Final Report of Traineeship Program 2025

On

"Analysis of Chemical Components"

MEDTOUREASY



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ABSTRACT

The cosmetic industry has witnessed remarkable growth in recent years, largely fueled by consumers who are increasingly seeking products tailored to their specific skin concerns and ingredient preferences. However, navigating the vast amount of information on product labels especially complex ingredient lists can be overwhelming, making it difficult for consumers to choose products that best suit their needs. This project aims to address this challenge by developing a content-based recommendation system that suggests cosmetic products based on ingredient similarity.

Using a dataset from Sephora, the system analyzes and processes ingredient lists to recommend alternative products with comparable compositions. To handle the high dimensional nature of ingredient data, the project utilizes t-Distributed Stochastic Neighbor Embedding (T-SNE) for dimensionality reduction, effectively mapping the data into a 2D space for easier visualization. Bokeh, an interactive visualization library, is employed to build an intuitive user interface, enabling users to explore and compare products based on their ingredient profiles.

The primary objective of this system is to help consumers make informed decisions by identifying suitable alternatives that match their skin type and avoid potentially harmful or undesirable ingredients. By simplifying the interpretation of complex ingredient information, the system offers a user-friendly platform for personalized skincare recommendations.

In particular, selecting new skincare products can be a daunting task for individuals with sensitive skin due to the technical nature of ingredient terminology. This project specifically focuses on moisturizers for dry skin, leveraging data from 1,472 products listed on Sephora. By analyzing ingredient content, the system provides tailored recommendations, making the product selection process more accessible and less intimidating. This report details the methodology, data preprocessing steps, dimensionality reduction techniques, system implementation, results, and potential future enhancements to improve the model's accuracy and usability.



I. Introduction

1.1 **About the Company**

MedTourEasy, a global healthcare company, provides you the informational resources needed to evaluate your global options. MedTourEasy provides analytical solutions to our partner healthcare providers globally.

1.2 **About the Project**

The Content-Based Recommendation System for Cosmetics project addresses a common challenge faced by consumers in selecting cosmetic products that align with their skin type and ingredient preferences. With an increasing awareness of the ingredients used in skincare products, many individuals seek transparency and guidance in their purchasing decisions. This project aims to simplify the decision-making process by providing a tool that recommends cosmetic products based on their ingredient similarities.

The project focuses on moisturizers suitable for dry skin, utilizing a dataset of 1,472 cosmetic products from Sephora. By analyzing the ingredient lists of these products, the system identifies potential alternatives that share similar compositions. This approach not only aids consumers in making informed choices but also enhances their understanding of cosmetic ingredients.

Key Features:

- 1. Data Collection and Preparation: The project starts with gathering and cleaning the dataset, ensuring ingredient information for each product is accurately structured and ready for analysis.
- 2. Ingredient Tokenization: Ingredient lists are broken down into individual tokens to build a meaningful and standardized vocabulary, which forms the foundation for comparing product similarities.
- 3. Construction of Document-Term Matrix (DTM): A matrix is created to numerically represent the presence and frequency of each ingredient across products, enabling structured comparison of ingredient compositions.
- 4. Visual Representation through t-SNE: To simplify the interpretation of high-dimensional ingredient data, t-SNE is applied to reduce the dimensions and map the products into a 2D space, making it easier to visualize and explore their similarities.



Objectives:

The primary objectives of this project include:

- Simplifying ingredient analysis by offering users clear, data-driven recommendations that demystify complex cosmetic formulations.
- Assisting users in identifying alternative products that align with their individual skin concerns, ingredient preferences, and sensitivity needs.
- Raising awareness of ingredient impact by promoting better understanding of how specific cosmetic ingredients influence skincare outcomes.
- This project functions as a practical decision-support tool for consumers navigating the skincare market. Through the use of data science and interactive visualizations, it streamlines the product selection process while empowering users to make well-informed and personalized choices. Potential future improvements include broadening the range of product types, incorporating user feedback and reviews, and integrating toxicity analysis to enhance recommendation precision and safety.



1.3 Objectives and Deliverables

Objectives:

1. Development of a Content-Based Recommendation System:

o Create a robust system that utilizes ingredient similarity to recommend cosmetic products. The recommendation engine will analyze the composition of various moisturizers, focusing on those suitable for dry skin, and suggest alternatives based on similar ingredients.

