

APPLICATION OF LINE GRAPHS AND COMPLETE HAMILTONIAN GRAPHS

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ABSTRACT

In 1856, Hamiltonian introduced the Hamiltonian Graph where a Graph which is covered all the vertices without repetition and end with starting vertex. In this Paper I would like to prove that Let “ G ” be a Complete graph with at least four vertices. Then, the line graph “ $L(G)$ ” is Complete Hamiltonian if and only if “ G ” is dominating trailable.

Key Words : Graph, Hamiltonian Graph, Complete Graph, Neighborhood, Locally Complete Graph.

Introduction :

Graphs, considered here, are finite, undirected and simple and complete Graphs being followed for terminology and notation. let $G = (V, E)$ be a graph, with V the set of vertices and E the set of edges. Suppose that W is a nonempty subset of V . The sub graph of G , whose vertex set is W and whose edge set is the set of those edges of G that have both ends in W , is called the sub

graph of G induced by W and is denoted by $G[W]$. For any vertex v in V , the *neighbour set* of v is the set of all vertices adjacent to v . This set is denoted by $N(v)$. For a graph $G = (V, E)$, we shall denote

$$\delta(G) = \min_{v \in V} |N(v)| \qquad \Delta(G) = \max_{v \in V} |N(v)|$$

a graph $G = (V, E)$ is locally complete, if for each vertex v the graph $G[N(v)]$ is complete. With every graph G , having at least one edge, there exists associated a graph $L(G)$, called the line graph of G , whose vertices, can be put in a one-to-one correspondence with the edges of G , in such a way that two vertices of $L(G)$ are adjacent if and only if the corresponding edges of G are adjacent.

The neighborhood is often denoted $N_G(v)$ or (when the graph is unambiguous) $N(v)$. The same neighborhood notation may also be used to refer to sets of adjacent vertices rather than the corresponding induced sub graphs. The neighborhood described above does not include v itself, and is more specifically the **open neighborhood** of v ; it is also possible to define a neighborhood in which v itself is included, called the **closed neighborhood** and denoted by $N_G[v]$. When stated without any qualification, a neighborhood is assumed to be open.

1.1 Definition: A graph – usually denoted $G(V,E)$ or $G = (V,E)$ – consists of set of vertices V together with a set of edges E . The number of vertices in a graph is usually denoted n while the number of edges is usually denoted m .

1.2 Definition: Vertices are also known as nodes, points and (in social networks) as actors, agents or players.

1.3 Definition: Edges are also known as lines and (in social networks) as ties or links. An edge $e = (u,v)$ is defined by the unordered pair of vertices that serve as its end points.

1.4 Example: The graph depicted in Figure 1 has vertex set $V=\{a,b,c,d,e,f\}$ and edge set

$E = \{(a,b),(b,c),(c,d),(c,e),(d,e),(e,f)\}$.

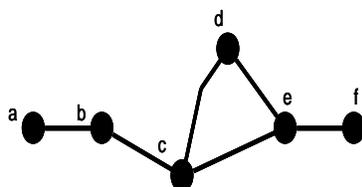


Figure 1.

1.5 Definition: Two vertices u and v are *adjacent* if there exists an edge (u,v) that connects them.

1.6 Definition: An edge (u,v) is said to be *incident* upon nodes u and v .

1.7 Definition: An edge $e = (u,u)$ that links a vertex to itself is known as a *self-loop* or *reflexive* tie.

1.8 Definition: Every graph has associated with it an *adjacency matrix*, which is a binary $n \times n$ matrix A in which $a_{ij} = 1$ and $a_{ji} = 1$ if vertex v_i is adjacent to vertex v_j , and $a_{ij} = 0$ and $a_{ji} = 0$ otherwise. The natural graphical representation of an adjacency matrix is a table, such as shown below.

	a	b	c	d	e	f
a	0	1	0	0	0	0
b	1	0	1	0	0	0
c	0	1	0	1	1	0
d	0	0	1	0	1	0
e	0	0	1	1	0	1
f	0	0	0	0	1	0

Adjacency matrix for graph in Figure 1.

1.9 Definition: Examining either Figure 1 or given adjacency Matrix, we can see that not every vertex is adjacent to every other. A graph in which all vertices are adjacent to all others is said to be *complete*.

