
coupling_material: NX_CHAR
distance: NX_FLOAT {units=NX_LENGTH}
temperature: NX_FLOAT {units=NX_TEMPERATURE}
type: NX_CHAR

(aperture): NXaperture
x_pixel_offset: NX_FLOAT {units=NX_LENGTH}

origin: NXgeometry
orientation: NXorientation
value[6]: NX_FLOAT
Six out of nine rotation parameters.

shape: NXshape
description: NX_CHAR
shape: NX_CHAR

size[3]: NX_FLOAT {units=NX_LENGTH}

translation: NXtranslation
distance[3]: NX_FLOAT {units=NX_LENGTH}

(attenuator): NXattenuator
distance: NX_FLOAT {units=NX_LENGTH}

(polarizer): NXpolarizer

(crystal): NXcrystal
type: NX_CHAR
wavelength: NX_FLOAT {units=NX_WAVELENGTH}

origin: NXgeometry
description: NX_CHAR
orientation: NXorientation
value[6]: NX_FLOAT
Six out of nine rotation parameters.

shape: NXshape
description: NX_CHAR
shape: NX_CHAR

size: NX_FLOAT {units=NX_LENGTH}

translation: NXtranslation
distance[3]: NX_FLOAT {units=NX_LENGTH}

(monitor): NXmonitor
NeXus: a common data format for neutron, x-ray, and muon science, Release 3.1

data[numtimechannels]:  \texttt{NX\_UINT} \\
\textbf{distance}:  \texttt{NX\_FLOAT} \ \{\texttt{units=NX\_LENGTH}\} \\
\textbf{mode}:  \texttt{NX\_CHAR} \\
\textbf{time\_of\_flight[numtimechannels + 1]}:  \texttt{NX\_FLOAT} \ \{\texttt{units=NX\_TIME}\} \\
\textbf{sample}:  \texttt{NX\_sample} \\
\textbf{changer\_position}:  \texttt{NX\_CHAR} \\
\textbf{holder}:  \texttt{NX\_CHAR} \\
\textbf{identifier}:  \texttt{NX\_CHAR} \\
\textbf{name}:  \texttt{NX\_CHAR} \\
\hspace{1cm} \text{Descriptive name of sample} \\
\textbf{nature}:  \texttt{NX\_CHAR} \\
\textbf{(user)}:  \texttt{NX\_user} \\
\textbf{facility\_user\_id}:  \texttt{NX\_CHAR} \\
\textbf{name}:  \texttt{NX\_CHAR} \\
\textbf{role}:  \texttt{NX\_CHAR} \\

\textbf{Source}:  Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXsnshisto.nxdl.xml

3.5.13 \textbf{NXsolenoid\_magnet}

\textbf{Status}: \\
\hspace{1cm} contributed definition, extends \texttt{NX\_object}, version 1.0

\textbf{Description}: \\
\hspace{1cm} definition for a solenoid magnet.

\textbf{Symbols}: 
\hspace{1cm} No symbol table

\textbf{Groups cited}:  \texttt{NX\_log}

\textbf{Structure}: 
\hspace{1cm} \textbf{description}:  \texttt{NX\_CHAR} \\
\hspace{1.5cm} extended description of the magnet. \\
\hspace{1cm} \textbf{beamline\_distance}:  \texttt{NX\_FLOAT} \ \{\texttt{units=NX\_LENGTH}\} \\
\hspace{1.5cm} define position of beamline element relative to production target \\
\hspace{1cm} \textbf{set\_current}:  \texttt{NX\_FLOAT} \ \{\texttt{units=NX\_CURRENT}\} \\
\hspace{1.5cm} current set on supply. \\
\hspace{1cm} \textbf{read\_current}:  \texttt{NX\_log} \\
\hspace{1.5cm} current read from supply. \\
\hspace{1cm} \textbf{value}:  \texttt{NX\_CHAR} \ \{\texttt{units=NX\_CURRENT}\} \\
\hspace{1cm} \textbf{read\_voltage}:  \texttt{NX\_log}

3.5. Contributed Definitions 269
3.5.14 NXspin_rotator

Status:

contributed definition, extends NXobject, version 1.0

Description:

definition for a spin rotator.

Symbols:

No symbol table

Groups cited: NXlog

Structure:

description: NX_CHAR

extended description of the spin rotator.

beamline_distance: NX_FLOAT {units=NX_LENGTH}

define position of beamline element relative to production target

set_Bfield_current: NX_FLOAT {units=NX_CURRENT}

current set on magnet supply.

set_Efield_voltage: NX_FLOAT {units=NX_VOLTAGE}

current set on HT supply.

read_Bfield_current: NXlog

current read from magnet supply.

value: NX_CHAR {units=NX_CURRENT}

read_Bfield_voltage: NXlog

voltage read from magnet supply.

value: NX_CHAR {units=NX_VOLTAGE}

read_Efield_current: NXlog

current read from HT supply.

value: NX_CHAR {units=NX_CURRENT}

read_Efield_voltage: NXlog

voltage read from HT supply.

value: NX_CHAR {units=NX_VOLTAGE}
3.5.15 NXtransformations

Status:
contributed definition, extends NXobject, version 1.0

Description:
Use NXtransformations to gather together any set of movable or fixed elements positioning the device described by the class that contains this.

Symbols:
No symbol table

Groups cited: none

Structure:

anonymous__NEEDS_XSD_CHANGE__: NX_NUMBER
Units need to be appropriate for translation or rotation

@transformation_type: NX_CHAR
Any of these values: translation|rotation

@vector: NX_NUMBER
Three values that define the axis for this transformation

@offset: NX_NUMBER
A fixed offset applied before the transformation (three vector components).

@offset_units: NX_CHAR
Units of the offset.

@depends_on: NX_CHAR
Points to the path of the next element in the geometry chain.

Source: Automatically generated from https://github.com/nexusformat/definitions/blob/master/contributed_definitions/NXtransformations.nxdl.xml
4.1 Status

This application program interface (API) was developed to support the reading and writing of NeXus files through unified function calls, regardless of the physical data format (XML, HDF4, HDF5).