2. Empower Consumers to Compare Products:

 Provide users with the ability to easily compare multiple cosmetic products side by side. This feature will allow consumers to make informed decisions by assessing key attributes such as ingredient profiles, effectiveness, and suitability for their skin type.

3. Enhance User Understanding of Ingredients:

 Educate consumers about the importance of cosmetic ingredients by highlighting their functions and potential benefits or risks. This will help users develop a better understanding of what they are applying to their skin.

4. Facilitate Personalized Recommendations:

Tailor the recommendation system to consider individual preferences, such as skin type and specific ingredient avoidance (e.g., allergens or irritants). This personalization will improve user satisfaction and trust in the recommendations provided.

5. Interactive Visualization of Similarities:

 Create an engaging and intuitive visual representation of product similarities using dimensionality reduction techniques. The visualization will help users grasp complex ingredient relationships more easily.



Deliverables:

1. **Document-Term Matrix (DTM)**:

 A structured matrix that captures the frequency of ingredients across various cosmetic products. This DTM will serve as the foundation for subsequent analysis, enabling the calculation of ingredient similarities.

2. Dimensionality Reduction Visualization:

o Implementation of the t-SNE algorithm to reduce the high-dimensional data of the DTM into a two-dimensional space. This will facilitate easier analysis and interpretation of ingredient similarities among products.

3. Interactive Scatter Plot:

- o A visually appealing scatter plot created using Bokeh, allowing users to interact with the data. Features will include:
 - **Hover Functionality**: Users can hover over points to view detailed information about each product, including its name, brand, price, and key ingredients.
 - Customization Options: Users may filter the visualization based on specific criteria, such as product type or brand.

4. Ingredient Comparison Tool:

A feature enabling users to select multiple products and compare their ingredient lists side by side. This will highlight similarities and differences, aiding users in making informed choices about their skincare products.

5. Comprehensive Project Documentation:

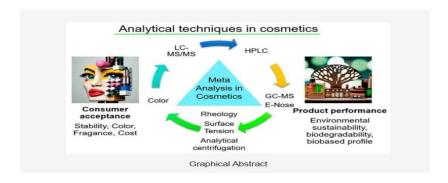
A detailed report that outlines the entire project process, including data collection, preprocessing, methodology, results, and analysis. This documentation will serve as a reference for future improvements and for those interested in replicating or building upon the project.

6. Future Scope Recommendations:

Suggestions for potential enhancements to the recommendation system, including integrating user reviews, expanding to additional product categories, and conducting toxicity analyses to flag harmful ingredients. This will provide a roadmap for future development efforts.



II. METHODOLOGY



2.1 Flow of the Project & Use Case Diagram

The project followed a well-defined workflow to ensure a systematic and efficient development process. It began with data acquisition, sourcing a cosmetics dataset from Sephora, followed by thorough preprocessing to clean and organize the information. The scope was specifically limited to moisturizers formulated for dry skin. ingredient lists were then tokenized, breaking them down into simpler, analyzable elements, which were used to construct a Document-Term Matrix (DTM) representing ingredient frequency across products. To better interpret the high-dimensional data, the t-Distributed Stochastic Neighbor Embedding (t-SNE) technique was employed for dimensionality reduction, enabling effective visualization of ingredient similarities. Finally, Bokeh was utilized to build an interactive visualization interface, allowing users to intuitively explore and compare products based on their ingredient profiles.

The system supports various use cases that cater to different stakeholders, including:

- **Consumers**: Individuals searching for alternative moisturizers with similar ingredients that align with their skin type and personal preferences.
- **Dermatologists**: Professionals analyzing common ingredients in a patient's skincare routine, allowing them to offer informed recommendations and insights.
- **Researchers**: Academics investigating the prevalence and impact of certain chemicals across skincare products, contributing to broader studies on cosmetic safety and efficacy.
- Retailers: Businesses aiming to enhance product offerings by understanding consumer

preferences based on ingredient analysis.

• This diverse range of use cases highlights the system's versatility and potential impact across different sectors.



