```
1/
```

file:/Users/dawidl/courses/xml/external Examples/periodic Table Source Cropped.xml

```
<?xml version="1.0"?>
<!-- Data provided by a third-party vendor in a messy format -->
<PERIODIC_TABLE>
  <ATOM>
    <NAME>Actinium</NAME>
    <ATOMIC_WEIGHT>227</ATOMIC_WEIGHT>
    <ATOMIC_NUMBER>89</ATOMIC_NUMBER>
    <OXIDATION_STATES>3</OXIDATION_STATES>
    <BOILING_POINT UNITS="Kelvin">3470</BOILING_POINT>
    <SYMBOL>Ac</SYMBOL>
    <DENSITY UNITS="grams/cubic centimeter">
      10.07
    </DENSITY>
    <ELECTRON_CONFIGURATION>[Rn] 6d1 7s2 </ELECTRON_CONFIGURATION>
    <ELECTRONEGATIVITY>1.1/ELECTRONEGATIVITY>
    <ATOMIC_RADIUS UNITS="Angstroms">1.88</ATOMIC_RADIUS>
    <ATOMIC_VOLUME UNITS="cubic centimeters/mole">
      22.5
    </ATOMIC_VOLUME>
    <SPECIFIC_HEAT_CAPACITY UNITS="Joules/gram/degree Kelvin">
    </SPECIFIC_HEAT_CAPACITY>
    <IONIZATION_POTENTIAL>5.17</IONIZATION_POTENTIAL>
    <THERMAL_CONDUCTIVITY UNITS="Watts/meter/degree Kelvin">
    </THERMAL_CONDUCTIVITY>
  </ATOM>
  <ATOM>
    <NAME>Aluminum</NAME>
    <ATOMIC_WEIGHT>26.98154</ATOMIC_WEIGHT>
    <ATOMIC_NUMBER>13</ATOMIC_NUMBER>
    <OXIDATION_STATES>3</OXIDATION_STATES>
    <BOILING_POINT UNITS="Kelvin">2740</BOILING_POINT>
    <MELTING_POINT UNITS="Kelvin">933.5/MELTING_POINT>
    <SYMBOL>Al</SYMBOL>
    <DENSITY UNITS="grams/cubic centimeter">
     2.7
    </DENSITY>
    <ELECTRON_CONFIGURATION>[Ne] 3s2 p1 </ELECTRON_CONFIGURATION>
    <COVALENT_RADIUS UNITS="Angstroms">1.18</COVALENT_RADIUS>
    <ELECTRONEGATIVITY>1.61/ELECTRONEGATIVITY>
    <ATOMIC_RADIUS UNITS="Angstroms">1.43</ATOMIC_RADIUS>
    <HEAT_OF_VAPORIZATION UNITS="kilojoules/mole">
      290.8
    </HEAT_OF_VAPORIZATION>
    <ATOMIC_VOLUME UNITS="cubic centimeters/mole">
     10
    </ATOMIC_VOLUME>
    <HEAT_OF_FUSION UNITS="kilojoules/mole">
      10.7
    </HEAT_OF_FUSION>
    <IONIZATION_POTENTIAL>5.986</IONIZATION_POTENTIAL>
    <SPECIFIC_HEAT_CAPACITY UNITS="Joules/gram/degree Kelvin">
     0.9
    </SPECIFIC_HEAT_CAPACITY>
    <THERMAL_CONDUCTIVITY UNITS="Watts/meter/degree Kelvin">
      237
    </THERMAL_CONDUCTIVITY>
  </ATOM>
  <!-- ... Rest omitted ... -->
</PERIODIC_TABLE>
```

```
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```

file:/Users/dawidl/courses/xml/external Examples/derived/periodic Minimal.xsd