1.10 Definition: While not every vertex in the graph in Figure 1 is adjacent, one can construct a sequence of adjacent vertices from any vertex to any other. Graphs with this property are called *connected*.

1.11 Note: Reachability. Similarly, any pair of vertices in which one vertex can reach the other via a sequence of adjacent vertices is called *reachable*. If we determine reachability for every pair of vertices, we can construct a reachability matrix R such as depicted in Figure 2. The matrix R can be thought of as the result of applying transitive closure to the adjacency matrix A .

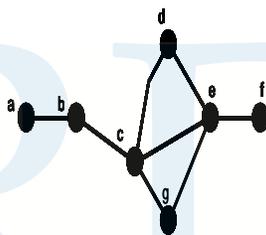


Figure: 2

1.12 Definition : A walk is closed if $v_0 = v_n$. *degree* of the vertex and is denoted $d(v)$.

1.13 Definition : A *tree* is a connected graph that contains no cycles. In a tree, every pair of points is connected by a unique path. That is, there is only one way to get from A to B.

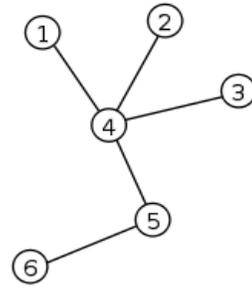


Figure 3: A labeled tree with 6 vertices and 5 edges

1.14 Definition: A *spanning tree* for a graph G is a sub-graph of G which is a tree that includes every vertex of G .

1.15 Definition: The length of a walk (and therefore a path or trail) is defined as the number of edges it contains. For example, in Figure 3, the path a,b,c,d,e has length 4.

1.16 Definition: The number of vertices adjacent to a given vertex is called the *degree* of the vertex and is denoted $d(v)$.

1.17 Definition : In the mathematical field of graph theory, a bipartite graph (or bigraph) is a graph whose vertices can be divided into two disjoint sets U and V such that every edge connects a vertex in U to one in V ; that is, U and V are independent sets. Equivalently, a bipartite graph is a graph that does not contain any odd-length cycles.

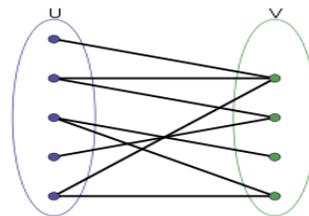


Figure 4: Example of a bipartite graph.

1.18 Definition : An Eulerian circuit in a graph G is circuit which includes every vertex and every edge of G . It may pass through a vertex more than once, but because it is a circuit it traverse each edge exactly once. A graph which has an Eulerian circuit is called an Eulerian graph. An Eulerian path in a graph G is a walk which passes through every vertex of G and which traverses each edge of G exactly once

1.19 Example : Königsberg bridge problem: The city of Königsberg (now Kaliningrad) had seven bridges on the Pregel River. People were wondering whether it would be possible to take a walk through the city passing exactly once on each bridge. Euler built the representative graph, observed that it had vertices of odd degree, and proved that this made such a walk impossible.

Does there exist a walk crossing each of the seven bridges of Königsberg exactly once?

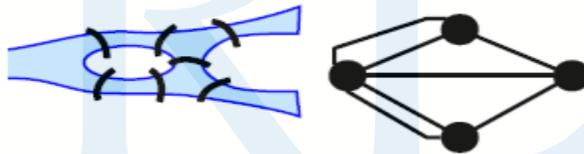


Figure 5: Königsberg problem

2. Complete Graphs, Locally Complete Graphs, Hamiltonian Graphs, Line Graphs

In this section we have to prove that main theorem using definitions.

2.1 Definition: A Hamilton circuit is a path that visits every vertex in the graph exactly once and return to the starting vertex. Determining whether such paths or circuits exist is an NP-complete problem. In the diagram below, an example Hamilton Circuit would be

2.2 Example :

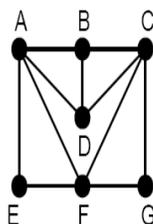


Figure: Hamilton Circuit would be AEFGCDBA.

2.3 Definition : Complete Graph: A simple graph in which there exists an edge between every pair of vertices is called a complete graph.

