

Development of a Preliminary Emission Inventory of Toxic Air Contaminants (TACs) for the Bay Area

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Introduction

Purpose of the TAC Emission Inventory:

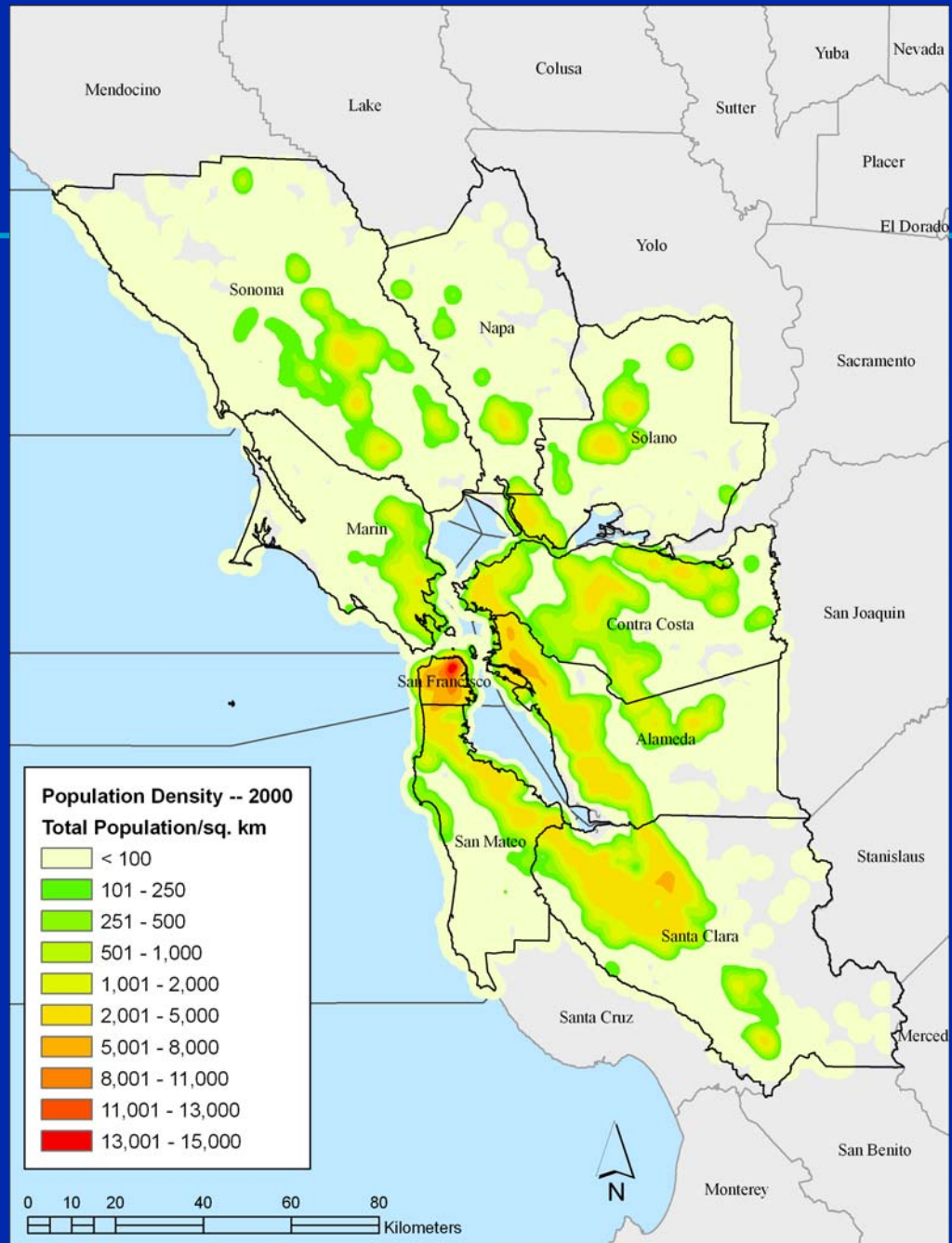
Support screening-level assessments of population exposures and selection of a study community.

Objective: Develop a first-draft emission inventory for the Bay Area using existing information suitable for use with dispersion and exposure models.