- cosmetics have been used to enhance a person's appearance and to improve the quality of their skin.
- Chemical analysis is often used to ensure both the purity and quality of cosmetics and other dermatological formulations.
- Many cosmetics and other dermatological products contain toxic ingredients that may be harmful.



2.2 Data Preprocessing & Techniques Used: t-SNE and Bokeh

The preprocessing phase involved several key steps crucial for ensuring the dataset was clean and ready for analysis:

- 1. **Filtering the Dataset**: The dataset was filtered to isolate products labeled as "Moisturizers" specifically targeting consumers with dry skin. This narrowed focus ensured the subsequent analysis was relevant to the intended audience.
- 2. **Resetting the Index**: The index of the Data Frame was reset to facilitate easier manipulation of the data, allowing for straightforward referencing and data handling during the analysis.
- 3. **Tokenization of Ingredients**: Ingredients were tokenized to break down complex ingredient lists into individual components. This process involved converting all text to lowercase and splitting the strings based on a specified delimiter (', '), resulting in a list of individual ingredients for each product.
- 4. **Data Cleaning**: Additional cleaning steps included removing any duplicates, handling missing values, and ensuring consistent formatting across the dataset. This comprehensive cleaning was essential for generating accurate and reliable recommendations.

The preprocessing efforts ensured that the data was not only clean but also well-structured for subsequent analysis, which is critical for generating accurate recommendation outputs.

The project employed two key techniques for analysis and visualization:

- t-SNE (t-Distributed Stochastic Neighbor Embedding): This technique was crucial for reducing the dimensionality of the ingredient space. By mapping each product to a two-dimensional space based on ingredient similarity, t-SNE facilitated the visualization of relationships between products, making it easier to identify clusters and similarities among moisturizers.
- **Bokeh**: Bokeh was utilized to create an interactive visualization platform. The library allowed for the development of an engaging scatter plot that displayed product details dynamically when users interacted with the data points.



2.3 Tokenization and Document-Term Matrix

To quantify the presence of each ingredient in a product, the following steps were implemented:

Tokenization: The ingredient list for each product was tokenized by first converting the text to lowercase and then splitting it into individual tokens using a predefined delimiter.

Document-Term Matrix (DTM) Initialization: A matrix of zeros was initialized to represent the presence (or absence) of each ingredient in each product. The dimensions of the matrix were determined by the number of products in the dataset (rows) and the total number of unique ingredients identified (columns). This matrix served as the foundation for subsequent ingredient-based analysis, allowing for the quantification of ingredient similarities across different products.



III. IMPLEMENTATION

The overall goal of the project is to build a system that can recommend cosmetic products (specifically moisturizers for dry skin) based on ingredient similarity. This system will allow consumers to compare products and find alternatives, especially when they are looking for specific ingredients in their skincare routine.

3.1 Importing and Inspecting the Dataset

Step Explanation:

The first step is loading the dataset into Python for analysis. The dataset contains product details from Sephora, including product name, brand, price, and a list of ingredients for each product.

- **Pandas Library**: We use the Pandas library to import the dataset and inspect its structure.
- **Data Inspection**: It's crucial to inspect the dataset by displaying a few rows and checking for any missing values or inconsistencies. We also check the different categories of products available, such as "Moisturizers," "Cleansers," etc., by counting the frequency of each product type.

Key Points:

- **Data Loading**: We use Pandas to read the CSV file into a Data Frame.
- **Basic Inspection**: We display the first few rows to understand how the data is structured.
- Understanding Labels: We count how many products belong to each label (e.g., moisturizer, cleanser, etc.) to ensure we have sufficient data for the analysis.

```
import pandas as pd
import numpy as np

# Load the cosmetics dataset
df = pd.read_csv("datasets/cosmetics.csv")

# Inspect the dataset - display first few rows
display(df.head())

# Check the structure of the dataset and types of columns
df.info()

# Count the number of unique product categories in the dataset
print(df['Label'].value_counts())
```

3.2 Filtering and Tokenizing Ingredients

Step Explanation:

In this step, we filter the dataset to focus on a specific product category in this case, "Moisturizers" that are suitable for dry skin. This helps narrow down our analysis and provides a more specific recommendation system.