In the meantime it has been decided that active development of NeXus definitions and tools will concentrate on HDF5 as the only supported physical data format. It is expected that most application developers will use standard HDF5 tools to read and write NeXus. Two examples are provided in *Example NeXus C programs using native HDF5 commands*. Therefore, the decision has been taken to freeze the NAPI. Maintenance is reduced to bug fixes.

4.2 Overview

The core routines have been written in C but wrappers are available for a number of other languages including C++, Fortran 77, Fortran 90, Java, Python and IDL. The API makes the reading and writing of NeXus files transparent; the user doesn’t even need to know the underlying format when reading a file since the API calls are the same.

More in-depth and up-to-date information about the NeXus Application Programming Interface for the various language backends is available on-line from http://download.nexusformat.org.

The NeXusIntern.pdf document (http://svn.nexusformat.org/code/trunk/doc/api/NeXusIntern.pdf) describes the internal workings of the NeXus-API. You are very welcome to read it, but it will not be of much use if all you want is to read and write files using the NAPI.

The NeXus Application Program Interface call routines in the appropriate backend (HDF4, HDF5 or XML) to read and write files with the correct structure. The API serves a number of purposes:

1. It simplifies the reading and writing of NeXus files.
2. It ensures a certain degree of compliance with the NeXus standard.
3. It hides the implementation details of the format. In particular, the API can read and write HDF4, HDF5, and XML files using the same routines.
4.3 Core API

The core API provides the basic routines for reading, writing and navigating NeXus files. Operations are performed using a handle that keeps a record of its current position in the file hierarchy. All are read or write requests are then implicitly performed on the currently open entity. This limits number of parameters that need to be passed to API calls, at the cost of forcing a certain mode of operation. It is very similar to navigating a directory hierarchy; NeXus groups are the directories, which can contain data sets and/or other directories.

The core API comprises the following functional groups:

- General initialization and shutdown: opening and closing the file, creating or opening an existing group or dataset, and closing them.
- Reading and writing data and attributes to previously opened datasets.
- Routines to obtain meta-data and to iterate over component datasets and attributes.
- Handling of linking and group hierarchy.
- Routines to handle memory allocation. (Not required in all language bindings.)

4.3.1 NAPI C and C++ Interface

Doxygen documentation is provided online:

- C : http://download.nexusformat.org/doxygen/html-c/
- C++ : http://download.nexusformat.org/doxygen/html-cpp/

4.3.2 NAPI Fortran 77 Interface

Doxygen documentation is provided for the f77 NAPI. (http://download.nexusformat.org/doxygen/html-f77/)

4.3.3 NAPI Fortran 90 Interface

The Fortran 90 interface is a wrapper to the C interface with nearly identical routine definitions. As with the Fortran 77 interface, it is necessary to reverse the order of indices in multidimensional arrays, compared to an equivalent C program, so that data are stored in the same order in the NeXus file.

Any program using the F90 API needs to put the following line at the top (after the PROGRAM statement):

use NXmodule

Use the following table to convert from the C data types listed with each routine to the Fortran 90 data types.

<table>
<thead>
<tr>
<th>C data type</th>
<th>F90 data type</th>
</tr>
</thead>
<tbody>
<tr>
<td>int, int</td>
<td>integer</td>
</tr>
<tr>
<td>char*</td>
<td>character(len=*)</td>
</tr>
<tr>
<td>NXhandle, NXhandle*</td>
<td>type(NXhandle)</td>
</tr>
<tr>
<td>NXstatus</td>
<td>integer</td>
</tr>
<tr>
<td>int[]</td>
<td>integer(:)</td>
</tr>
<tr>
<td>void*</td>
<td>real(:)</td>
</tr>
<tr>
<td>NXlink a, NXlink* a</td>
<td>type(NXlink)</td>
</tr>
</tbody>
</table>

The parameters in the next table, defined in NXmodule, may be used in defining variables.
<table>
<thead>
<tr>
<th>Name</th>
<th>Description</th>
<th>Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>NX_MAXRANK</td>
<td>Maximum number of dimensions</td>
<td>32</td>
</tr>
<tr>
<td>NX_MAXNAMELEN</td>
<td>Maximum length of NeXus name</td>
<td>64</td>
</tr>
<tr>
<td>NXi1</td>
<td>Kind parameter for a 1-byte integer</td>
<td>selected_int_kind(2)</td>
</tr>
<tr>
<td>NXi2</td>
<td>Kind parameter for a 2-byte integer</td>
<td>selected_int_kind(4)</td>
</tr>
<tr>
<td>NXi4</td>
<td>Kind parameter for a 4-byte integer</td>
<td>selected_int_kind(8)</td>
</tr>
<tr>
<td>NXr4</td>
<td>Kind parameter for a 4-byte real</td>
<td>kind(1.0)</td>
</tr>
<tr>
<td>NXr8</td>
<td>Kind parameter for an 8-byte real</td>
<td>kind(1.0D0)</td>
</tr>
</tbody>
</table>

Also see the doxygen documentation. (http://download.nexusformat.org/doxygen/html-f90/)

### 4.3.4 NAPI Java Interface

This section includes installation notes, instructions for running NeXus for Java programs and a brief introduction to the API.

The Java API for NeXus (jnexus) was implemented through the Java Native Interface (JNI) to call on to the native C library. This has a number of disadvantages over using pure Java, however the most popular file backend HDF5 is only available using a JNI wrapper anyway.

**Acknowledgement**

This implementation uses classes and native methods from NCSA’s Java HDF Interface project. Basically all conversions from native types to Java types is done through code from the NCSA HDF group. Without this code the implementation of this API would have taken much longer. See NCSA’s copyright for more information.

**Installation**

**Requirements**

**Caution:** Documentation is old and may need revision.

For running an application with jnexus an recent Java runtime environment (JRE) will do.

In order to compile the Java API for NeXus a Java Development Kit is required on top of the build requirements for the C API.