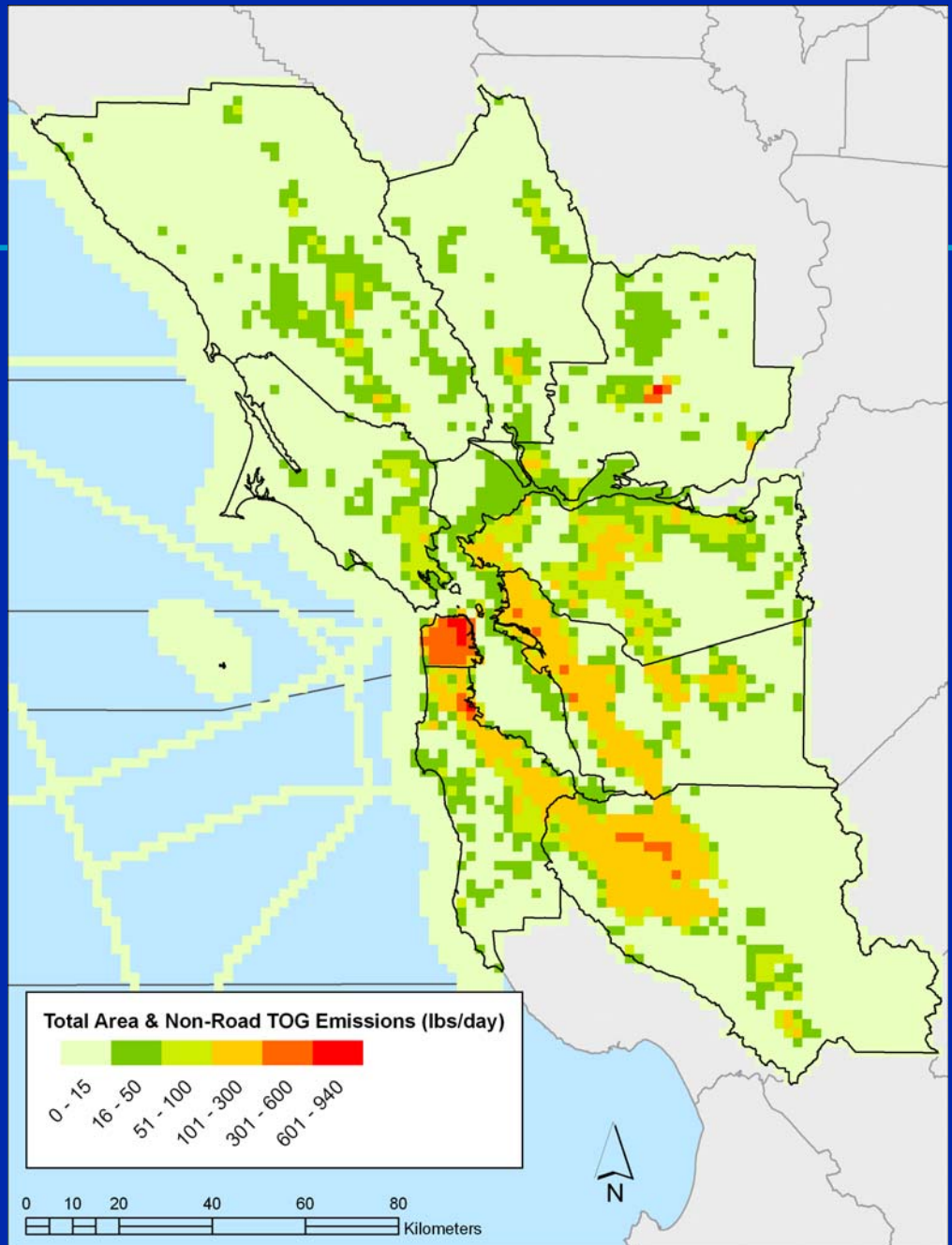
General Approach

1. Begin with existing criteria pollutant inventories.
2. Apply available chemical speciation profiles.
3. Apply available cancer and non-cancer unit risk factors.
4. Spatially allocate emissions.

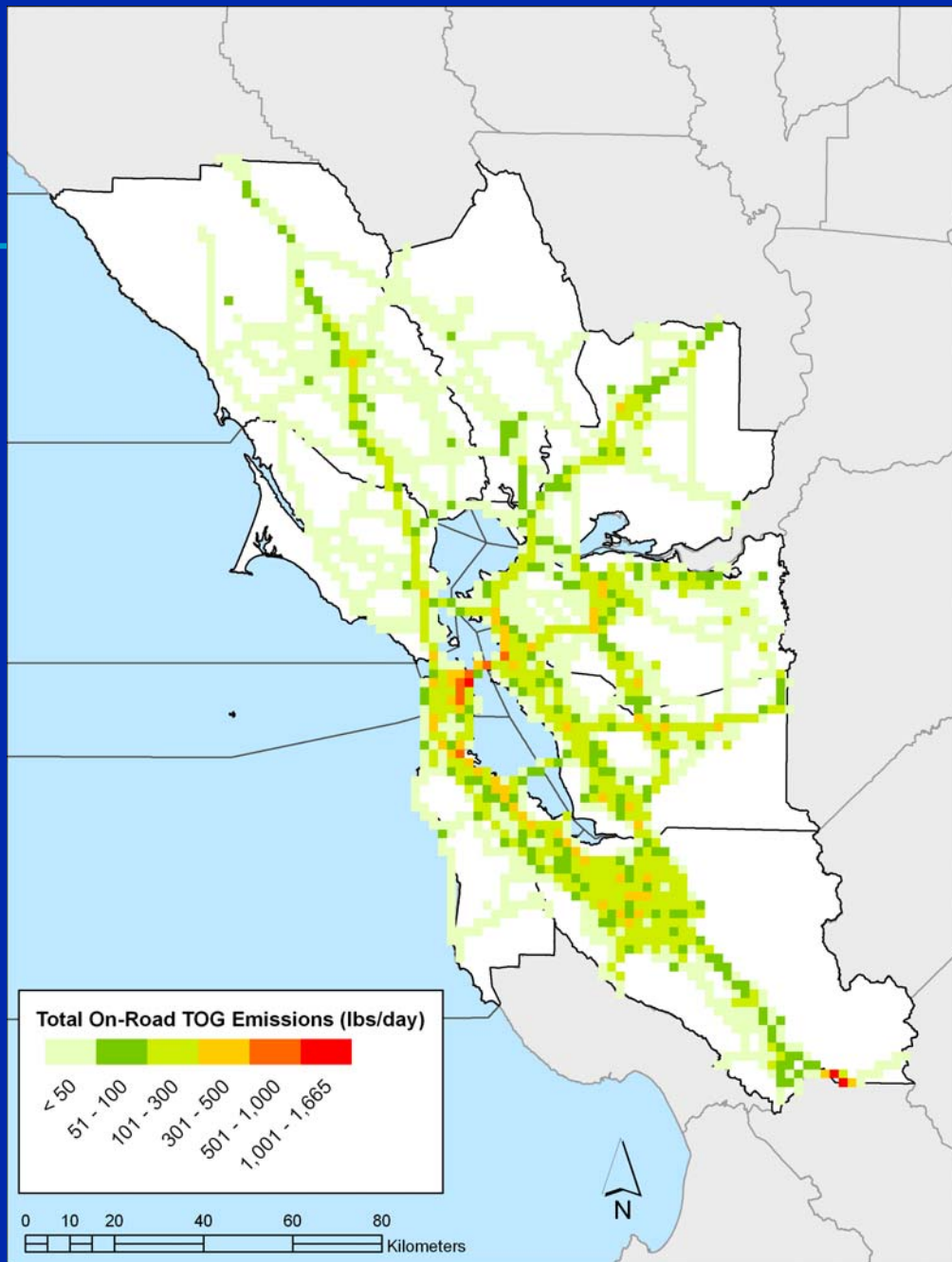
Bay Area Population Density (2000)