- **Filtering**: Filtering is performed to extract only the products labeled as "Moisturizer" for dry skin, as we aim to build a recommendation system for this specific category.
- **Tokenizing Ingredients:** The ingredient list of each product was split into individual components to create a structured list. This process is crucial for transforming unstructured text data into an analyzable format. We used the. split() function with a comma as the delimiter to separate the ingredients.

```
# Filter dataset for moisturizers targeting dry skin
moisturizers = df[df['Label'] == 'Moisturizer']
moisturizers_dry = moisturizers[moisturizers['Dry'] == 1].reset_index(drop=True)

# Tokenize the ingredients
corpus = []
for product in moisturizers_dry['Ingredients']:
    tokens = product.lower().split(', ') # Split ingredients into individual components
    corpus.append(tokens)

# Display a sample of the tokenized ingredients
print(corpus[:2]) # Display first two tokenized ingredient lists
```

Key Points:

- **Filtering**: We focus only on moisturizers labeled for dry skin by using condition-based filtering.
- **Tokenization**: Breaking the ingredient list into individual components allows us to treat each ingredient as a feature for analysis.

The dataset was filtered to focus on a specific category of products, such as "Moisturizers" for dry skin. The ingredients for each product were then tokenized, converting each list of ingredients into individual components to form a corpus for further analysis.

3.3 Document-Term Matrix Initialization

Step Explanation:

The **Document-Term Matrix (DTM)** is a fundamental part of text analysis. In our case, each product's ingredient list is like a "document," and each ingredient is a "term." The DTM represents which ingredients are present in each product.

- Matrix Representation: We initialize a matrix where each row represents a product and each column represents a unique ingredient across all products. A "1" in the matrix means the ingredient is present in the product, while a "0" indicates absence.
- **Count Vectorizer**: We use Count Vectorizer from the scikit-learn library to build this matrix automatically. It tokenizes the text and creates a matrix where each column represents a unique ingredient and each row represents a product.

Key Points:

- **Document-Term Matrix (DTM)**: This matrix is the foundation for comparing products based on ingredient similarity.
- **Binary Representation**: The matrix encodes whether an ingredient is present or not in each product, enabling quantitative analysis.

```
python

from sklearn.feature_extraction.text import CountVectorizer

# Create a document-term matrix
vectorizer = CountVectorizer(tokenizer=lambda x: x.split(', '))
dtm = vectorizer.fit_transform(moisturizers_dry['Ingredients'])

# Convert the matrix to a dense array
ingredient_matrix = dtm.toarray()

# Get the feature names (ingredients)
ingredients = vectorizer.get_feature_names_out()
print(ingredients[:10]) # Display the first 10 unique ingredients
```

3.4 Dimensionality Reduction Using t-SNE

Step Explanation:

A document-term matrix for cosmetics can have hundreds of dimensions because of the vast number of unique ingredients. Visualizing this in its raw form would be difficult. To simplify the data and make it easier to visualize, we reduce its dimensionality using T-SNE (t-distributed Stochastic Neighbor Embedding).

- **t-SNE** is a popular machine learning algorithm for reducing data from a high-dimensional space to two or three dimensions while preserving the relationships between data points (in this case, products).
- **Purpose**: This allows us to plot products on a 2D plane, where products with similar ingredients are placed closer together.

Key Points:

- **Dimensionality Reduction**: We use t-SNE to reduce the number of dimensions in the data, which simplifies visualization.
- **Product Similarity**: Products that are close to each other in this 2D space have similar ingredients, making this step critical for our recommendation system.

```
python

from sklearn.manifold import TSNE

# Apply t-SNE to reduce dimensions
model = TSNE(n_components=2, random_state=42)
tsne_features = model.fit_transform(ingredient_matrix)

# Add the t-SNE features back to the DataFrame
moisturizers_dry['X'] = tsne_features[:, 0] # X-coordinate
moisturizers_dry['Y'] = tsne_features[:, 1] # Y-coordinate
```

3.5 Visualizing Ingredient Similarity Using Bokeh

Step Explanation:

Once we have the 2D coordinates from t-SNE, we use Bokeh, a Python library for creating interactive visualizations, to plot the products in a scatter plot. Each point represents a product, and its position is based on its ingredient composition.