**Installation under Windows**

1. Copy the HDF DLL’s and the file jnexus.dll to a directory in your path. For instance C:\Windows\system32.
2. Copy the jnexus.jar to the place where you usually keep library jar files.

Note that the location or the naming of these files in the binary Nexus distributions have changed over the years. In the Nexus 4.3.0 Windows 64-bit distribution (http://download.nexusformat.org/kits/4.3.0/win64/), By default, the DLL is at: C:\Program Files\NeXus Data Format\bin\libjnexus-0.dll. Please rename this file to jnexus.dll before making it available in your path. This is important, otherwise, JVM runtime will not be able to locate this file.

For the same distribution, the location of jnexus.jar is at: C:\Program Files\NeXus Data Format\share\Java.
Installation under Unix

The jnexus.so shared library as well as all required file backend .so libraries are required as well as the jnexus.jar file holding the required Java classes. Copy them wherever you like and see below for instructions how to run programs using jnexus.

Running Programs with the NeXus API for Java

In order to successfully run a program with jnexus, the Java runtime systems needs to locate two items:

1. The shared library implementing the native methods.
2. The nexus.jar file in order to find the Java classes.

Locating the shared libraries

The methods for locating a shared library differ between systems. Under Windows32 systems the best method is to copy the jnexus.dll and the HDF4, HDF5 and/or XML-library DLL files into a directory in your path.

On a UNIX system, the problem can be solved in three different ways:

1. Make your system administrator copy the jnexus.so file into the systems default shared library directory (usually /usr/lib or /usr/local/lib).
2. Put the jnexus.so file wherever you see fit and set the LD_LIBRARY_PATH environment variable to point to the directory of your choice.
3. Specify the full pathname of the jnexus shared library on the java command line with the -Dorg.nexusformat.JNEXUSLIB=full-path-2-shared-library option.

Locating jnexus.jar

This is easier, just add the the full pathname to jnexus.jar to the classpath when starting java. Here are examples for a UNIX shell and the Windows shell.

UNIX example shell script to start jnexus.jar

```
#!/sbin/sh
java -classpath /usr/lib/classes.zip:../jnexus.jar:. -Dorg.nexusformat.JNEXUSLIB=../libjnexus.so TestJapi
```

Windows 32 example batch file to start jnexus.jar

```
set JL=-Dorg.nexusformat.JNEXUSLIB=..
nexus\bin\win32\jnexus.dll
java -classpath C:\jre\lib\classes.zip;..\jnexus.jar; %JL% TestJapi
```

Programming with the NeXus API for Java

The NeXus C-API is good enough but for Java a few adaptions of the API have been made in order to match the API better to the idioms used by Java programmers. In order to understand the Java-API, it is useful to study the NeXus
C-API because many methods work in the same way as their C equivalents. A full API documentation is available in Java documentation format. For full reference look especially at:

- The interface NeXusFileInterface first. It gives an uncluttered view of the API.
- The implementation NexusFile which gives more details about constructors and constants. However this documentation is interspersed with information about native methods which should not be called by an application programmer as they are not part of the standard and might change in future.

See the following code example for opening a file, opening a vGroup and closing the file again in order to get a feeling for the API:

```java
fragment for opening and closing
try{
    NexusFile nf = new NexusFile(filename, NexusFile.NXACC_READ);
    nf.opengroup("entry1","NXentry");
    nf.finalize();
}catch(NexusException ne) {
    // Something was wrong!
}
```

Some notes on this little example:

- Each NeXus file is represented by a NexusFile object which is created through the constructor.
- The NexusFile object takes care of all file handles for you. So there is no need to pass in a handle anymore to each method as in the C language API.
- All error handling is done through the Java exception handling mechanism. This saves all the code checking return values in the C language API. Most API functions return void.
- Closing files is tricky. The Java garbage collector is supposed to call the finalize method for each object it decides to delete. In order to enable this mechanism, the NXclose() function was replaced by the finalize() method. In practice it seems not to be guaranteed that the garbage collector calls the finalize() method. It is safer to call finalize() yourself in order to properly close a file. Multiple calls to the finalize() method for the same object are safe and do no harm.

**Data Writing and Reading**

Again a code sample which shows how this looks like:

```java
fragment for writing and reading
int idata[][] = new idata[10][20];
int iDim[] = new int[2];
// put some data into idata.......
// write idata
iDim[0] = 10;
iDim[1] = 20;
nf.makedata("idata",NexusFile.NX_INT32,2,iDim);
nf.opendata("idata");
nf.putdata(idata);
```

4.3. Core API
The dataset is created as usual with `makedata()` and opened with `putdata()`. The trick is in `putdata()`. Java
is meant to be type safe. One would think then that a `putdata()` method would be required for each Java data type.
In order to avoid this, the data to `write()` is passed into `putdata()` as type `Object`. Then the API proceeds to
analyze this object through the Java introspection API and convert the data to a byte stream for writing through the
native method call. This is an elegant solution with one drawback: An array is needed at all times. Even if only a
single data value is written (or read) an array of length one and an appropriate type is the required argument.

Another issue are strings. Strings are first class objects in Java. HDF (and NeXus) sees them as dumb arrays of bytes.
Thus strings have to be converted to and from bytes when reading string data. See a writing example:

```
String writing
String ame = "Alle meine Entchen";
nf.makedata("string_data",NexusFile.NX_CHAR,
1,ame.length()+2);
nf.opendata("string_data");
nf.putdata(ame.getBytes());
```

And reading:

```
String reading
String ame = "Alle meine Entchen";
nf.makedata("string_data",NexusFile.NX_CHAR,
1,ame.length()+2);
nf.opendata("string_data");
nf.putdata(ame.getBytes());
```

The aforementioned holds for all strings written as SDS content or as an attribute. SDS or vGroup names do not need
this treatment.

**Inquiry Routines**

Let us compare the C-API and Java-API signatures of the `getinfo()` routine (C) or method (Java):

**C API signature of `getinfo()`**

```
/* C -API */
NXstatus NXgetinfo(NXhandle handle, int *rank, int iDim[],
int *datatype);
```

**Java API signature of `getinfo()`**

```
// Java
void getinfo(int iDim[], int args[]);
```
The problem is that Java passes arguments only by value, which means they cannot be modified by the method. Only array arguments can be modified. Thus **args** in the `getinfo()` method holds the rank and datatype information passed in separate items in the C-API version. For resolving which one is which, consult a debugger or the API-reference.