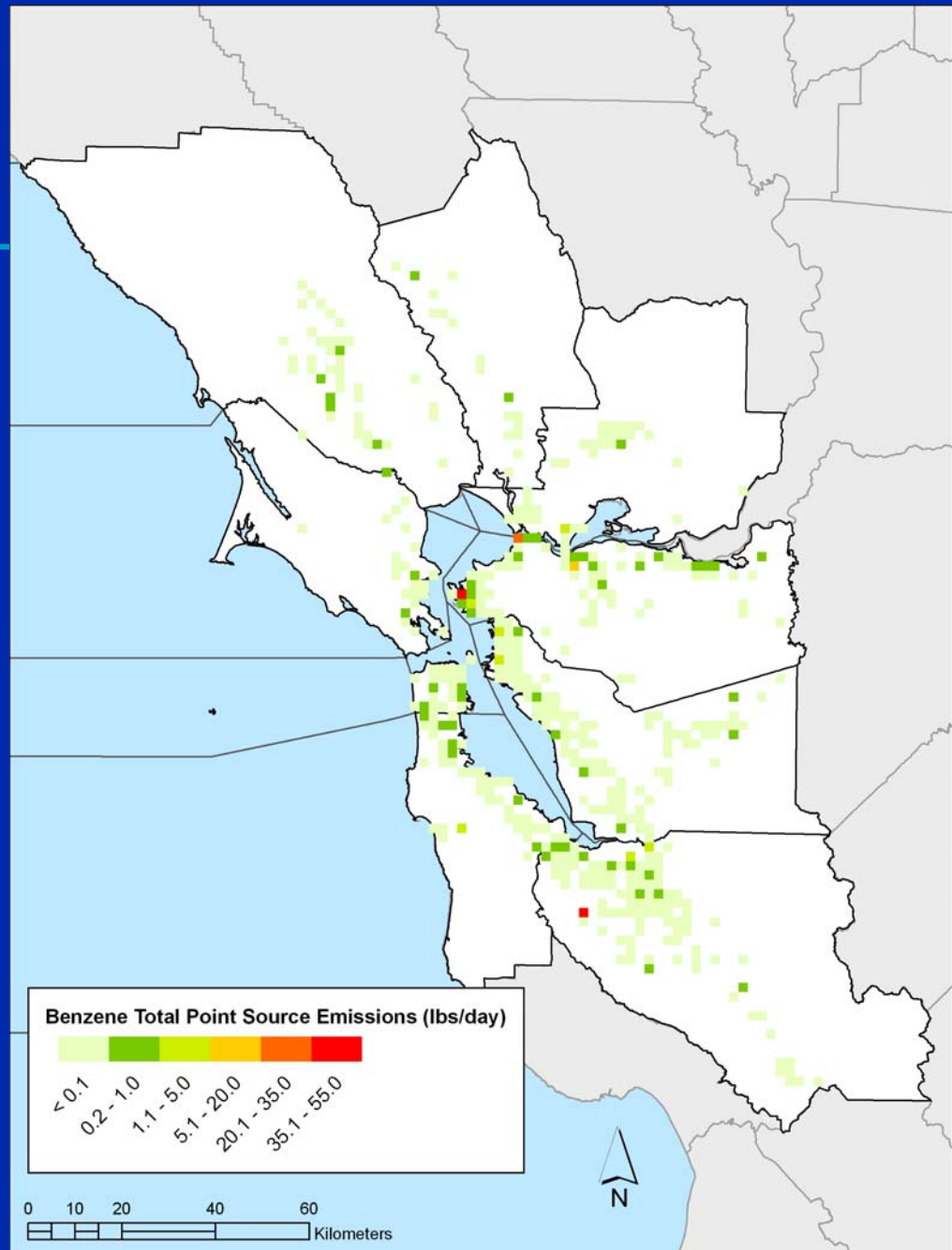
2000 Area & Non-Road TOG Emissions



2000 On-Road TOG Emissions



2000 Benzene Emissions from Point Sources



Chemical Speciation

Existing speciation profile libraries

- ARB
- EPA (Speciate 3.2)
- Desert Research Institute (DRI)

Chemical Speciation

CAS No.	Pollutant	Percent
67641	ACETONE	2.14
123864	N-BUTYL ACETATE	2.6
71363	N-BUTYL ALCOHOL	1.82
	C-9 CYCLOPARAFFINS	7.3
	C10 PARAFFINS	16.78
	C4 SUBSTITUTED CYCLOHEXANE	17.04
	C7-C16 PARAFFINS	2.82
	C9 PARAFFIN	5.32
141786	ETHYL ACETATE	0.74
64175	ETHYL ALCOHOL	2.44
67630	ISOPROPYL ALCOHOL	2.84
67561	METHYL ALCOHOL	0.08
78933	METHYL ETHYL KETONE	4.16
108101	METHYL ISOBUTYL KETONE	1.4
	PARAFFINS/OLEFINS (C12-C16)	6.7
109604	N-PROPYL ACETATE	0.18
108883	TOLUENE	6.7
71556	1,1,1-TRICHLOROETHANE	1.76
	ISOMERS OF UNDECANE	9.62
	UNIDENTIFIED	4.44
1330207	ISOMERS OF XYLENE	3.12

EPA TOG
Profile #6003:
Architectural
Coatings

Chemical Speciation

Cross-Reference Table

EIC Code	Description	Profile#
52052091050000	Primers & Sealers	6003
53053032250000	Pesticides	1000
54059004000000	Asphalt Roofing	2400

Speciation Profile Library

Profile#	CAS#	Pollutant	Percent
6003	67641	ACETONE	2.15
6003	123864	N-BUTYL ACETATE	2.60
6003	71363	N-BUTYL ALCOHOL	1.82

Unit Risk Factors



Cancer risk is approximated as a linear function of ambient concentration

URF Units of Measure =

No. Cases of Effects per Hundred Thousand
Persons Exposed \div Ambient Concentration

Example – Benzene

- Ambient concentration = $2 \mu\text{g}/\text{m}^3$
- $\text{URF} = 2.9 \times 10^{-5} \cdot 10^{-5} \cdot (\mu\text{g}/\text{m}^3)^{-1}$
- Risk = concentration \times URF
= 5.8×10^{-10} or “58 in 100 billion”

Reference Concentrations



Reference Concentration (RfC) is a regulatory definition.

RfC = Concentration at which no adverse effects are expected,

plus a safety margin for uncertainty,

plus a safety margin based on expert toxicologists' judgment.

RfC Units of Measure = Concentration

Example – Benzene

- Ambient concentration = $2 \mu\text{g}/\text{m}^3$
- RfC = $60 \mu\text{g}/\text{m}^3$ for chronic effects and $1300 \mu\text{g}/\text{m}^3$ for acute effects
- Risk = negligible at the observed concentration

Sources of Information

Sources of Inhalation URFs and RfCs

- ARB-Approved Risk Assessment Health Values
- EPA Office of Environmental Health Hazard Assessment (OEHHA)
- EPA Integrated Risk Information System (IRIS)
- Risk Assessment Information System

Risk-weighted Emissions

- Best available URFs and RfCs were selected for each TAC in the speciated inventories.
- Risk-weighted emissions were calculated.
- Uncertainty ranges were documented where applicable.
- A database of risk-weighted emissions by TAC and source category was prepared.

Risk-weighted Emissions (“RWE”)

$$\text{Cancer RWE}_i = E_i \times \text{UR}_i \div \text{UR}_X$$

$$\text{Chronic or Acute RWE}_i = E_i \div \text{RfC}_i \times \text{RfC}_Y$$

i denotes TAC species *i*

E = Mass-based emissions; lbs/day

UR = Unit risk factor; $10^{-5} \cdot (\mu\text{g}/\text{m}^3)^{-1}$

RfC = Reference concentration; $\mu\text{g}/\text{m}^3$

UR_X = UR for hypothetical species “X”
 $\equiv 1 \cdot 10^{-5} \cdot (\mu\text{g}/\text{m}^3)^{-1}$ (similar to dioxins or furans,
which are extremely toxic)

RfC_Y = RfC for hypothetical species “Y”
 $\equiv 1 \mu\text{g}/\text{m}^3$ (similar to aniline or acrylic acid, which
are moderately toxic)

Spatial Allocation (1 of 2)

Area and Non-road Sources —

Apply spatial allocation factors (SAFs)