- **Scatter Plot**: The scatter plot visualizes products as points on a plane. Products that are closer together have more similar ingredients. Users can interact with the plot to explore products visually.
- **Hover Tool**: We add interactivity by enabling hover functionality so that when a user hovers over a point, they can see the product name, brand, and price.

```
python
                                                                               Copy code
from bokeh.plotting import figure, show, ColumnDataSource
from bokeh.models import HoverTool
# Create a ColumnDataSource for Bokeh
source = ColumnDataSource(moisturizers dry)
# Create a scatter plot
plot = figure(title="Cosmetic Ingredient Similarity",
              x_axis_label='T-SNE 1', y_axis_label='T-SNE 2',
              plot width=800, plot height=800)
# Add circles to represent products
plot.circle(x='X', y='Y', size=10, source=source, color="navy", alpha=0.6)
# Add hover tool to show product details
hover = HoverTool(tooltips=[
   ("Product", "@`Product Name`"),
   ("Brand", "@Brand"),
   ("Price", "@Price")
1)
plot.add tools(hover)
```

Key Points:

- **Interactive Visualization**: Bokeh allows users to explore the data interactively, making the recommendation system more user-friendly.
- **Ingredient-Based Product Similarity**: The position of each point (product) on the scatter plot is determined by the similarity of its ingredients to other products.

3.6 Adding Hover Functionality

Step Explanation:

To enhance the usability of the scatter plot, we add hover functionality, allowing users to interact with the plot in a meaningful way. When a user hovers over a product point, they can see more information, such as the product name, brand, and price.

• **User Experience**: This step significantly improves the user experience by providing detailed information directly on the plot without requiring additional navigation.

3.7 Result Analysis

Step Explanation:

Once the visualization is in place, we can analyze the relationships between products based on their ingredient similarity. Products that are located close to each other in the scatter plot have similar ingredients, which means they can be used as alternatives to one another.

• **Product Comparison**: We also implemented a function to identify the nearest products to a given product based on ingredient similarity. This would provide concrete recommendations when a user selects a specific product.

```
Copy code
python
# Example of how we can analyze nearest products using distances
from sklearn.metrics.pairwise import cosine similarity
# Calculate similarity between products
similarity matrix = cosine similarity(ingredient matrix)
# Function to find top similar products for a given product
def recommend similar products(product idx, top n=5):
    similarity scores = similarity matrix[product idx]
    similar indices = similarity scores.argsort()[::-1][1:top n+1] # Top N similar product
    return moisturizers dry.iloc[similar indices][['Product Name', 'Brand', 'Price']]
# Example: Recommend products similar to the first product
recommend similar products(0)
```

Key Points:

- Cosine Similarity: We use the cosine similarity metric to compare products based on their ingredient lists.
- **Top N Recommendations**: This function allows us to recommend the top N products most similar to a given product.

we created a content-based recommendation system for moisturizers using the ingredient lists of cosmetic products. By applying text processing, dimensionality reduction, and machine learning techniques, we developed an interactive system that enables users to find product alternatives based on their ingredient composition. This can help users find products with desired ingredients or avoid specific components.

IV. Results and Analysis

In this section, we delve deeply into the outcomes of the content-based recommendation system that was built to suggest alternative cosmetic products based on their ingredient composition. The project is designed specifically for users looking for moisturizers targeted at dry skin, but the methodology can easily be expanded to include other types of cosmetic products. Below is a detailed breakdown of each aspect of the results and their significance.

4.1 Mapping the Cosmetic Items

Mapping cosmetic items was a key step in analyzing how different products relate to each other based on their ingredient composition. Each product contains a unique blend of ingredients, making direct comparisons in their original high-dimensional format (often involving hundreds of unique ingredients) both computationally intensive and difficult to visualize. To address this, we employed t-distributed Stochastic Neighbor Embedding (t-SNE), a powerful technique for dimensionality reduction.

Dimensionality Reduction

Cosmetic products were represented as high dimensional vectors, where each dimension corresponds to a specific ingredient. For example, with 300 unique ingredients across all moisturizers, each product becomes a 300dimensional binary vector where 1 indicates the presence of an ingredient and 0 its absence. Although this representation captures rich ingredient information, it is not practical for direct interpretation or visualization. Dimensionality reduction simplifies this complexity.

t-SNE Output

t-SNE reduces these high-dimensional vectors into a 2-dimensional space while preserving the relative distances between products. This means that products with similar ingredient compositions appear closer together on the 2D plane. The result is a transformed dataset with two components t-SNE 1 and t-SNE 2 which can be plotted for easy visual interpretation.

