

2. CONCEPTS AND SPECIFICATIONS

```

data_cell_length_
  loop__name      ' _cell_length_a'
                  ' _cell_length_b'
                  ' _cell_length_c'
  _category       cell
  _type           numb
  _type_conditions esd
  _enumeration_range 0.0:
  _units          A
  _units_detail   'angstroms'
  _definition
; Unit-cell lengths in angstroms corresponding to
  the structure reported.
;

```

Fig. 2.5.5.3. DDL1 definition of irreducible data items.

```

data_atom_site_attached_hydrogens
_name            ' _atom_site_attached_hydrogens'
  _category       atom_site
  _type           numb
  _list           yes
  _list_reference ' _atom_site_label'
  _enumeration_range 0:8
  _enumeration_default 0
  _definition
; The number of hydrogen atoms attached to the atom
  at this site excluding any hydrogen atoms for
  which coordinates (measured or calculated) are
  given.
;

```

Fig. 2.5.5.4. DDL1 definition of a 'list' data item.

blocks. The name of a data block matches the defined data name or the initial portion thereof. Each definition data block contains a sequence of data declarations, one for each attribute. The attributes within a data block constitute the total definition information, and only those attributes appropriate to a given defined item need be specified.

2.5.5.3. Definition example 3: irreducible data items

Some data items are closely related to other data items. Fig. 2.5.5.3 shows the definition of the length components `_cell_length_a`, `*_b` and `*_c` of the crystal unit-cell metric tensor. Because of their different enumeration and units attributes, the definition blocks for cell lengths and angles are specified separately. This definition shows how data items closely related in function may be defined in the same data block. The `loop__` command is used to list the multiple `_name` values. The attribute `_type_conditions` is used to indicate that these values are measurements and can have a standard uncertainty ('s.u.') value appended in parentheses *e.g.* 7.254(2). (The label 'esd' reflects the historical but inaccurate terminology 'estimated standard deviation'.) The `_units` and `_units_detail` attributes stipulate that the measurement unit of length is ångströms.

2.5.5.4. Definition example 4: list data

The definition of the number of hydrogen atoms attached to an atom site is shown in Fig. 2.5.5.4. The attributes `_list` and `_list_reference` signal whether the item is used in a list and what key item is in that list. As described in Chapters 2.1 and 2.2, a CIF list is a two-dimensional table where data items are in the columns (headed by the data names) and the values are the rows. The attribute `_list` is `yes` if a data item may be used in a list. The attribute `_list_reference` specifies the key item as

`_atom_site_label`. Each value of a key item must be unique so that the row of associated items can be accessed unambiguously. The attribute `_enumeration_default` specifies a default value for the defined item if it is absent from the data file.

2.5.5.5. Definition example 5: category information

Each CIF dictionary also contains example applications of defined data items, grouped according to category. Fig. 2.5.5.5 shows these for the category `ATOM_SITE`, as defined in the core dictionary (see Chapter 3.2). A `_type` attribute with a value of `null` specifies that a data block contains no definition information. The particular examples shown here illustrate that the reference key for the `ATOM_SITE` category is `_atom_site_label`. This data item is present in both examples shown, and in each packet the label is unique.

2.5.5.6. Definition example 6: mandatory and linked items

The example definition in Fig. 2.5.5.6 shows the attributes for the item `_atom_site_label`. As illustrated in Section 2.5.5.5, this item is the reference key to a list of items belonging to the category `ATOM_SITE`. The attribute `_list_mandatory`, which is set to a value of `yes`, specifies that this item is a mandatory item to a list of category `ATOM_SITE` items. The attribute `_list_link_child` identifies items in other categories that are 'linked' by derivation to this item and therefore share the same data values. This dependency is known as a child dependency. The definition shows that data items describing the labels of atom sites in a list of angular geometry, *i.e.* `_geom_angle_atom_site_label_1` and `_geom_angle_atom_site_label_2`, are the same labels described by `_atom_site_label` in the atom-site list. This is because the geometry is derived directly from the atom-site data. Note that this dependency requires that the `ATOM_SITE` list be present in the same CIF as the `GEOM_ANGLE` list, otherwise the molecular geometry information cannot be linked to the three-dimensional structural information.

2.5.5.7. Definition example 7: joinable lists

In the example definition shown in Fig. 2.5.5.7, the key item `_atom_site_aniso_label` is shown to have a special relationship with the key item `_atom_site_label`. This is because the data items in the list category `ATOM_SITE_ANISO` may be merged with (*i.e.* joined to) items in the list category `ATOM_SITE`. If this happens, the item `_atom_site_label` assumes the role as the key to the merged packets and this is signalled using the attribute `_list_link_parent`. When these two categories appear in separate lists, the category `ATOM_SITE_ANISO` data require the category `ATOM_SITE` data to be present in the data instantiation, but not *vice versa*. Note that the parent relationship, unlike that in the example of Section 2.5.5.6 arising from a derivation dependency, is because `ATOM_SITE_ANISO` is a subcategory of `ATOM_SITE`.

2.5.5.8. Definition example 8: equivalent items

As with the preceding example definition, this example concerns anisotropic atomic displacement parameters. Fig. 2.5.5.8 shows the definition of the component items `_atom_site_aniso_u_` that comprise an irreducible set of matrix elements. The key item in a list of `ATOM_SITE_ANISO` items is `_atom_site_aniso_label`. This definition illustrates how the attributes `_related_item` and `_related_function` are used to identify the similar data items `_atom_site_aniso_B_` that are atomic displacement parameters related by a simple conversion factor. Other attributes used to specify data relationships are described in Section 2.5.6 below.