Practical Graph Isomorphism, II

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Isomorphism and Automorphism

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\[
\begin{align*}
(1) \\
(1 4)(2 3)(5 8)(6 7) \\
(1 8)(2 7)(3 6)(4 5) \\
(1 5)(2 6)(3 7)(4 8)
\end{align*}
\]
Ubiquity of graph isomorphism

Very many isomorphism problems can be efficiently expressed as graph isomorphism problems. For convenience, we use coloured graphs, in which the vertices (and possibly the edges) are coloured. Isomorphisms must preserve vertex and edge colour.

\[
\begin{pmatrix}
1 & 3 & 5 & 4 \\
1 & 5 & 4 & 3 \\
3 & 4 & 5 & 1
\end{pmatrix}
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Hadamard equivalence

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Group isomorphism

\[
\begin{bmatrix}
  e & g_1 & g_2 & g_3 \\
  g_1 & e & g_3 & g_2 \\
  g_2 & g_3 & e & g_1 \\
  g_3 & g_2 & g_1 & e
\end{bmatrix}
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The group elements are edge colours and vertices. The black vertex is the identity element. Preserve vertex colours but allow permutation of edge colours.
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Permutation equivalence of linear codes

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\begin{pmatrix}
0 & 1 & 1 & 0 & 1 & 0 & 0 & 0 & 0 \\
0 & 1 & 1 & 0 & 1 & 1 & 1 & 0 & 0 \\
1 & 0 & 0 & 1 & 0 & 0 & 1 & 1 & 1 \\
1 & 0 & 1 & 0 & 1 & 1 & 0 & 1 & 1 \\
0 & 0 & 1 & 0 & 0 & 0 & 0 & 1 & 0
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0 & 0 & 1 & 0 & 0 & 0 & 0 & 1 & 0
\end{pmatrix}
\]

Suppose isomorphism of two 0-1 matrices means that their columns can be permuted so that the rows generate the same vector space over $GF(2)$.

Nobody knows how to make an equivalent graph problem of size which is polynomial in the size of the generator matrix.
Theoretical complexity

The graph isomorphism problem is:

**GI:** Given two graphs, determine whether they are isomorphic.

- The fastest known algorithm for GI is \(2^{O(\sqrt{n \log n})}\) time; no polynomial time algorithm is known (despite many claims).
- GI is not known to be NP-complete.
- GI is not known to be in co-NP.
- There are polynomial time algorithms for many special classes of graphs (bounded degree, bounded genus, classes with excluded minors or topological subgraphs). Few of these are practical, planar graphs being a notable exception.

It is possible that group isomorphism is easier than GI and that permutation equivalence of linear codes is harder than GI. But nobody can prove it.
Canonical labelling

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Two labelled objects which are isomorphic become identical when they are canonically labelled. Since identity of objects is usually far easier to check than isomorphism, canonical labelling is the preferred method of sorting a collection of objects into isomorphism classes.
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A possible definition of the canonical member of an isomorphism class would be the member which maximises some linear representation (such as a list of edges).

In practice, most programs compute a canonical member which is easy for computers to find rather than easy for humans to define.
Software for graph isomorphism

This task was already described as a disease in 1970, and programs still appear regularly. The most successful programs still supported are:

- **nauty** (McKay, 1976–) canonical label and automorphism group
- **VF2** (Cordella, Foggia, Sansone and Vento, 1999–) comparison of two graphs
- **saucy** (Darga, Sakallah and Markov, 2004–) automorphism group
- **bliss** (Juntilla and Kaski, 2007–) canonical label and automorphism group
- **Traces** (Piperno, 2008–) canonical label and automorphism group
- **conauto** (López-Presa, Anta and Chiroque, 2009–) automorphism group and comparison of two graphs

Many similar principles appear in these programs. **saucy** and **bliss** began life as reimplementations of **nauty** but have since diverged.
Partition refinement

A key concept used by nauty is partition refinement. ("partition" = "colouring")

An equitable partition is one where every two vertices of the same colour are adjacent to the same number of vertices of each colour.

Two examples of equitable partitions:
Refinement of a partition means to subdivide its cells.

Given a partition (colouring) $\pi$, there is a unique equitable partition that is a refinement of $\pi$ and has the least number of colours.
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![Diagram showing the process of refining a partition and counting neighbours for different colours.](image-url)
Search tree

All competitive isomorphism programs generate a tree based on partition refinement. The nodes of the tree correspond to partitions (colourings).

The root of the tree corresponds to the initial colouring, refined.

If a node corresponds to a discrete partition (each vertex with a different colour), it has no children and is a leaf.

Otherwise we choose a colour used more than once (the target cell), individualize one of those vertices by giving it a new unique colour, and refine to get a child.
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However the tree can be much too large to generate completely.
Node invariants

A node invariant is a value $\phi(\nu)$ attached to each node in the tree that depends only on the combinatorial properties of $\nu$ and its ancestors in the search tree (not on the labelling of the graph).

Moreover, the order of node invariants must be preserved by descendants: Let $\nu_1, \nu_2$ be nodes at the same level in the search tree. If $\nu'_1$ is a descendant of $\nu_1$, and $\nu'_2$ is a descendant of $\nu_2$, then

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For example,

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Let $\phi^*$ be the greatest node invariant of a leaf. Then the lexicographically greatest labelled graph corresponding to a leaf $\nu$ with $\phi(\nu) = \phi^*$ is a canonical graph.
Pruning operations

Node invariants, together with automorphisms, allow us to remove parts of the search tree without generating them.