The attribute and vGroup search routines have been simplified using HashTables. The **Hashtable** returned by `groupdir()` holds the name of the item as a key and the classname or the string SDS as the stored object for the key. Thus the code for a vGroup search looks like this:

**vGroup search**

```java
1. nf.opengroup(group, nxclass);
2. h = nf.groupdir();
3. e = h.keys();
4. System.out.println("Found in vGroup entry:");
5. while (e.hasMoreElements())
6. {
7.   vname = (String)e.nextElement();
8.   vclass = (String)h.get(vname);
9.   System.out.println("  Item: " + vname + "  class: " + vclass);
10. }
```

For an attribute search both at global or SDS level the returned **Hashtable** will hold the name as the key and a little class holding the type and size information as value. Thus an attribute search looks like this in the Java-API:

**attribute search**

```java
1. Hashtable h = nf.attrdir();
2. Enumeration e = h.keys();
3. while (e.hasMoreElements())
4. {
5.   attname = (String)e.nextElement();
6.   atten = (AttributeEntry)h.get(attname);
7.   System.out.println("Found global attribute: " + attname +
8.   "  type: " + atten.type + ",length: " + atten.length);
9. }
```

For more information about the usage of the API routines see the reference or the NeXus C-API reference pages. Another good source of information is the source code of the test program which exercises each API routine.

**Known Problems**

These are a couple of known problems which you might run into:

**Memory** As the Java API for NeXus has to convert between native and Java number types a copy of the data must be made in the process. This means that if you want to read or write 200MB of data your memory requirement will be 400MB! This can be reduced by using multiple `getslab()`/`putslab()` to perform data transfers in smaller chunks.

**Java.lang.OutOfMemoryException** By default the Java runtime has a low default value for the maximum amount of memory it will use. This ceiling can be increased through the `-mxXXm` option to the Java runtime. An example: `java -mx512m ...` starts the Java runtime with a memory ceiling of 512MB.
Maximum 8192 files open  The NeXus API for Java has a fixed buffer for file handles which allows only 8192 NeXus files to be open at the same time. If you ever hit this limit, increase the MAXHANDLE define in native/handle.h and recompile everything.

On-line Documentation

The following documentation is browsable online:

1. The Doxygen API documentation
2. A verbose tutorial for the NeXus for Java API.
3. The API Reference.
4. Finally, the source code for the test driver for the API which also serves as a documented usage example.

4.3.5 NAPI Python Interface

Documentation available in pydoc and doxygen. (http://download.nexusformat.org/doxygen/html-python)

4.3.6 NAPI IDL Interface

IDL is an interactive data evaluation environment developed by Research Systems - it is an interpreted language for data manipulation and visualization. The NeXus IDL bindings allow access to the NeXus API from within IDL - they are installed when NeXus is compiled from source after being configured with the following options:

```bash
configure \
   --with-idlroot=/path/to/idl/installation \
   --with-idldlm=/path/to/install/dlm/files/to
```

For further details see the README (http://htmlpreview.github.com/?https://github.com/nexusformat/code/blob/master/bindings/idl/README.html) for the NeXus IDL binding.

4.4 Utility API

The NeXus F90 Utility API provides a number of routines that combine the operations of various core API routines in order to simplify the reading and writing of NeXus files. At present, they are only available as a Fortran 90 module but a C version is in preparation.

The utility API comprises the following functional groups:

- Routines to read or write data.
- Routines to find whether or not groups, data, or attributes exist, and to find data with specific signal or axis attributes, i.e. to identify valid data or axes.
- Routines to open other groups to which NXdata items are linked, and to return again.

1 http://download.nexusformat.org/doxygen/html-java/
line required for use with F90 API

Any program using the F90 Utility API needs to put the following line near the top of the program:

```fortran
use NXUmodule
```

**Note:** Do not put `USE` statements for both `NXmodule` and `NXUmodule`. The former is included in the latter.

### 4.4.1 List of F90 Utility Routines

<table>
<thead>
<tr>
<th>name</th>
<th>description</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>Reading and Writing</strong></td>
<td></td>
</tr>
<tr>
<td>NXUwriteglobals</td>
<td>Writes all the valid global attributes of a file.</td>
</tr>
<tr>
<td>NXUwritegroup</td>
<td>Opens a group (creating it if necessary).</td>
</tr>
<tr>
<td>NXUwritedata</td>
<td>Opens a data item (creating it if necessary) and writes data and its units.</td>
</tr>
<tr>
<td>NXUreaddata</td>
<td>Opens and reads a data item and its units.</td>
</tr>
<tr>
<td>NXUwritehistogram</td>
<td>Opens one dimensional data item (creating it if necessary) and writes histogram centers and their units.</td>
</tr>
<tr>
<td>NXUreadhistogram</td>
<td>Opens and reads a one dimensional data item and converts it to histogram bin boundaries.</td>
</tr>
<tr>
<td>NXUsetcompress</td>
<td>Defines the compression algorithm and minimum dataset size for subsequent write operations.</td>
</tr>
<tr>
<td><strong>Finding Groups, Data, and Attributes</strong></td>
<td></td>
</tr>
<tr>
<td>NXUfindclass</td>
<td>Returns the name of a group of the specified class if it is contained within the currently open group.</td>
</tr>
<tr>
<td>NXUfinddata</td>
<td>Checks whether a data item of the specified name is contained within the currently open group.</td>
</tr>
<tr>
<td>NXUfindattr</td>
<td>Checks whether the currently open data item has the specified attribute.</td>
</tr>
<tr>
<td>NXUfindsignal</td>
<td>Searches the currently open group for a data item with the specified SIGNAL attribute.</td>
</tr>
<tr>
<td>NXUfindaxis</td>
<td>Searches the currently open group for a data item with the specified AXIS attribute.</td>
</tr>
<tr>
<td><strong>Finding Linked Groups</strong></td>
<td></td>
</tr>
<tr>
<td>NXUfindlink</td>
<td>Finds another link to the specified NeXus data item and opens the group it is in.</td>
</tr>
<tr>
<td>NXUresume link</td>
<td>Reopens the original group from which <code>NXUfindlink</code> was used.</td>
</tr>
</tbody>
</table>

Currently, the F90 utility API will only write character strings, 4-byte integers and reals, and 8-byte reals. It can read other integer sizes into four-byte integers, but does not differentiate between signed and unsigned integers.