```
<?xml version="1.0" encoding="UTF-8"?>
<xs:schema</pre>
    xmlns:xs="http://www.w3.org/2001/XMLSchema"
    xmlns:p="urn:physics:elements"
    targetNamespace="urn:physics:elements"
    elementFormDefault="qualified"
    attributeFormDefault="unqualified">
    <xs:annotation><xs:documentation>
        This schema contains types which could be used to construct
        a minimal, but strict, periodic table of the elements.
    </xs:documentation></xs:annotation>
    <xs:complexType name="Element">
        <xs:annotation><xs:documentation>
            A minimal representation of a periodic element, including its
            state at room temperature or ('natural state'), assumed as
            a temperature of 300 Kelvin.
        </xs:documentation></xs:annotation>
        <xs:sequence>
            <xs:element name="name" type="xs:string"/>
            <xs:element name="symbol" type="p:AtomicSymbol"/>
            <xs:element name="atomicNumber" type="xs:positiveInteger"/>
            <xs:element name="atomicWeight" type="xs:decimal" min0ccurs="0"/>
        </xs:sequence>
        <xs:attribute name="naturalState" type="p:MaterialState" use="optional"/>
    </xs:complexType>
    <xs:complexType name="CollectionOfElements">
        <xs:annotation><xs:documentation>
            A sequence of elements that could represent a periodic table
        </xs:documentation></xs:annotation>
        <xs:sequence>
            <xs:element name="element" type="p:Element" maxOccurs="unbounded"/>
        </xs:sequence>
    </xs:complexType>
    <xs:simpleType name="MaterialState">
        <xs:annotation><xs:documentation>
            The state of an element: 'gas', 'liquid' or 'solid'.
        </xs:documentation></xs:annotation>
        <xs:restriction base="xs:string">
            <xs:enumeration value="gas"/>
            <xs:enumeration value="liquid"/>
            <xs:enumeration value="solid"/>
        </xs:restriction>
    </xs:simpleType>
    <xs:simpleType name="AtomicSymbol">
        <xs:annotation><xs:documentation>
            The symbol of an element, e.g. 'Al' or 'H'.
        </xs:documentation></xs:annotation>
        <xs:restriction base="xs:string">
            <xs:pattern value="[A-Z][a-z]\{0,2\}"/>
        </xs:restriction>
    </xs:simpleType>
    <xs:element name="periodicTable" type="p:CollectionOfElements">
        <xs:annotation><xs:documentation>
            A periodic table of elements
        </xs:documentation></xs:annotation>
    </xs:element>
</xs:schema>
```

```
file:/Users/dawidl/courses/xml/extemalExamples/derived/simplification.xsl
```

```
<?xml version="1.0" encoding="UTF-8"?>
<!-- An XSLT Transformation that transforms the complete periodic table,
    as received from a vendor in a messy format, into a minimal
    table which conforms to our XML schema. Performs the following
    key tasks:
    - Infers the state of the elements (gas, liquid, solid).
    - Translates to a namespaced vocabulary, and inserts a link to the schema
    - Inserts a link to a CSS stylesheet for visual rendering in a browser
    - Orders by atomic number
<xsl:stylesheet</pre>
    xmlns:xsl="http://www.w3.org/1999/XSL/Transform"
    xmlns:p="urn:physics:elements"
    xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance"
    version="1.0">
    <!-- Room temperature, in Kelvin (used to determine element state) -->
    <xsl:variable name="roomTemp">300</xsl:variable>
    <!-- The URI of the schema to which our output document conforms.
    In practical use, this should be a publically accessible, HTTP URL-->
    <xsl:variable name="schemaURI">periodicMinimal.xsd</xsl:variable>
    <!-- The URI of the CSS style sheet to attach to the resulting document -->
    <xsl:variable name="cssURI">periodicMinimal.css</xsl:variable>
    <!-- From the old document root, create a new (simplified) one which has links
    to the schema, etc -->
    <xsl:template match="PERIODIC_TABLE">
        <!-- Indicate that we want this processing instruction in the output document, which
        contains a link to a CSS style sheet. This enables the resulting document to be
        opened in any web browser for a pleasing view, whilst still containing all the
        information which applications / technical users might require. -->
        <xsl:processing-instruction name="xml-stylesheet">type="text/css" href="<xsl:value-of"</pre>
select="$cssURI"/>"</xsl:processing-instruction>
        <!-- Insert a comment in the output -->
        <xsl:comment>This simplification automatically produced with XSLT</xsl:comment>
        <!-- The root of our new (output) document. Note the inclusion of the 'schemaLocation'
            attribute, which means our output document is instantly validatable. -->
        <p:periodicTable xsi:schemaLocation="urn:physics:elements {$schemaURI}">
            <xsl:apply-templates>
                <!-- Sort by atomic number -->
                <xsl:sort select="ATOMIC_NUMBER" data-type="number" order="ascending"/>
            </xsl:apply-templates>
        </p:periodicTable>
    </xsl:template>
    <!-- Transform each atom to it's new (simplified) form -->
    <xsl:template match="ATOM">
        <p:element>
            <!-- If we have sufficient information, determine the element's
            natural state -->
            <xsl:if test="BOILING_POINT and MELTING_POINT">
                <xsl:attribute name="naturalState">
                    <xsl:call-template name="determineState">
                        <xsl:with-param name="atom" select="."/>
                    </xsl:call-template>
                </xsl:attribute>
            </xsl:if>
            <!-- Other fields (name, symbol, etc) simply copied through -->
            <p:name><xsl:value-of select="NAME"/></p:name>
```

file:/Users/dawidl/courses/xml/extemalExamples/derived/simplification.xsl