On-road Mobile Sources —

Already resolved to 2-km x 2-km grid

Point Sources —

Assign to facility location coordinates

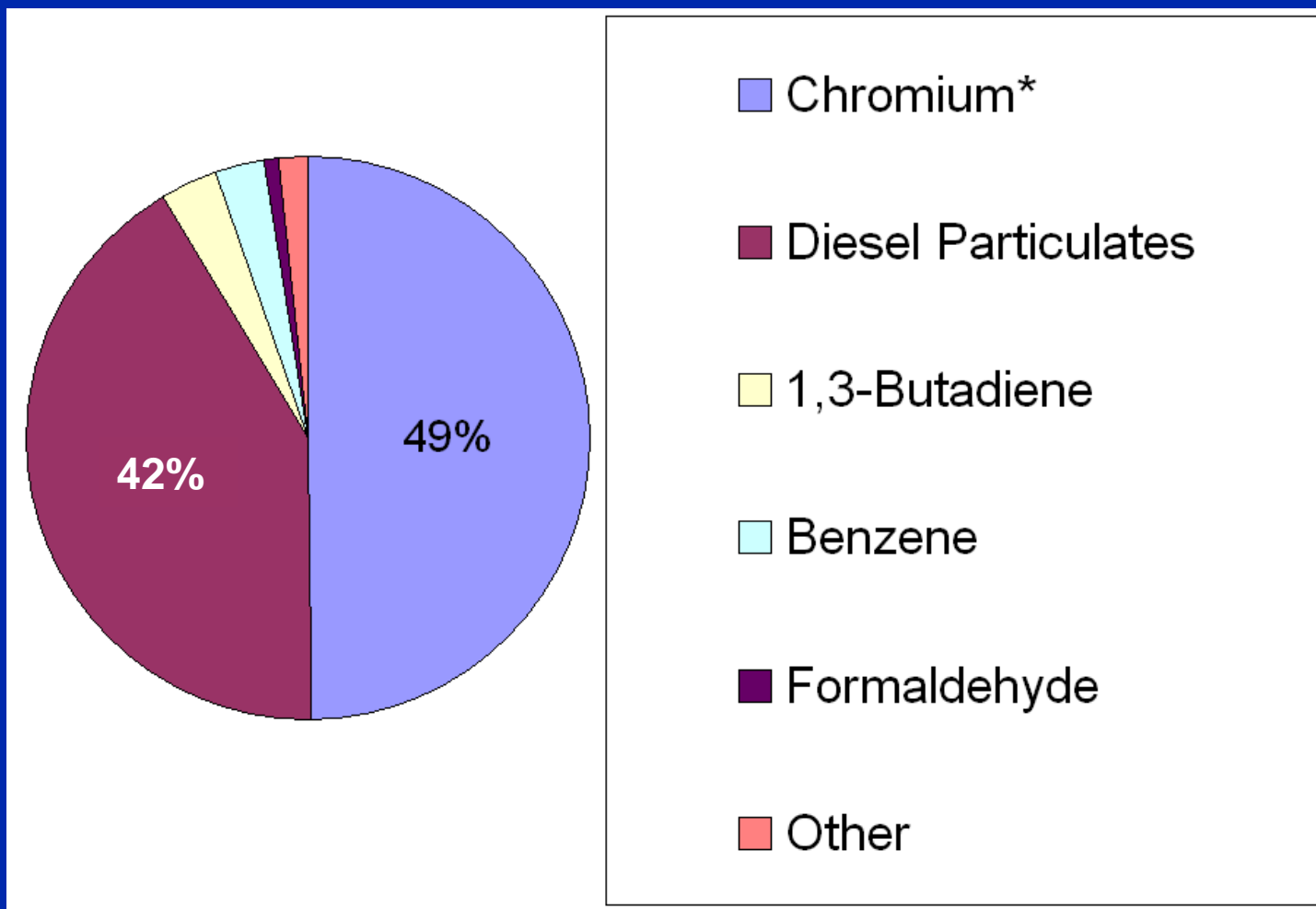
Spatial Allocation (2 of 2)

SAFs were developed from geographic information systems (GIS) databases.

- Demographic data (e.g., population density)
- Landuse/landcover data sets (e.g., residential versus agricultural land use)
- Line length (railroad tracks)
- Facility locations
- Other available GIS databases