### 4.5 Building Programs

The install kit provides a utility call `nxbuild` that can be used to build simple programs:

```bash
nxbuild -o test test.c
```

This script links in the various libraries for you and reading its contents would provide the necessary information for creating a separate Makefile. You can also use `nxbuild` with the example files in the NeXus distribution kit which are installed into `/usr/local/nexus/examples`

Note that the executable name is important in this case as the test program uses it internally to determine the `NXACC_CREATE*` argument to pass to `NXopen`.  

---

**4.5. Building Programs** 281
building and running a simple NeXus program

# builds HDF5 specific test
nxbuild -o napi_test-hdf5 napi_test.c

# runs the test
./napi_test-hdf5

NeXus is also set up for pkg-config so the build can be done as:
gcc 'pkg-config --cflags' 'pkg-config --libs' -o test test.c

4.6 Reporting Bugs in the NeXus API

If you encounter any bugs in the installation or running of the NeXus API, please report them online using our Issue Reporting system. (http://www.nexusformat.org/IssueReporting)
NeXus began as a group of scientists with the goal of defining a common data storage format to exchange experimental results and to exchange ideas about how to analyze them.

The NeXus Scientific Community provides the scientific data, advice, and continued involvement with the NeXus standard. NeXus provides a forum for the scientific community to exchange ideas in data storage through the NeXus wiki.

The NeXus International Advisory Committee (NIAC) supervises the development and maintenance of the NeXus common data format for neutron, X-ray, and muon science. The NIAC supervises a technical committee to oversee the NeXus Application Programmer Interface (NAPI) and the NeXus class definitions.

There are several mechanisms in place in order to coordinate the development of NeXus with the larger community.

5.1 NeXus Wiki

First of all, there is the NeXus wiki, http://wiki.nexusformat.org/, which provides all kinds of information, including membership, minutes, and discussions from the meetings of the NIAC and Technical Committee Code Camps, proposed designs for consideration by NeXus, the NeXus logo, as well as some legacy documentation that we have not quite managed to move into the manual.

5.2 Contributed Definitions

The community is encouraged to provide new definitions (Base Class Definitions or Application Definitions) for consideration in the NeXus standard. These community contributions will be entered in the Contributed Definitions and will be curated according to procedures set forth by the NIAC: The NeXus International Advisory Committee.

5.3 Other Ways NeXus Coordinates with the Scientific Community

5.3.1 NIAC: The NeXus International Advisory Committee

The purpose of the NeXus International Advisory Committee (NIAC) is to supervise the development and maintenance of the NeXus common data format for neutron, X-ray, and muon science. This purpose includes, but is not limited to, the following activities.

\[1\] For more details about the NIAC constitution, procedures, and meetings, refer to the NIAC wiki page: http://wiki.nexusformat.org/NIAC The members of the NIAC may be reached by email: nexus-committee@nexusformat.org
1. To establish policies concerning the definition, use, and promotion of the NeXus format.

2. To ensure that the specification of the NeXus format is sufficiently complete and clear for its use in the exchange and archival of neutron, X-ray, and muon data.

3. To receive and examine all proposed amendments and extensions to the NeXus format. In particular, to ratify proposed instrument and group class definitions, to ensure that the data structures conform to the basic NeXus specification, and to ensure that the definitions of data items are clear and unambiguous and conform to accepted scientific usage.

4. To ensure that documentation of the NeXus format is sufficient, current, and available to potential users both on the internet and in other forms.

5. To coordinate with the developers of the NeXus Application Programming Interface to ensure that it supports the use of the NeXus format in the neutron, X-ray, and muon communities, and to promote other software development that will benefit users of the NeXus format.

6. To coordinate with other organizations that maintain and develop related data formats to ensure maximum compatibility.

The committee will meet at least once every other calendar year according to the following plan:

- In years coinciding with the NOBUGS series of conferences (once every two years), members of the entire NIAC will meet as a satellite meeting to NOBUGS, along with interested members of the community.
- In intervening years, the executive officers of the NIAC will attend, along with interested members of the NIAC. This is intended to be a working meeting with a small group.

### 5.3.2 NeXus Mailing List

We invite anyone who is associated with neutron and/or X-ray synchrotron science and who wishes to be involved in the development and testing of the NeXus format to subscribe to this list. It is for the free discussion of all aspects of the design and operation of the NeXus format.

- **List Address:** nexus@nexusformat.org
- **Subscriptions:** http://lists.nexusformat.org/mailman/listinfo/nexus
- **Archive:** http://lists.nexusformat.org/pipermail/nexus

### 5.3.3 NeXus International Advisory Committee (NIAC) Mailing List

This list contains discussions of the **NIAC: The NeXus International Advisory Committee**, which oversees the development of the NeXus data format. Its members represent many of the major neutron and synchrotron scattering sources in the world. Membership and posting to this list are limited to the committee members, but the archives are public.

- **List Address:** nexus-committee@nexusformat.org
- **Subscriptions:** http://lists.nexusformat.org/mailman/listinfo/nexus-committee
- **Archive:** http://lists.nexusformat.org/pipermail/nexus-committee

### 5.3.4 NeXus Developers Mailing List (retired)

This mailing list was for discussions concerning the technical development of NeXus (the Definitions, NXDL, and the NeXus Application Program Interface). There was, however, much overlap with the general NeXus mailing list and so this separate list was closed in October 2012, but the archive of previous posting is still available.

