Mining Potential Chemical Co-exposures from Consumer Product Purchasing and Ingredient Data

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The views expressed in this presentation are those of the authors and do not necessarily reflect the views or policies of the U.S. EPA
• Addressing risks associated with chemical mixtures is a challenge
  • Too many chemicals and too many co-exposures
• EPA’s ToxCast Program has screened thousands of chemicals for bioactivity in high-throughput \textit{in vitro} assays
• HTS and mixtures
  1. Predict activity from component chemical responses using modeling
  2. Test whole mixtures (can inform #1)
• But which mixtures to test?
• \textit{In ExpoCast, we are developing tools that allow us to identify relevant chemicals with potential real-world co-exposures}
Approaches for Identifying Co-Exposures

• **Modeling approaches**
  - Multiple sources, pathways, and routes of exposure can be considered
  - Uncertainties associated with estimating external versus internal exposure (dose) – timing of exposures and consideration of absorption, distribution, metabolism, and excretion (ADME) processes
  - Impacted by data gaps in behavior (e.g., consumer habits and practices), source information (e.g., chemical use or ingredient data), or toxicokinetics

• **Biomonitoring**
  - Can identify both parent chemicals and metabolites in blood or urine
  - Aggregate over time (e.g., bioaccumulating compounds)
  - Limited number of chemicals (expensive, need standard analytical methods)
Kapraun et al. (2017) mined biomonitoring data from the NHANES study to identify prevalent chemical combinations.

- Measured concentrations were discretized to presence/absence using a fixed threshold.
- Examined co-occurrence within three groups of chemicals measured in unique subsamples of the study population, using frequent itemset mining (FIM).
- Identified 90 chemical combinations consisting of relatively few chemicals that occur in at least 30% of the U.S. population.
- Identified three “supercombinations” of chemicals that occurred in a smaller fraction of the population.

Kapraun et al., Environmental Health Perspectives, 125:8, 2017
Current Approach

- Integrate large datasets of consumer product ingredient and product purchasing information to develop a dataset that can be mined for chemical co-exposures

- Apply FIM to identify prevalent co-occurring chemicals within household-months

- Stratified results by household demographics to characterize variability in co-exposure patterns and identify potential chemical combinations associated with sensitive populations, such as families with young children and women of childbearing age
EPA-ORD’s Chemicals and Products Database (CPDat)

• EPA ORD database containing curated chemical use and consumer product ingredient data

• Public version of the dataset contains ingredient data for over 60,000 products, mapped to standardized product categories for use in exposure assessment and modeling

• Also recently extracted ingredient data from 230,407 retailer-provided product safety data sheets (SDSs), including product name, category, universal product code (UPC), and chemical identifiers

https://www.epa.gov/chemical-research/chemical-and-products-database-cpdat
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- Chemical identifiers curated to harmonized EPA Distributed Structure-Searchable Toxicity (DSSTox) Substance Identifiers (DTXSIDs)

Consumer Product Purchasing Study

• EPA initiated a collaboration with Nielsen in 2013
• Shared data from the National Consumer Panel (NCP)
• Formerly called “Homescan” project
Purchasing Data

• 60,000 U.S. households for 1 year (2012)

• Demographic information for each household
  • Income, number of household members, Nielsen market (metro area), county size, race, presence and age of children, age and occupation of female head of household

• All purchases for product categories of interest to Nielsen
  • 29 broad categories called “Groups” (e.g., Household Cleaners, Cosmetics, Fresheners and Deodorizers).
  • Date of purchase, UPC, brand, number of units, size
  • ~4.6 million individual product purchase records

• 133,966 unique product UPCs

• Recent publication: Tornero-Velez et. al (2020) examined product co-purchases which gave us some idea about chemical co-exposure from previous ingredient data; the ability to link individual purchases to specific chemicals is a major step forward.

Tornero-Velez et al., Risk Analysis, 125:8, 2017
## Data Integration

### Product-Chemical Data

<table>
<thead>
<tr>
<th>UPC</th>
<th>Chemicals</th>
</tr>
</thead>
<tbody>
<tr>
<td>UPC1</td>
<td>DTXSID1, DTXSID2, DTXSID4</td>
</tr>
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</tr>
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### Purchasing Data

<table>
<thead>
<tr>
<th>Date</th>
<th>Household (HHLD)</th>
<th>UPC (12 digits)</th>
<th>Product Variables</th>
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<td>2012-01-01</td>
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<tr>
<td>2012-02-09</td>
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#### Monthly Transaction Matrix

<table>
<thead>
<tr>
<th>HHLD-month</th>
<th>Chemicals</th>
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<tbody>
<tr>
<td>HHLD1-01</td>
<td>DTXSID1, DTXSID2, DTXSID4, DTXSID5, DTXSID6</td>
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<tr>
<td>HHLD2-02</td>
<td>DTXSID2, DTXSID3</td>
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Direct and Fuzzy Matching by UPC

Could match ~50.3% purchases
Data Integration

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<tr>
<td>...</td>
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</table>