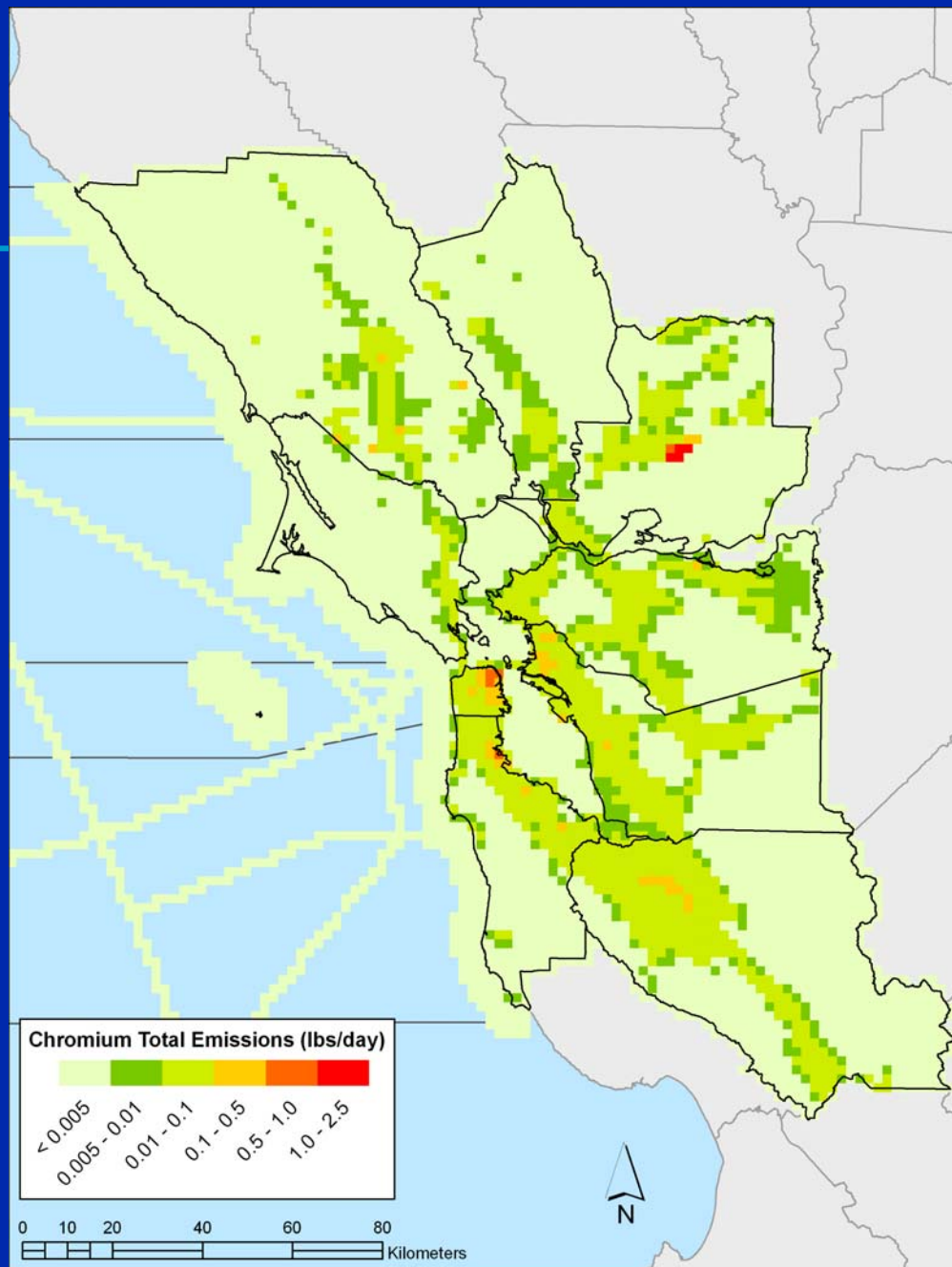
Cancer Risk-weighted Emissions

16 equivalent pounds/day

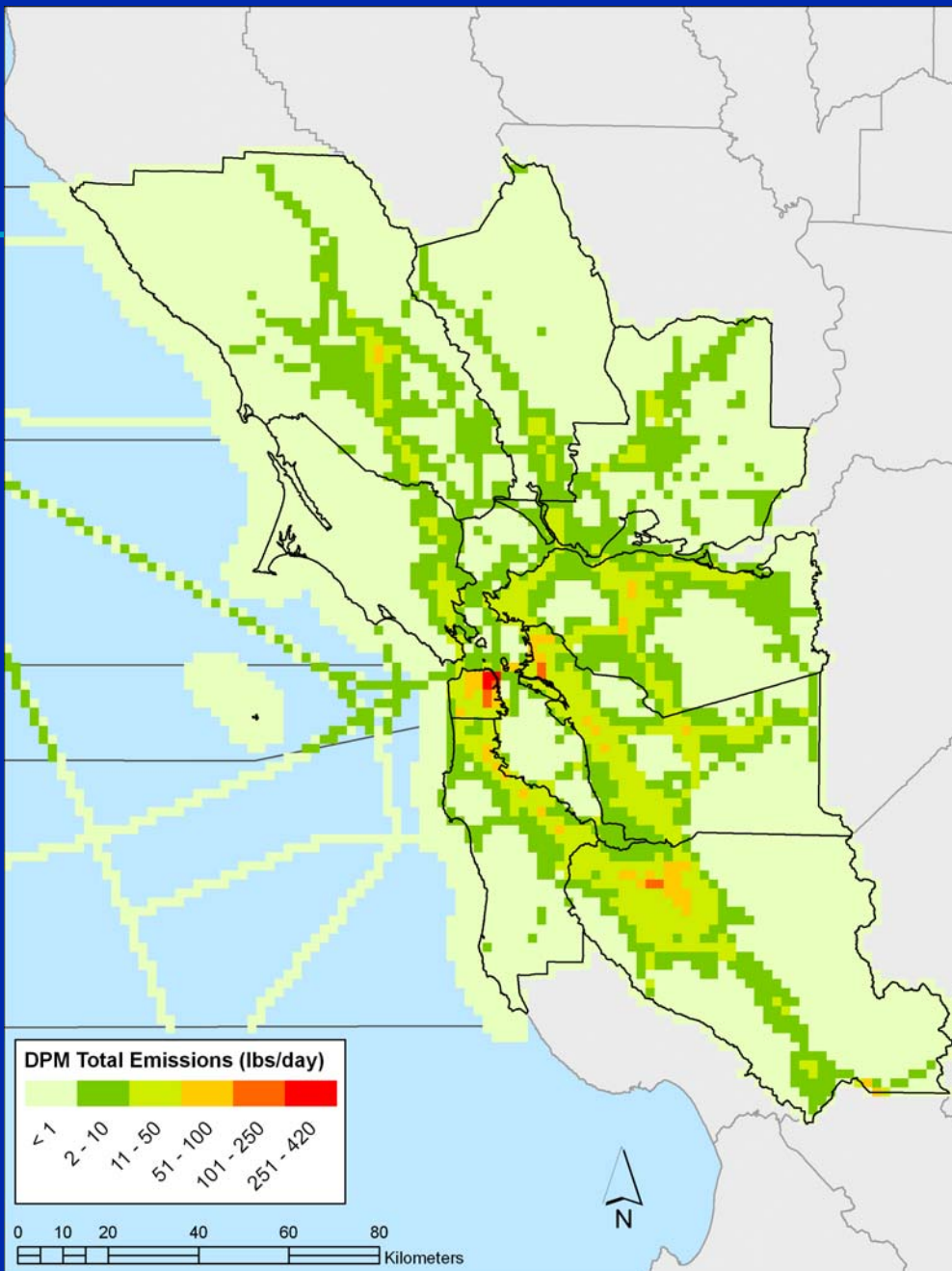


* denotes TACs with large uncertainties

Geographic Distribution of Total Chromium Emissions

