

Air Toxics Data Analysis Madeleine Strum EPA/OAQPS Air Quality Assessment Division

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Uses of Ambient Data of Air Toxics

- Confirm risk model results
 - Assess risk predicted by model in key locations (i.e., schools, communities)
- Model Evaluation
- Trends
 - Tract progress of air toxic reduction efforts for key pollutants (i.e., benzene, formaldehyde)
- Local assessments
 - Near roadway, fenceline monitoring downwind from industrial facilities

http://www.epa.gov/national-air-toxics-assessment

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National Air Toxics Assessment



On December 17, 2015, EPA released the most recent update to the National Air Toxics Assessment (NATA). NATA contains emissions data from 2011 and uses models to make broad estimates of health risks over geographic areas of the country.

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Learn more

NATA Overview

- Limitations
- Glossary of Terms
- Frequent Questions



- 2011 Assessment Results
- <u>2011 NATA Map</u>
- 2011 Assessment Methods

Quick Links

- Previous versions of NATA
- Other environmental screening tools
- Learn about risk assessment
- <u>Hazardous Air Pollutants website</u>
- <u>Urban Air Toxics website</u>

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Monitoring Data Included in the NATA Map App





Monitoring Data Included in the NATA Map App

		- /
AMA SITE CODE	482010036	
Location	NA	
STATE	TX	
Latitude	29.776100	
Longitude	-95.105103	
Location Type	SUBURBAN	
Setting	NA	
NATA HAP Name	BENZENE	
AQS PARAMETER CODE	45201	
AQS PARAMETER NAME	Benzene	
Pollutant Code Description	Benzene	
Duration Description	24 HOURS	
2011 Max Conc (µg/m3)	10.670840	÷.,
2011 Mara Cara (0.037540	



Monitoring Data Included in the NATA Map App





Air Toxics Challenges

- What is level of completeness to require to creating annual average/statistics: (11 measurements vs 12 per quarter)
- Differences in measurements across POCS (monitors) at same monitoring site
 - Average across all POCS or pick a POC?
- Treatment of data below MDL and ND
 - cancer benchmarks below MDL



Pollutants where MDL > 1 in 1-million cancer benchmark



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Data analysis challenges: data below MDL and non detects

- MDL statistical construct, estimate of the concentration at which there is 99% confidence that the analyte, when positively identified, is present
- ND analyte not identified based on set criteria such as signal to noise
 - Does not mean the concentration is zero
 - We see NDs where remote concentration (background) is nonzero

Don't see these distinguished in literature on handling data below detect



What does below MDL/ND mean: not there? vs can't measure it?

	REMOTE (ug/m3)	Range of Site Minimum MDL 2014	Sites with NDs In 2014
Carbon tetrachloride	0.547	0.006 to 0.53	6 sties
Chloroform	0.058	0.01 to 0.4	12 sites
Benzene	0.116	0.006 to 0.27	5 sites
Methyl chloroform	0.06	0.01 to 0.46	7 sites
Methyl bromide	0.0294	0.02 to 0.4	16 sites
Dichloromethane (methylene chloride)	0.146	0.01 to 7	6 sites
Trichloroethylene	0.0041	0.01 -0.46	3 sites
Tetrachloroethylene	0.0131	0.01-0.6	10 sites



ND versus below MDL



Figure 2. Analyte Concentration vs. Reporting Certainty. (The y axis represents signal strength, in units of the standard deviation (σ) used to determine the MDL.) Adapted from Keith, 1991 by Johnson, 2001.



What to do with data below MDL/Nondetect

- When using censored data: Helsel (Chemosphere, 2006) – substituting is fabricating. Use statistical techniques.
- When reporting the data: many scientists including EPA's Science Advisory Board; Analytical Methods Committee, Royal Society of Chemistry: Report the result found, with statement of its uncertainty



Sensitivities to different treatment at individual sites

- Ethylene dichloride
- 1,3 butadiene

First, we look at MDLs for across the NATTS sites



Ethylene Dichloride MDLs – 2014





1,3 Butadiene MDLs – 2014





Can treat below MDL and ND many different ways

- Substitution approaches
 - Data below MDL = $\frac{1}{2}$ MDL
 - Data below MDL = 0
 - Use data as-is. No change to data below MDL; ND=0
 - Data below MDL=1/2 MDL; ND=0
- Statistical Approaches
 - Use Regression on Order statistics (ROS) for ND
 - Use Kaplan-Meir (KM) for ND
 - Use ROS for all data below MDL
 - Use K-M for all data below MDL



Ethylene dichloride- sensitivity to substitutions Treat <MDL; ND as the same





Ethylene dichloride- sensitivity to substitutions Treat <MDL; ND differently





Ethylene dichloride– sensitivity to substitutions/statistical approaches









1,3 butadiene – sensitivity to approaches for below MDL and ND





Sensitivities to different treatments - NATTS wide, pollutant wide – 2005 and later

RATIO OF MEAN: $(\frac{1}{2} \text{ MDL}; \text{ ND}= 0) / (\text{ND}=\text{ROS})$

- Looked at 2005 and later
- sites with between 20% and 80% data below MDL





Sensitivities to different treatments - NATTS wide, pollutant wide – 2014





Analysis of data below MDL

- Just beginning –lots of questions- e.g.,
 - distinguish between ND and MDL?
 - What is the percent of data below MDL above which the site should not be used for model evaluation or trends?
- Deciding what to do with values below MDL should not be datareporting dependent; it should be done at the analysis stage not the reporting/recording stage
- Once values are removed or replaced, we cannot get them back
 - Can utilize values to learn about uncertainty/distribution of data below MDL if you have the actual values
 - these data could help us determine best practices for data analysis
- Substituting a value (0, MDL/2, MDL) will bias the result
- Foster consistency