Understanding Clusters

The 2D scatter plot generated using t-SNE reveals meaningful clusters of cosmetic products. Each point represents a product, and groups of nearby points indicate products that share similar ingredient profiles. For instance, moisturizers that all contain ingredients like hyaluronic acid and ceramides tend to cluster together. These clusters provide valuable insights into product similarity, potential shared effects, formulation strategies, and targeted skin benefits.

Key Insights:

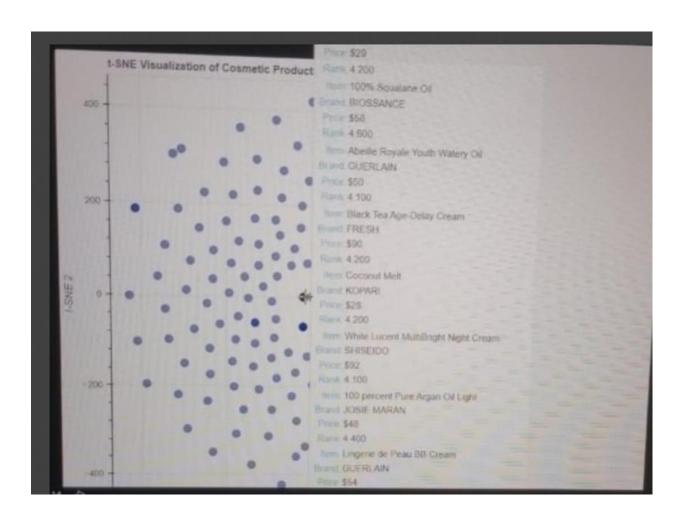
- **Identifying Substitute Products**: Products that are positioned closer together in the t-SNE space are likely to serve as substitutes or alternatives. This is useful for consumers who may want to switch from one brand to another without drastically changing the ingredients they are using.
- **Segmentation by Formulation**: The clusters on the plot might correspond to different types of formulations, such as products that are oil-based versus water-based. Such clusters can reveal market segmentation and product positioning within the cosmetics industry.
- Understanding Ingredient Trends: By looking at clusters, we can infer broader ingredient trends, such as a group of moisturizers that avoid potentially irritating preservatives like parabens or products focused on natural or organic ingredients.

4.2 Ingredient Similarity Visualization

The core of the project is the ability to visualize and interact with product data based on ingredient similarity. The power of this system lies in its ability to transform raw data into actionable insights through intuitive and interactive visualizations. For this purpose, we utilized Bokeh, a powerful Python library for creating interactive plots.

- Creating a Scatter Plot: Once the t-SNE results were computed, we used Bokeh to create an interactive scatter plot, where each point represents a moisturizer. The axes represent the two t-SNE dimensions, and the positions of the points reveal ingredient similarities. Products with similar ingredients appear closer together, while products with very different ingredient compositions are more distant.
- **Hover Tool for Interactivity**: One of the primary advantages of using Bokeh was its support for interactive elements, such as the hover tool. When the user hovers over a data point (representing a product), they can view additional information about the product, including the brand name, product name, and price. This feature enhances the user experience by providing a rich, contextual understanding of each product's details without cluttering the scatter plot.

- **User Exploration**: The interactive scatter plot is not just a static representation but allows users to explore and navigate the data themselves. A user can easily hover over multiple products in the same cluster to identify similar items, compare their details, and make informed decisions. The hover tool also helps users understand why certain products might be clustered together by showing how their price points or brands compare.
- Customization of Visuals: The scatter plot's appearance was further customized for usability. The size and color of the points were adjusted to enhance visibility and distinguish overlapping data points. For example, products from different brands were color-coded, or those in different price ranges were represented with varying point sizes. Such visual cues further enrich the user's ability to intuitively grasp complex relationships.