- **Archive:** http://lists.nexusformat.org/pipermail/nexus-developers
5.3.5 NeXus Repositories

NeXus NXDL class definitions (both base classes and application definitions) and the NeXus code library source are held in a pair of git repositories on GitHub.

The repositories are world readable. You can browse them directly:

- **NeXus code library and applications**  https://github.com/nexusformat/code
- **NeXus NXDL class definitions**  https://github.com/nexusformat/definitions

If you would like to contribute (thank you!), the normal GitHub procedure of forking the repository and generating pull requests should be used.

Please report any problems via the *Issue Reporting* system.

5.3.6 NeXus Issue Reporting

NeXus is using GitHub (https://github.com) as source code repository and for problem reporting. The issue reports (see *View current issues* below) are used to guide the NeXus developers in resolving problems as well as implementing new features.

**NeXus Code (NAPI, Library, and Applications)**

- **View current issues**  https://github.com/nexusformat/code/issues
- **Timeline (recent ticket and code changes)**  https://github.com/nexusformat/code/pulse

**NeXus Definitions (NXDL base classes and application definitions)**

- **View current issues**  https://github.com/nexusformat/definitions/issues
- **Timeline (recent ticket and definition changes)**  https://github.com/nexusformat/definitions/pulse
This section describes how to install the NeXus API and details the requirements. The NeXus API is distributed under the terms of the GNU Lesser General Public License version 3.

The source code and binary versions for some popular platforms can be found on http://download.nexusformat.org/kits/. Up to date instructions can be found on the NeXus Wiki Download page (http://www.nexusformat.org/Download). In case you need help, feel free to contact the NeXus mailing list: http://lists.nexusformat.org/mailman/listinfo/nexus

### 6.1 Precompiled Binary Installation

#### 6.1.1 Prerequisites

**HDF5/HDF4**

*Note:* HDF5 is the preferred format to use for NeXus.

NeXus uses HDF5 as the main underlying binary format. (HDF4 is supported as a legacy underlying binary format but is not recommended for new use.) It is necessary first to install the HDF subroutine libraries and include files before compiling the NeXus API. It is not usually necessary to download the HDF source code since precompiled object libraries exist for a variety of operating systems including Windows, Mac OS X, Linux, and various other flavors of Unix. Check the HDF web pages for more information: http://www.hdfgroup.org/

Packages for HDF4 and HDF5 are available for both Fedora (hdf, hdf5, hdf-devel, hdf5-devel) and Ubuntu/Debian (libhdf4g, libhdf5).

**XML**

*Note:* XML is not the preferred format to use for NeXus.

The NeXus API also supports using XML as a legacy underlying on-disk format. This uses the Mini-XML library, developed by Michael Sweet, which is also available as a precompiled binary library for several operating systems. Check the Mini-XML web pages for more information: http://www=minixml.org/

Packages for MXML are available for both Fedora (mxml, mxml-devel) and Ubuntu/Debian (libxmlml1).
6.1.2 Linux RPM Distribution Kits

An installation kit (source or binary) can be downloaded from: http://download.nexusformat.org/kits/

A NeXus binary RPM (nexus-*.i386.rpm) contains ready compiled NeXus libraries whereas a source RPM (nexus-*.src.rpm) needs to be compiled into a binary RPM before it can be installed. In general, a binary RPM is installed using the command

```
rpm -Uvh file.i386.rpm
```

or, to change installation location from the default (e.g. /usr/local) area, using

```
rpm -Uvh --prefix /alternative/directory file.i386.rpm
```

If the binary RPMS are not the correct architecture for you (e.g. you need x86_64 rather than i386) or the binary RPM requires libraries (e.g. HDF4) that you do not have, you can instead rebuild a source RPM (.src.rpm) to generate the correct binary RPM for your machine. Download the source RPM file and then run

```
rpmbuild --rebuild file.src.rpm
```

This should generate a binary RPM file which you can install as above. Be careful if you think about specifying an alternative buildroot for rpmbuild by using `--buildroot` option as the “buildroot” directory tree will get remove (so `--buildroot /` is a really bad idea). Only change buildroot it if the default area turns out not to be big enough to compile the package.

If you are using Fedora, then you can install all the dependencies by typing

```
yum install hdf hdf-devel hdf5 hdf5-devel mxml mxml-devel
```

6.1.3 Microsoft Windows Installation Kit

A Windows MSI based installation kit is available and can be downloaded from: http://download.nexusformat.org/kits/windows/

6.1.4 Mac OS X Installation Kit

An installation disk image (.dmg) can be downloaded from: http://download.nexusformat.org/kits/macosx/

6.2 Source Installation

6.2.1 NeXus Source Code Distribution

The build uses autoconf (so autools are required) to determine what features will be available by your system. You must have the development libraries installed for all the file backends you want support for (see above). If you intend to build more than the C language bindings, you need to have the respective build support in a place where autoconf will pick them up (i.e. python development files, a Java Development Kit, etc.).

For more information see the README in the toplevel of the source distribution. In case you need help, feel free to contact the NeXus Mailing List:

- **Archives** http://lists.nexusformat.org/mailman/listinfo/nexus
- **email** nexus@nexusformat.org
Download the appropriate gzipped tar file, unpack it, and run the standard configure procedure from the resulting
nexus directory. For example, for version 4.2.1:

```bash
$ tar zxvf nexus-4.2.1.tar.gz
$ cd nexus-4.2.1
$ ./configure
```

To find out how to customize the installation, e.g., to choose different installation directories, type

```bash
$ ./configure --help
```

Carefully check the final output of the configure run. Make sure all features requested are actually enabled.

```bash
$ make
$ make install
```

See the README file for further instructions.

### 6.2.2 Cygwin Kits

HDF4 is not supported under CYGWIN - both HDF5 and MXML are supported and can be downloaded and built as
usual. When configuring HDF5 you should explicitly pass a prefix to the configure script to make sure the libraries
are installed in a “usual” location i.e.

```bash
./configure --prefix=/usr/local/hdf5
```

Otherwise you will have to use the `--with-hdf5=/path/to/hdf5` option later when configuring NeXus to tell
it where to look for hdf5. After building hdf5, configure and build NeXus using the instructions for source code
distribution above.
There are many utilities available to read, browse, write, and use NeXus data files. Some are provided by the NeXus technical group while others are provided by the community. Still, other tools listed here can read or write one of the low-level file formats used by NeXus (HDF4, HDF5, or XML).