2.4 Definition : Let $\{v_1, v_2, \dots, v_n\}$ be the vertex set of a graph G , and for each ' α '.

let N_i^* denote the closed neighborhood of v_a . Let N_a be any subset of N_a^* containing v_a which generates a complete subgraph C_a of G . Then C_a is called a complete sub neighborhood of v_a , and the indexed family $C^* = \{C_1, C_2, \dots, C_n\}$ is called a complete family for G if $G = \bigcup C^*$. A graph G is called locally complete iff G has at least one complete family.

2.5 Examples : It is easily seen that complete graphs, trees, and unicyclic graphs are also locally complete.

The complete bigraph $K_{3,2}$ is the smallest (nontrivial, connected) graph which fails to be locally complete.

2.6 Proposition : If G is Hamiltonian, then $L(G)$ is Hamiltonian.

Proof : This is a nice, basic result to see if a line graph is Hamiltonian.

A graph is Hamiltonian if there exists a Hamiltonian cycle in the graph.

It may be easier to find a Hamiltonian cycle in G than $L(G)$, but from this proposition, we would get that $L(G)$ is Hamiltonian.

2.7 Theorem : Let ' G ' be a complete graph having $n \geq 3$ vertices then $L(G)$ is Complete Hamiltonian

2.8 Theorem : Let G be a Complete graph with at least four vertices. Then, the line graph $L(G)$ is Complete Hamiltonian if and only if G is dominating trail able.

Proof. We begin by assuming $L(G)$ is Complete Hamiltonian.

So, between any two vertices, x and y , in $L(G)$,

we have a Hamiltonian path written as

$x = x_0, x_1, x_2, \dots, x_n = y$, where $n + 1$ is the number of vertices in $L(G)$.

Since the vertices of $L(G)$ correspond to the edges of G , then $x_0, x_1, x_2, \dots, x_n$ is a sequence of edges in G , where $x_i \in E(G)$ for $i = 0, 1, \dots, n$.

Let v_i be the common vertex between x_i and x_{i+1} in G and create a list of vertices as

v_0, v_1, \dots, v_n .

Now, in this list of vertices v_0, v_1, \dots, v_n , some vertices may appear more than once.

So, create a subset w_1, w_2, \dots, w_k , where each vertex appears only once and $k \leq n$.

In creating this, once a vertex is listed, we won't list it again.

Now, for two vertices w_i and w_{i+1} , where $i = 1, 2, \dots, k$, list the corresponding edge between these two as e_i .

Then, $x_1, w_1, e_1, w_2, e_2, \dots, w_k, e_k, y$ is a dominating trail in G between edges x and y ,

since every edge in G is incident with one of w_1, w_2, \dots, w_k .

Since this trail works for all edges in G ,

we can say that G is dominating trailable.

Conversely, we can assume G is dominating trailable, and let x and y be edges of G .

Then, there exists a dominating trail between x and y written as $x, v_1, x_1, v_2, \dots, v_n, x_n = y$, where $x_i \in E(G)$ for all $i = 1, 2, \dots, n$.

So, n is the number of internal vertices of the trail. For the remaining edges not listed in the dominating trail, we will partition in the following way.

Create n sets, labeled S_1, S_2, \dots, S_n .

Next, for an edge incident with v_i , place that edge in the corresponding set, S_i .

Then, start this process with v_1 , and once an edge is placed in a set, do not consider that edge again. Notice that some sets may be empty, and some sets may have more than one element.

Define the elements of S_i as $s_{i,1}, 1, s_{i,2}, 2, \dots, s_{i,r}, r$ where r is the length of S_i .

Then, consider the list $x, S_1, x_1, S_2, \dots, S_n, y$ written as

$x, s_{1,1}, 1, s_{1,2}, 2, \dots, s_{1,r}, r, x_1, s_{2,1}, 1, \dots, s_{n,r}, r, y$.

Since the edges of G correspond to

This the vertices of $L(G)$, we now classify this sequence as a list of vertices in $L(G)$. This sequence is a path, since it consists of distinct vertices of $L(G)$ with each vertex in the list adjacent to the one before and after it. By construction, we have accounted

for every edge in G , and thus every vertex in $L(G)$. makes the path

$x, s_{1,1}, 1, s_{1,2}, 2, \dots, s_{1,r}, r, x_1, s_{2,1}, 1, \dots, s_{n,r}, r, y$ a Hamiltonian path in $L(G)$.

Since this is true for any x and $y \in E(G)$,

$L(G)$ is Complete Hamiltonian

Hence The Theorem.

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