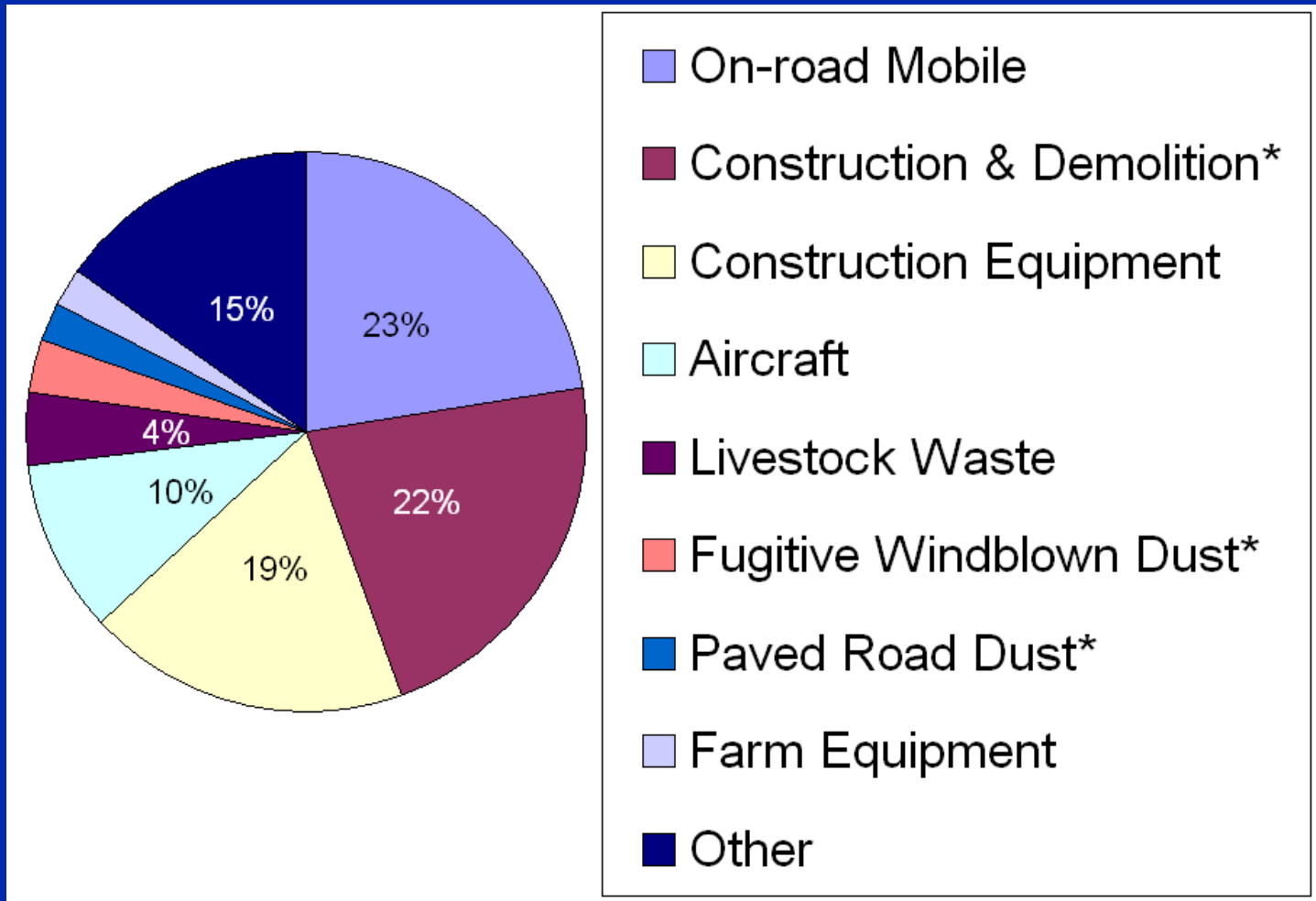


Geographic Distribution of Total DPM Emissions



Cancer Risk-weighted Emissions

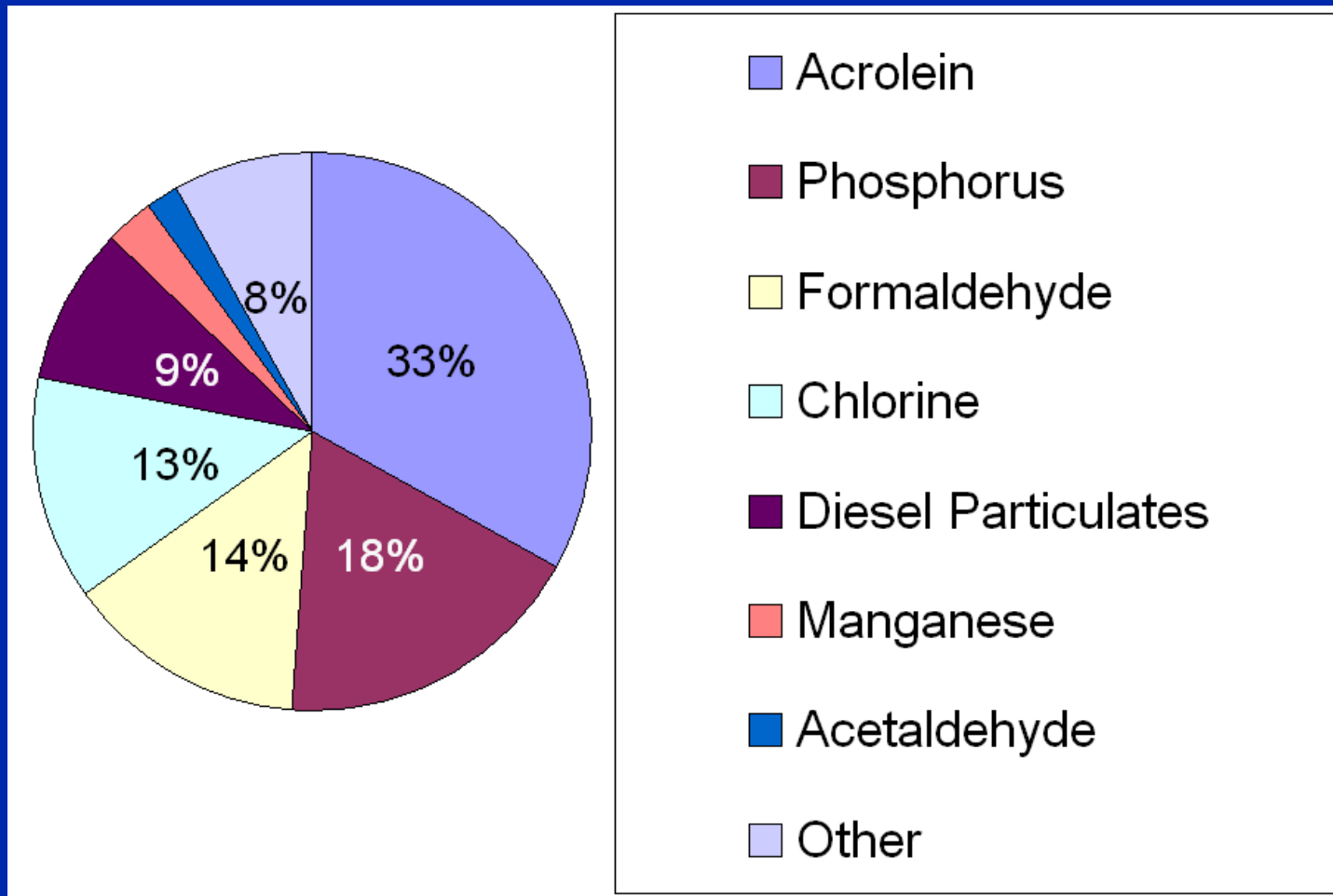
16 equivalent pounds/day