### Data Summary

<table>
<thead>
<tr>
<th>Data</th>
<th>Count</th>
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<tbody>
<tr>
<td>Transactions</td>
<td>539,857</td>
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<tr>
<td>Households</td>
<td>53,525</td>
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<tr>
<td>Products</td>
<td>31,375</td>
</tr>
<tr>
<td>Chemicals</td>
<td>783</td>
</tr>
</tbody>
</table>
Chemical Lists

- Analysis of co-occurring chemicals was restricted to chemicals of regulatory or biological interest in order to avoid identification of prevalent chemical combinations containing common substances having little relevance to risk assessment (e.g., water)

- **Broad Chemical List:** Active public chemical inventory of the Toxic Substances Control Act (TSCA)
  - 649 chemicals in the consumer product transaction dataset
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- **Broad Chemical List**: Active public chemical inventory of the Toxic Substances Control Act (TSCA)
  - 649 chemicals in the consumer product transaction dataset

- **Case-Study: Potential Endocrine Active Chemicals (EACs)**

<table>
<thead>
<tr>
<th>Source</th>
<th>Investigated Biological Action</th>
<th>Chemicals Predicted to be Active</th>
<th>Chemicals Mapped to Purchased Products</th>
</tr>
</thead>
<tbody>
<tr>
<td>Collaborative Estrogen Receptor Activity Prediction Project (CERAPP)¹</td>
<td>Estrogen Disruptors</td>
<td>1,142</td>
<td>10</td>
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<tr>
<td>Collaborative Modeling Project for Androgen Receptor Activity (COMPARA)²</td>
<td>Androgen Disruptors</td>
<td>16,112</td>
<td>42</td>
</tr>
<tr>
<td>Additional potential EACs from Literature Sources³</td>
<td>Multiple</td>
<td></td>
<td>17</td>
</tr>
<tr>
<td><strong>Total (unique)</strong></td>
<td></td>
<td></td>
<td><strong>65</strong></td>
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</tbody>
</table>

³Dodson et. al. Environmental health perspectives. 120:935-943.
**Frequent Itemset Mining**

- **Itemset**
  - A collection of one or more items
  - Example: \{**DTXSID1**, **DTXSID4**, **DTXSID5**\}
  - \(k\)-itemset
    - An itemset that contains \(k\) items

- **Relative support/prevalence (**\(\sigma**)****
  - Fraction of transactions that contain an itemset
  - E.g. \(\sigma(\{**DTXSID1**, **DTXSID4**, **DTXSID5**\}) = 2/5\)

- **Frequent itemset**
  - An itemset whose prevalence is greater than or equal to a *minimum support* threshold

Apply to transaction data to identify prevalent combinations

<table>
<thead>
<tr>
<th>Transaction ID</th>
<th>Items</th>
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<tbody>
<tr>
<td>HHLD-01</td>
<td><strong>DTXSID1</strong>, <strong>DTXSID4</strong></td>
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<td>HHLD-02</td>
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</tbody>
</table>
Frequent Itemset Mining Analyses

• Performed using the *ECLAT* (Equivalence Class Clustering and bottom-up Lattice Traversal) function of the *Arules* R package

• Performed identification of prevalent individual chemicals and combinations
  • For TSCA chemicals and EACs, based on a threshold prevalence for chemical group that provided a manageable number of itemsets
  • Within product groups
  • Within demographics, including:
    • Women of childbearing age
    • Different income ranges
    • Race of female head of household
    • Education level
    • Different family sizes/ages of children

• Interpreted chemicals within prevalent combinations by examining chemical functions
  • Harmonized functional uses defined by Phillips et al. (2017).
    • Dataset of 14,000+ reported chemical-function pairs

Results: Orientation

<table>
<thead>
<tr>
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<th>0</th>
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Results: Orientation

Demographic

Deviation of the rank of the chemical or combination for each demographic from the rank in the population overall

- Elevated rank relative to overall population
- Potentially higher exposure for the demographic
- Reduced rank relative to overall population
- Potentially lower exposure for the demographic
Results: Orientation

Demographic

{Chemical or Chemical Combination}

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- Elevated rank relative to overall population
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- Reduced rank relative to overall population
  - Potentially lower exposure for the demographic
Results: Orientation

Chemical Function Information

Receptor Information

Deviation of the rank of the chemical or combination for each demographic from the rank in the population overall

Elevated rank relative to overall population

Potentially higher exposure for the demographic

Reduced rank relative to overall population

Potentially lower exposure for the demographic
### Most Prevalent Single Chemicals

**Group 1 (Broad TSCA Inventory)**

- 20 overall most prevalent individual chemicals
- Top 5 chemicals are what were termed “ubiquitous function” chemicals - perform a variety of functions in products
Most Prevalent Single Chemicals

Group 1 (Broad TSCA Inventory)

- 20 overall most prevalent individual chemicals
- Top 5 chemicals are what were termed “ubiquitous function” chemicals - perform a variety of functions in products
- Differences by demographic can be observed

e.g., titanium dioxide has a higher prevalence rank for Asian female head of household