• Key Insights:

• Understanding Product Similarities:

• The interactive plot helps users identify products with similar ingredient compositions. This makes it easier to explore alternatives that may differ in brand or price but offer comparable formulations.

Revealing Ingredient Trends:

• The visualization uncovers patterns such as luxury and budget-friendly products sharing similar ingredients. This can provide valuable insights into how pricing strategies and marketing may be influenced more by branding than formulation.

• Exploring Ingredient-Based Groupings:

The scatter plot reveals natural groupings of products based on their ingredients. For
instance, products designed for sensitive skin often cluster together due to common
soothing components like aloe vera or colloidal oatmeal, allowing users to infer product
intent based on shared formulations.

4.3 Comparing Similar Cosmetic Products

After mapping the products based on their ingredient similarity and visualizing their relationships, the next step in the system's functionality is comparing specific products based on their ingredients. This was achieved using cosine similarity, which measures the similarity between two ingredient lists.

- Cosine Similarity Calculation: Cosine similarity is a metric that calculates the cosine of the angle between two vectors. In this case, the vectors represent the presence or absence of ingredients in two different products. The cosine similarity score ranges from 0 to 1, where 1 means the products have identical ingredients, and 0 means they have no ingredients in common. This metric allows us to numerically quantify how similar two products are based on their ingredients.
- Top N Similar Products: For each product in the dataset, we computed its cosine similarity score with every other product, ranking them based on the similarity scores. This allowed us to generate a list of the top N most similar products for any given item. The list provides users with highly relevant alternatives that they might consider if they are interested in finding similar products.



V. CONCLUSION

Conclusion:

The project undertaken to develop a content-based recommendation system for cosmetic products marks a significant step forward in how ingredient-conscious consumers can explore, compare, and discover alternative products. By focusing specifically on moisturizers for dry skin, we have built a robust framework that can easily be expanded to other cosmetic categories. Through the careful application of natural language processing (NLP), machine learning, and data visualization techniques, the system provides personalized, ingredient-based recommendations—addressing a growing demand in the beauty industry for transparency and specificity.

Problem Addressed: One of the most pressing concerns for consumers today, especially in the skincare and cosmetics domain, is understanding what they are putting on their skin. With so many products in the market containing complex, and often unfamiliar, ingredient lists, consumers find it challenging to make informed choices. Often, they are swayed by marketing claims, pricing, or brand loyalty without understanding the actual efficacy or suitability of a product for their specific skin needs.

- Lack of Ingredient Knowledge: Consumers are increasingly aware of the importance of ingredients but lack the tools to easily compare products at a granular level based on ingredient composition. Many products share a significant overlap in their formulations, but without accessible ingredient data and comparison tools, it becomes difficult to identify alternatives that may offer similar benefits at a lower cost or with fewer irritants.
- Customization of Skincare Needs: Another major challenge is personalization. Each individual's skin has unique characteristics—some have specific needs like avoiding allergens, while others are seeking targeted treatments like anti-aging or hydration. There is no universal solution in skincare, and consumers must sift through countless products to find one that meets their requirements.
- This project directly addresses these challenges by building a data-driven system that uses the ingredients as the basis for comparison.

VI. FUTURE SCOPE

Future Scope:

This system has significant potential for extension and improvement, including:

- User Ratings and Reviews: Integrating user feedback to refine the recommendation process further. Incorporating ratings and reviews can enhance the system's accuracy in recommending products that not only have similar ingredients but are also highly rated by consumers.
- Support for Additional Product Categories: Expanding the system to include a broader range of products, such as cleansers, sunscreens, and serums. This would provide a comprehensive resource for consumers looking to evaluate various skincare products based on their ingredient compositions.
- Ingredient Toxicity Analysis: Implementing analyses to flag harmful or controversial components in cosmetic products. By evaluating the safety of ingredients, the system could guide users in avoiding potentially harmful substances and promoting safer skincare choices.
- Enhanced User Personalization: Developing features that allow users to input specific skin concerns or preferences (e.g., sensitivity, allergy avoidance) could lead to more tailored recommendations. This level of personalization would significantly enhance user satisfaction and trust in the recommendations provided.
- Collaborative Filtering Techniques: Exploring the integration of collaborative filtering methods to recommend products based on user behavior and preferences, in addition to content-based recommendations.



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