* denotes categories with large uncertainties

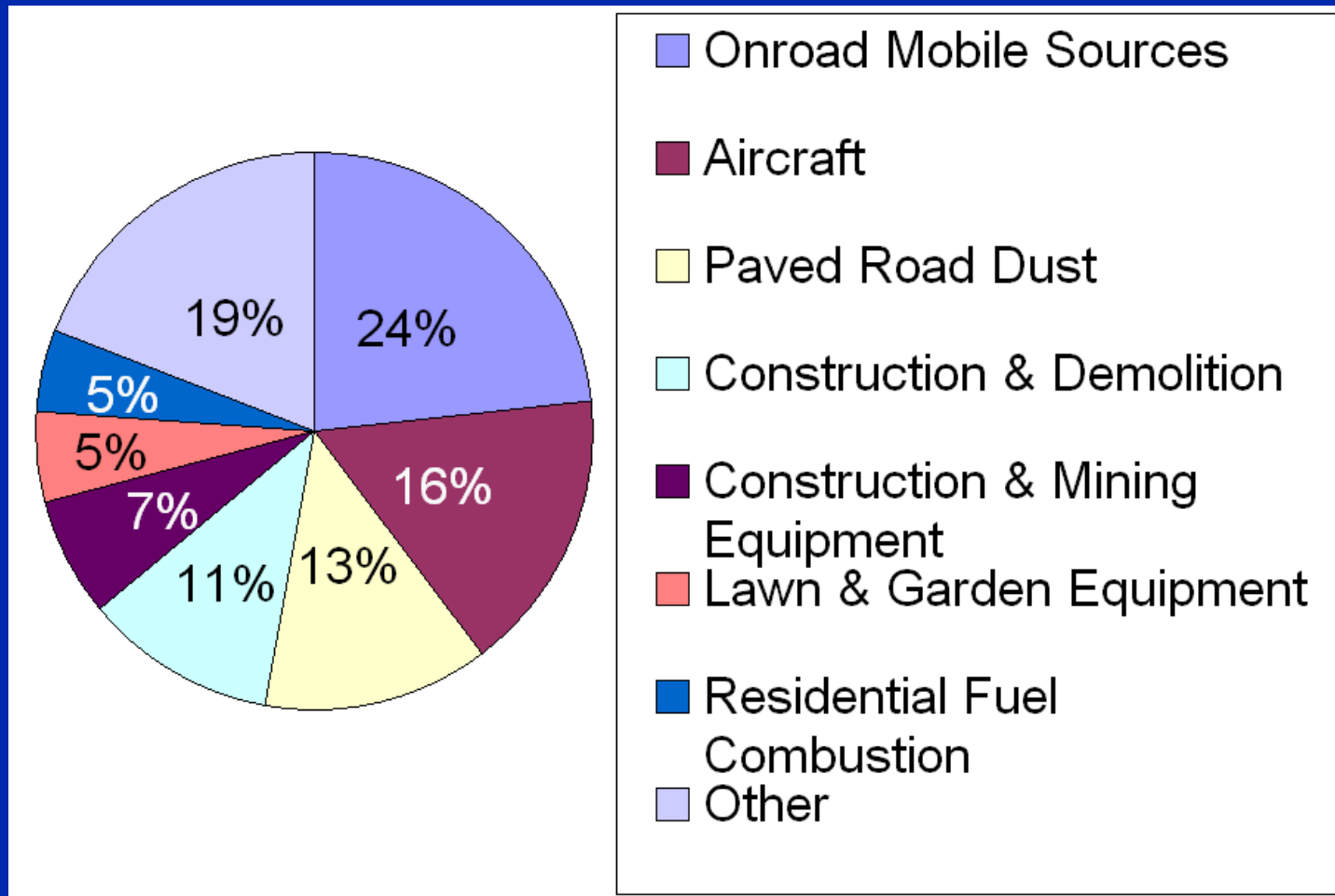
Chronic Risk-weighted Emissions

25 equivalent tons/day



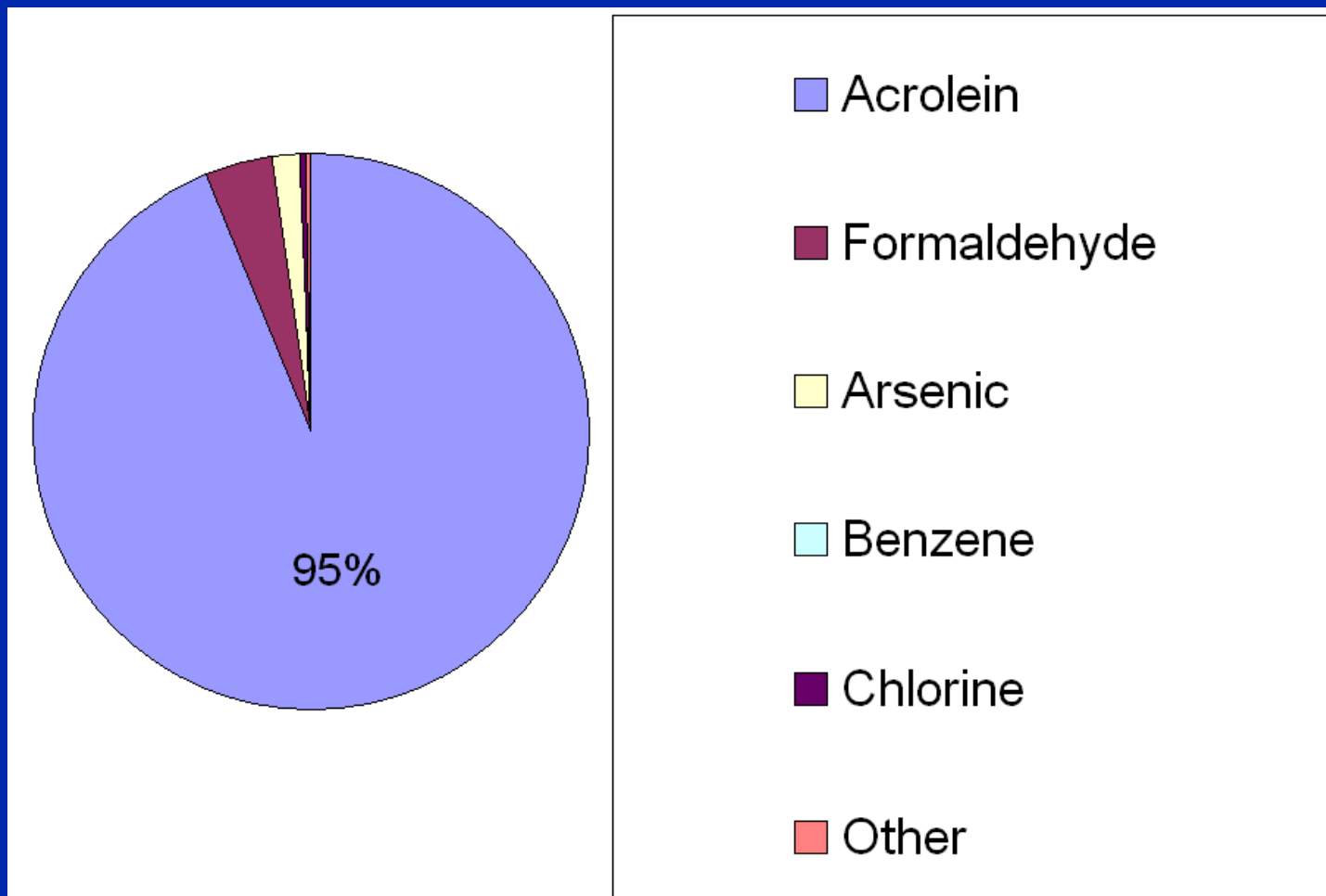
Chronic Risk-weighted Emissions