Two common cleaning product ingredients had reduced prevalence houses where the female head had post-college education
Most Prevalent Single Chemicals

**Endocrine Active Chemicals**

- Many of the most prevalent EAC chemicals were fragrances (or categorized as such due to presence in fragrance formulations)
- Many of these chemicals were present in a variety of personal care products
Most Prevalent Single Chemicals

**Endocrine Active Chemicals**

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- Many of these chemicals were present in a variety of personal care products
- **Benzethonium chloride** and **diazolidinyl urea**, which were ranked 2 or 3 places higher in households with children, are commonly used as topical antimicrobial agents in baby wipes, bubble baths, cosmetics, and skin care products
**Most Prevalent Single Chemicals**

**Endocrine Active Chemicals**

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- Many of these chemicals were present in a variety of personal care products.
- Benzethonium chloride and diazolidinyl urea, which were ranked 2 or 3 places higher in households with children, are commonly used as topical antimicrobial agents in baby wipes, bubble baths, cosmetics, and skin care products.
- Households with children under 6 have a higher ranking for quaternary ammonium compounds, di-c14-18-alkyldimethyl, me sulfates, which are commonly used in disinfectants and hand soaps.
Prevalent Combinations

Group 1 (Broad TSCA Inventory)

- Here demographics and chemical sets are clustered to indicate the similarity of rankings of chemical combinations
- Set A: ubiquitous consumer product chemicals present in households with children, higher income, and more highly educated, representing generally high consumer product use
- Set B: elevated difference in rank in lower to middle income demographics and African American households, and reduced rank differences in Asian households and females with post-college education and females of childbearing age; these three sets contained antimicrobials and surfactants found in cleaning products

Minimum prevalence= 2.5% HHLD-Months
Prevalent Combinations

**Endocrine Active Chemicals**

- One itemset \(\{dl\)-tocopherol mixture | phytonadione\}, contained two chemicals that targeted the same receptor (AR).

Minimum prevalence= 0.1% HHLD-Months
Prevalent Combinations

- One itemset \{dl-tocopherol mixture | phytonadione\}, contained two chemicals that targeted the same receptor (AR).

- The highest positive rank departure for households with children occurred for the itemset \{decamethylcyclopentasiloxane | limonene\}.

Endocrine Active Chemicals

- Minimum prevalence = 0.1% HHLD-Months
Prevalent Combinations

Endocrine Active Chemicals

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- Households with a female head of Asian race had the highest positive rank departure for the combination of limonene and linalool, the latter of which is used as a scent in many perfumed hygiene products and cleaning agents.

Minimum prevalence = 0.1% HHLD-Months
Prevalent Combinations

- One itemset \{\textit{dl}-tocopherol mixture \mid \textit{phytonadione}\}, contained two chemicals that targeted the same receptor (AR).

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- Households with a female head of Asian race had the highest positive rank departure for the combination of \textit{limonene} and \textit{linalool}, the latter of which is used as a scent in many perfumed hygiene products and cleaning agents.

- African American households had a positive rank departure of 6 for the combination \{\textit{linalool} \mid 2-phenylethanol\}; the second chemical is a floral fragrance primarily present in air fresheners.

Minimum prevalence= 0.1% HHLD-Months
• Collectively across all products and by product group, results indicated that households with children, households headed by women of color, and lower income households exhibited divergence from the general population in the chemical combinations they encounter most frequently.

  • This may be due to a need for different types of personal care products designed specifically for given races or ethnicities, brand or regional preferences, or simply the need for a wider variety of products in households with multiple children.

  • These patterns may reflect differential experiences and thus differential exposures among demographics.

• Lists of most prevalent combinations (overall and for various demographics) can be evaluated for feasibility for testing in \textit{in vitro} assay systems, and further prioritized based on single-chemical activity or exposure-related factors.

• New non-targeted analysis (NTA) studies of biological media such as blood or urine can complement and evaluate predictions of co-exposures associated with consumer products.

  • Such studies also have the potential for identifying mixtures containing metabolites of consumer product chemicals.
Testing Chemical Mixtures

• Use existing ToxCast data and amenability to screening as pilot information for mixture design

• Construct chemical mixtures (each with 3 chemicals) and screen each constituent

• General hypothesis: the point-of-departure for mixture bioactivity can be predicted from the concentration response data for the chemical constituents using models of concentration addition, independent action, or a joint model
Humans are exposed to thousands of chemicals from the products they purchase and use within the household.

Assessing every possible set of chemicals for toxicity is an impossible task but also an unnecessary one as the number of chemical mixtures that are prevalent and occur in real-world scenarios is drastically less.

We have presented here a novel approach that applies FIM on a dataset describing the chemicals entering households through purchased consumer products to identify a manageable number of chemical combinations that regularly occur in homes across the US.

These identified combinations can inform the prioritization of chemical combinations for toxicity testing.
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CPHEA
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*Trainees

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Silent Spring Institute
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Southwest Research Institute
Alice Yau
Kristin Favela
Summit Toxicology
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