1. If $\nu_1, \nu_2$ are nodes at the same level in the search tree and $\phi(\nu_1) < \phi(\nu_2)$, then no canonical labelling is descended from $\nu_1$.

2. If $\nu_1, \nu_2$ are nodes at the same level in the search tree and $\phi(\nu_1) \neq \phi(\nu_2)$, then no labelled graph descended from $\nu_1$ is the same as one descended from $\nu_2$.

3. If $\nu_1, \nu_2$ are nodes with $\nu_2 = \nu_1^g$ for an automorphism $g$, then $g$ maps the subtree descended from $\nu_1$ onto the subtree descended from $\nu_2$. 
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Each of these observations enables us to prune parts of the search tree. The hope is that the remaining parts will be small enough to compute.
Variations between programs

The competing programs vary from each other in ways that include:

1. Data structures
2. Strength of the refinement procedure
3. Order of traversal of the search tree
4. Means of discovering automorphisms
5. Processing of automorphisms
Stronger refinement — \textit{nauty} “invariants”

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Example of stronger refinement performance

Consider random quartic graphs with 1000 vertices.

- `nauty` (dense data structures) 5.6 seconds
- `nauty` (sparse data structures) 0.5 seconds
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The big problem with these methods is that the best refinement technique varies from one graph class to another.

The user has to know which one to choose, which few do.
Tree traversal order — a Traces innovation
Tree traversal order — a \textit{Traces} innovation
Tree traversal order — a *Traces* innovation

Classical order: depth-first search
Tree traversal order — a Traces innovation

Traces order: breadth-first search
Tree traversal order — a Traces innovation

Traces order: breadth-first search

The problem with BFS is that automorphisms are discovered at leaves.
Tree traversal order — a *Traces* innovation
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Traces: experimental paths

Experimental paths allow automorphism detection during breadth-first search.
Refinement traces – a Traces innovation

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To alleviate this problem, *Traces* defines the node invariant to be a vector whose components are computed incrementally during the refinement. Then the refinement can often be aborted early. (This extends a technique introduced by *bliss.*)
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A possible trace (not Traces’ trace) is the number of colours at each step: (2,3,5,6).
Sparse automorphism detection — a saucy innovation

equitable partition
(no other blue or green)
Sparse automorphism detection — a *saucy* innovation

Equitable partition
(no other blue or green)

Individualize $v$

Individualize $w$
Sparse automorphism detection — a saucy innovation

This discovers the automorphism \((v \: w)(x \: y)\) without comparing leaves.
Automorphism group handling

Programs using depth-first search naturally produce automorphisms in the form of a base and strong generating set.

Other programs, like Traces, produce less predictable generators.

In order to perform automorphism-based pruning of the search tree, we need to efficiently (usually) determine if two sequences of vertices are equivalent under the group generated by the automorphisms found so far.
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Traces uses a technique called the random Schreier method to perform this computation probabilistically.

nauty also uses this in circumstances where the basic group handling is insufficient.
Fixing 1, 5, 9, 13: everything fixed
Fixing 1, 5, 9. Orbit of 13 is \{13, 15\}.
   (13 15)
Fixing 1, 5. Orbit of 9 is \{9, 11\}.
   (9 11)
Fixing 1. Orbit of 5 is \{5, 7, 13, 15\}.
   (5 7)
   (2 4)(5 13)(6 16)(7 15)(8 14)(10 12)
Fixing nothing. Orbit of 1 is \{1, 3, 5, 7, 9, 11, 13, 15\}.
   (1 3)
Group size = 2 \times 2 \times 4 \times 8 = 128.

This is a base and strong generating set.

There are at most \(n - 1\) generators, even though the group size can be up to \(n!\).
<table>
<thead>
<tr>
<th>Vertices</th>
<th>Time (sec)</th>
<th>Automorphism group</th>
<th>Canonical label</th>
</tr>
</thead>
<tbody>
<tr>
<td></td>
<td></td>
<td>Bliss 7.2</td>
<td>Nauty 2.5</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
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</tr>
</tbody>
</table>

**Figure 1:** Random graphs with $p = \frac{1}{2}$
<table>
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<tbody>
<tr>
<td>1</td>
<td>10^{-1}</td>
</tr>
<tr>
<td>10</td>
<td>10^{-2}</td>
</tr>
<tr>
<td>100</td>
<td>10^{-3}</td>
</tr>
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</tr>
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**Automorphism group**

**Canonical label**

Figure 2: Random graphs with average degree $\sqrt{n}$
Figure 3: Random cubic graphs
Figure 4: Random trees
Figure 5: Hypercubes
Figure 6: Miscellaneous vertex-transitive graphs
Figure 7: 2-dimensional and 3-dimensional grids
Figure 8: Latin square graphs
Figure 9: Strongly regular graphs
Figure 10: Hadamard matrix graphs
Figure 11: Projective planes of order 16
Figure 12: Cai-Fürer-Immermann graphs
Figure 13: Miyazaki graphs