```
<p:symbol><xsl:value-of select="SYMBOL"/></p:symbol>
            <p:atomicNumber><xsl:value-of select="ATOMIC_NUMBER"/>p:atomicNumber>
            <!-- Optional (if we have it) atomic weight -->
            <xsl:if test="ATOMIC_WEIGHT">
                <p:atomicWeight><xsl:value-of select="ATOMIC_WEIGHT"/></p:atomicWeight>
            </xsl:if>
        </p:element>
    </xsl:template>
   <!-- A template to determine the state of an element (liquid/gas/solid)
   based on an assumed room temperature. -->
    <xsl:template name="determineState">
        <xsl:param name="atom"/>
        <xsl:choose>
            <xsl:when test="$atom/BOILING_POINT &lt;= $roomTemp">gas</xsl:when>
            <xsl:when test="$atom/MELTING_POINT &lt;= $roomTemp and $atom/BOILING_POINT &gt;=</pre>
$roomTemp">liquid</xsl:when>
            <xsl:when test="$atom/MELTING_POINT &gt;= $roomTemp and $atom/BOILING_POINT &gt;=</pre>
$roomTemp">solid</xsl:when>
        </xsl:choose>
    </xsl:template>
</xsl:stylesheet>
```

file:/Users/dawidl/courses/xml/external Examples/derived/simple Periodic Table Cropped.xml

```
<?xml version="1.0"?>
<?xml-stylesheet type="text/css" href="periodicMinimal.css"?>
<!--This simplification automatically produced with XSLT-->
<p:periodicTable xmlns:p="urn:physics:elements"</pre>
    xmlns:xsi="http://www.w3.org/2001/XMLSchema-instance"
    xsi:schemaLocation="urn:physics:elements periodicMinimal.xsd">
    <p:element naturalState="gas">
        <p:name>Hydrogen</p:name>
        <p:symbol>H</p:symbol>
        <p:atomicNumber>1</p:atomicNumber>
        <p:atomicWeight>1.00794</p:atomicWeight>
    </p:element>
    <p:element naturalState="gas">
        <p:name>Helium</p:name>
        <p:symbol>He</p:symbol>
        <p:atomicNumber>2</p:atomicNumber>
        <p:atomicWeight>4.0026</p:atomicWeight>
    </p:element>
    <p:element naturalState="solid">
        <p:name>Lithium</p:name>
        <p:symbol>Li</p:symbol>
        <p:atomicNumber>3</p:atomicNumber>
        <p:atomicWeight>6.941</p:atomicWeight>
    </p:element>
    <p:element naturalState="solid">
        <p:name>Beryllium</p:name>
        <p:symbol>Be</p:symbol>
        <p:atomicNumber>4</p:atomicNumber>
        <p:atomicWeight>9.01218</p:atomicWeight>
    </p:element>
  <!-- ... SEVERAL REMOVED ... -->
    <p:element>
        <p:name>ununnilium</p:name>
        <p:symbol>Uun</p:symbol>
        <p:atomicNumber>110</p:atomicNumber>
        <p:atomicWeight>269</p:atomicWeight>
    </p:element>
    <p:element>
        <p:name>unununium</p:name>
        <p:symbol>Uuu</p:symbol>
        <p:atomicNumber>111</p:atomicNumber>
        <p:atomicWeight>272</p:atomicWeight>
    </p:element>
    <p:element>
        <p:name>ununbium</p:name>
        <p:symbol>Uub</p:symbol>
        <p:atomicNumber>112</p:atomicNumber>
        <p:atomicWeight>277</p:atomicWeight>
    </p:element>
</p:periodicTable>
```

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element atomicNumber

file:/Users/dawidl/courses/xml/external Examples/derived/periodic Minimal.css

```
/* This CSS styles a periodic table in the "urn:physics:elements" namespace for
simple interactive display in a browser. Elements are rendered as coloured blocks
(based on them being liquid/solid/gas).
/* Periodic table as a whole */
periodicTable
   display: block;
    color: white;
    background-color: black;
    font-family: sans-serif;
    font-size: 10pt;
/* Individual atoms (display as small floating blocks,
by default as grey - state unknown) */
element
   display: block;
   width: 5em;
   height: 5em;
   margin: 0.2em;
   padding: 1em;
   border: 2px solid #BBB;
   background-color: #999;
    float: left;
    text-align: center;
/* Color according to states */
element[naturalState=gas]
    background-color: #339966;
element[naturalState=liquid]
    background-color: #336699;
element[naturalState=solid]
    background-color: #993366;
/* If the user's mouse moves over, change border */
element:hover
    border-color: white;
/* Contents of the box */
element name
    display: block;
    font-style: italic;
element symbol
    display: block;
    font-size: 2em;
    font-weight: bold;
   margin: 0.2em;
```

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file:/Users/dawidl/courses/xml/external Examples/derived/periodic Minimal.css

```
display: block;
}

/* Do not show atomic weight by default */
element atomicWeight
{
    display: none;
    font-family: monospace;
    color: #CCC;
}

/* ...only when user's mouse hovers over */
element:hover atomicWeight
{
    display: block;
}
```