### 7.1 Utilities supplied with NeXus

Most of these utility programs are run from the command line. It will be noted if a program provides a graphical user interface (GUI). Short descriptions are provided here with links to further information, as available.

**nxbrowse**  NeXus Browser

**nxconvert**  Utility to convert a NeXus file into HDF4/HDF5/XML/...

**nxdir**  *nxdir* is a utility for querying a NeXus file about its contents. Full documentation can be found by running this command:

```
nxdir -h
```

**nxingest**  *nxingest* extracts the metadata from a NeXus file to create an XML file according to a mapping file. The mapping file defines the structure (names and hierarchy) and content (from either the NeXus file, the mapping file or the current time) of the output file. See the man page for a description of the mapping file. This tool uses the NAPI. Thus, any of the supported formats (HDF4, HDF5 and XML) can be read.

**nxsummary**  Use *nxsummary* to generate summary of a NeXus file. This program relies heavily on a configuration file. Each *item* tag in the file describes a node to print from the NeXus file. The *path* attribute describes where in the NeXus file to get information from. The *label* attribute will be printed when showing the value of the specified field. The optional *operation* attribute provides for certain operations to be performed on the data before printing out the result. See the source code documentation for more details.

**nxtranslate**  *nxtranslate* is an anything to NeXus converter. This is accomplished by using translation files and a plugin style of architecture where *nxtranslate* can read from new formats as plugins become available. The documentation for *nxtranslate* describes its usage by three types of individuals:

- the person using existing translation files to create NeXus files
- the person creating translation files
- the person writing new *retrievers*

All of these concepts are discussed in detail in the documentation provided with the source code.

**nxvalidate**  From the source code documentation:

“Utility to convert a NeXus file into HDF4/HDF5/XML/...”
Note: this command-line tool is different than the newer Java GUI program: NXvalidate.

**NXplot** 
An extendable utility for plotting any NeXus file. **NXplot** is an Eclipse-based GUI project in Java to plot data in NeXus files. (The project was started at the first NeXus Code Camp in 2009.)

### 7.2 Data Analysis

The list of applications below are some of the utilities that have been developed (or modified) to read/write NeXus files as a data format. It is not intended to be a complete list of all available packages.

**DAVE** ([http://www.ncnr.nist.gov/dave/](http://www.ncnr.nist.gov/dave/)) DAVE is an integrated environment for the reduction, visualization and analysis of inelastic neutron scattering data. It is built using IDL (Interactive Data Language) from ITT Visual Information Solutions.

**DAWN** ([http://www.dawnsci.org](http://www.dawnsci.org)) The Data Analysis Workbench (DAWN) project is an eclipse based workbench for doing scientific data analysis. It offers generic visualisation, and domain specific processing.

**GDA** ([http://www.opengda.org](http://www.opengda.org)) The GDA project is an open-source framework for creating customised data acquisition software for science facilities such as neutron and X-ray sources. It has elements of the DAWN analysis workbench built in.

**Gumtree** ([http://docs.codehaus.org/display/GUMTREE](http://docs.codehaus.org/display/GUMTREE)) Gumtree is an open source project, providing a graphical user interface for instrument status and control, data acquisition and data reduction.

**IDL** ([http://www.ittvis.com/](http://www.ittvis.com/)) IDL is a high-level technical computing language and interactive environment for algorithm development, data visualization, data analysis, and numeric computation.

**IgorPro** ([http://www.wavemetrics.com/](http://www.wavemetrics.com/)) IGOR Pro is an extraordinarily powerful and extensible scientific graphing, data analysis, image processing and programming software tool for scientists and engineers.

**ISAW** ([ftp://ftp.sns.gov/ISAW/](ftp://ftp.sns.gov/ISAW/)) The Integrated Spectral Analysis Workbench software project (ISAW) is a Platform-Independent system Data Reduction/Visualization. ISAW can be used to read, manipulate, view, and save neutron scattering data. It reads data from IPNS run files or NeXus files and can merge and sort data from separate measurements.

**LAMP** ([http://www.ill.eu/data_treat/lamp/](http://www.ill.eu/data_treat/lamp/)) LAMP (Large Array Manipulation Program) is designed for the treatment of data obtained from neutron scattering experiments at the Institut Laue-Langevin. However, LAMP is now a more general purpose application which can be seen as a GUI-laboratory for data analysis based on the IDL language.

**Mantid** ([http://www.mantidproject.org/](http://www.mantidproject.org/)) The Mantid project provides a platform that supports high-performance computing on neutron and muon data. It is being developed as a collaboration between Rutherford Appleton Laboratory and Oak Ridge National Laboratory.

**MATLAB** ([http://www.mathworks.com/](http://www.mathworks.com/)) MATLAB is a high-level technical computing language and interactive environment for algorithm development, data visualization, data analysis, and numeric computation.

**NeXpy** ([http://nexpy.github.io/nexpy/](http://nexpy.github.io/nexpy/)) The goal of NeXpy is to provide a simple graphical environment, coupled with Python scripting capabilities, for the analysis of X-Ray and neutron scattering data. (It was decided at the NIAC 2010 meeting that a large portion of this code would be adopted in the future by NeXus and be part of the distribution)

**OpenGENIE** ([http://www.opengenie.org/](http://www.opengenie.org/)) A general purpose data analysis and visualisation package primarily developed at the ISIS Facility, Rutherford Appleton Laboratory.

**PyMCA** ([http://pymca.sourceforge.net/](http://pymca.sourceforge.net/)) PyMca is a ready-to-use, and in many aspects state-of-the-art, set of applications implementing most of the needs of X-ray fluorescence data analysis. It also provides a Python toolkit for visualization and analysis of energy-dispersive X-ray fluorescence data. Reads, browses, and plots data from NeXus HDF5 files.
spec2nexus (http://spec2nexus.readthedocs.org/) (Python code) Converts SPEC data files and scans into NeXus HDF5 files. Provides h5toText utility program to inspect HDF5 file content. Provides libraries:

- *spec2nexus.prjPySpec*: python binding to read SPEC [#] data files
- *spec2nexus.eznx*: (Easy NeXus) supports writing NeXus HDF5 files using h5py

### 7.3 HDF Tools

Here are some of the generic tools that are available to work with HDF files. In addition to the software listed here there are also APIs for many programming languages that will allow low level programmatic access to the data structures.

**HDF Group command line tools** (http://www.hdfgroup.org/products/hdf5_tools/#h5dist/) There are various command line tools that are available from the HDF Group, these are usually shipped with the HDF5 kits but are also available for download separately.

**HDFexplorer** (http://www.space-research.org/) A data visualization program that reads Hierarchical Data Format files (HDF, HDF-EOS and HDF5) and also netCDF data files.

**HDFview** (http://www.hdfgroup.org) A Java based GUI for browsing (and some basic plotting) of HDF files.

#### 7.3.1 Language APIs

**h5py** (http://code.google.com/p/h5py/) HDF5 for Python (h5py) is a general-purpose Python interface to HDF5.
BRIEF HISTORY OF NEXUS

Two things to note about the development and history of NeXus:

- All efforts on NeXus have been voluntary except for one year when we had one full-time worker.
- The NIAC has already discussed many matters related to the format.

**June 1994** Mark Könnecke (then ISIS, now PSI) made a proposal using netCDF for the European neutron scattering community while working at ISIS

**August 1994** Jonathan Tischler (ORNL) proposed an HDF-based format as a standard for data storage at APS

**October 1994** Ray Osborn convened a series of three workshops called SoftNeSS. In the first meeting, Mark Könnecke and Jon Tischler were invited to meet with representatives from all the major U.S. neutron scattering laboratories at Argonne National Laboratory to discuss future software development for the analysis and visualization of neutron data. One of the main recommendations of SoftNeSS’94 was that a common data format should be developed.

**September 1995** At SoftNeSS 1995 (at NIST), three individual data format proposals by Przemek Klosowski (NIST), Mark Könnecke (then ISIS), and Jonathan Tischler (ORNL and APS/ANL) were joined to form the basis of the current NeXus format. At this workshop, the name NeXus was chosen.

**August 1996** The HDF-4 API is quite complex. Thus a NeXus Abstract Programmer Interface NAPI was released which simplified reading and writing NeXus files.

**October 1996** At SoftNeSS 1996 (at ANL), after reviewing the different scientific data formats discussed, it was decided to use HDF4 as it provided the best grouping support. The basic structure of a NeXus file was agreed upon. The various data format proposals were combined into a single document by Przemek Klosowski (NIST), Mark Könnecke (then ISIS), Jonathan Tischler (ORNL and APS/ANL), and Ray Osborn (IPNS/ANL) coauthored the first proposal for the NeXus scientific data standard.

**July 1997** SINQ at PSI started writing NeXus files to store raw data.

**Summer 2001** MLNSC at LANL started writing NeXus files to store raw data

**September 2002** NeXus API version 2.0.0 is released. This version brought support for the new version of HDF; HDF5, released by the HDF group. HDF4 imposed limits on file sizes and the number of objects in a file. These issues were resolved with HDF5. The NeXus API abstracted the difference between the two physical file formats away from the user.

**June 2003** Przemek Klosowski, Ray Osborn, and Richard Riedel received the only known grant explicitly for working on NeXus from the Systems Integration for Manufacturing Applications (SIMA)
program of the National Institute of Standards and Technology (NIST). The grant funded a person for one year to work on community wide infrastructure in NeXus.

**October 2003** In 2003, NeXus had arrived at a stage where informal gatherings of a group of people were no longer good enough to oversee the development of NeXus. This lead to the formation of the NeXus International Advisory Committee (NIAC) which strives to include representatives of all major stake holders in NeXus. A first meeting was held at CalTech. Since 2003, the NIAC meets every year to discuss all matters NeXus.

**July 2005** The community asked the NeXus team to provide an ASCII based physical file format which allows them to edit their scientific results in emacs. This lead to the development of a XML NeXus physical format. This was released with NeXus API version 3.0.0.

**May 2007** NeXus API version 4.0.0 is released with broader support for scripting languages and the feature to link with external files.

**October 2007** NeXus API version 4.1.0 is released with many bug-fixes.

**October 2008** *NXDL: The NeXus Definition Language* is defined. Until now, NeXus used another XML format, meta-DTD, for defining base classes and application definitions. There were several problems with meta-DTD, the biggest one being that it was not easy to validate against it. NXDL was designed to circumvent these problems. All current base classes and application definitions were ported into the NXDL.

**April 2009** NeXus API version 4.2.0 is released with additional C++, IDL, and python/numpy interfaces.

**September 2009** NXDL and draft *NXsas* presented to canSAS at SAS2009 conference

**January 2010** NXDL presented to ESRF HDF5 workshop on hyperspectral data

**May 2012** first release (3.1.0) of NXDL (NeXus Definition Language)
CHAPTER NINE

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9.2 Colophon

These docs (manual and reference) were produced using Sphinx (http://sphinx.pocoo.org). The source for the manual shows many examples of the structures used to create the manual. If you have any questions about how to contribute to this manual, please contact the NeXus Documentation Editor (Pete Jemian <jemian@anl.gov>).

Note: The indentation is very important to the syntax of the restructured text manual source. Be careful not to mix tabs and spaces in the indentation or the manual may not build properly.

9.3 Revision History

Browse the most recent Issues on the GitHub repository: https://github.com/nexusformat/definitions/pulse/weekly
9.4 Copyright and Licenses

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Publishing Information

This manual built 2014-10-05 17:22:12 CDT.

See Also:
This document is available in different formats:

PDF  nexus.pdf (available via online HTML link above)

A very brief overview is also available (separate from the manual).

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