25 equivalent tons/day



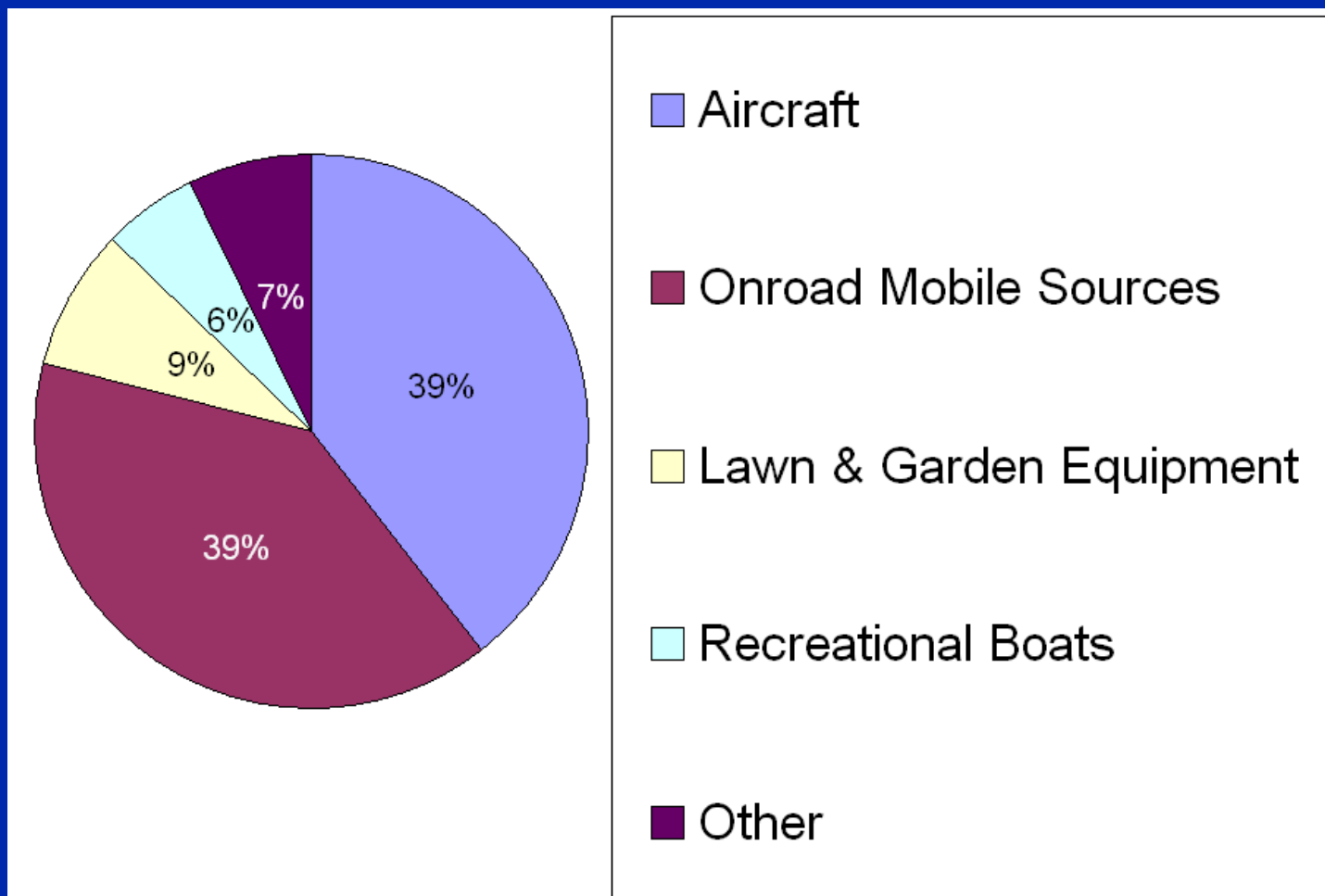
Acute Risk-weighted Emissions

2.8 equivalent tons/day

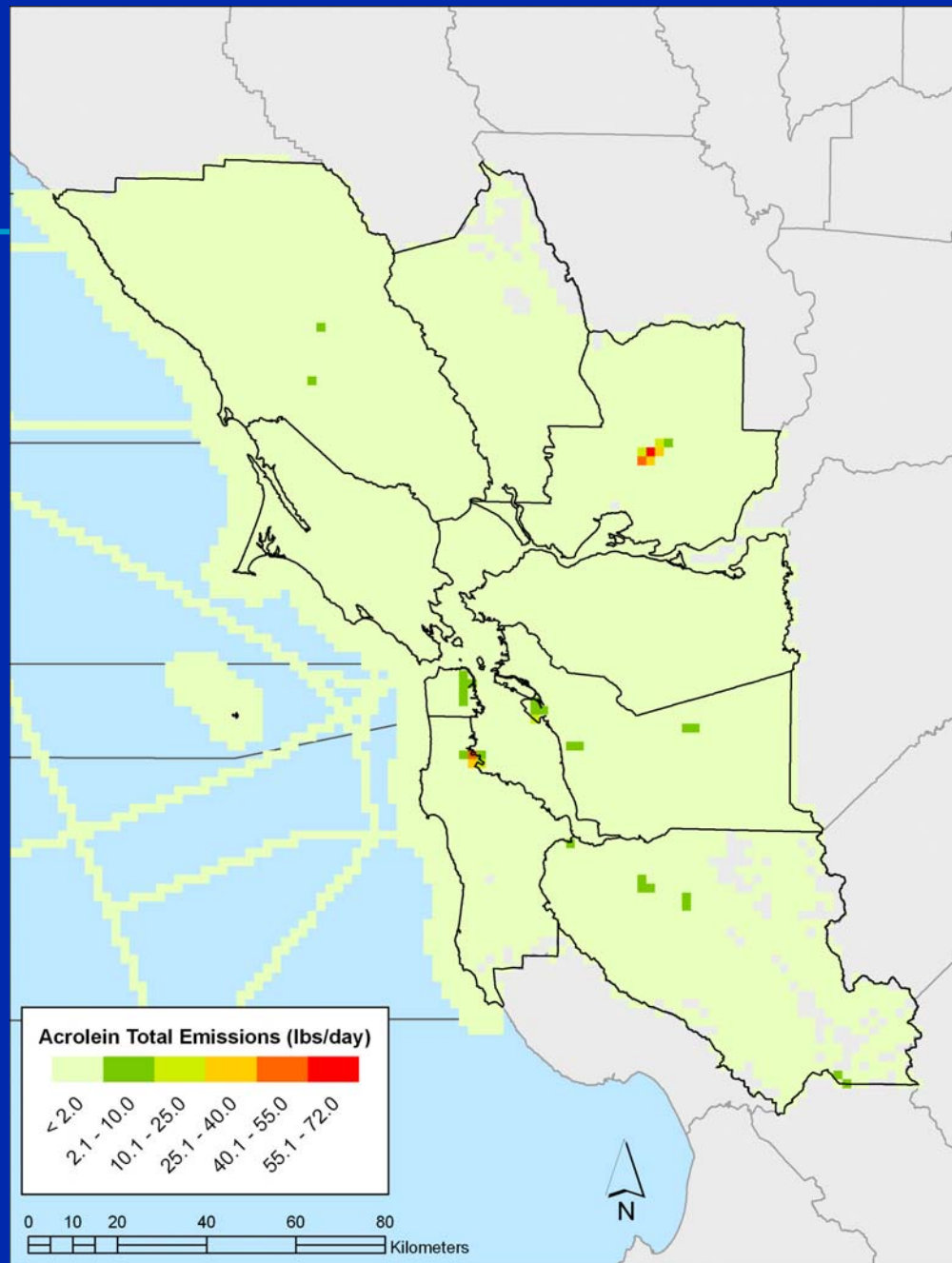


Acute Risk-weighted Emissions

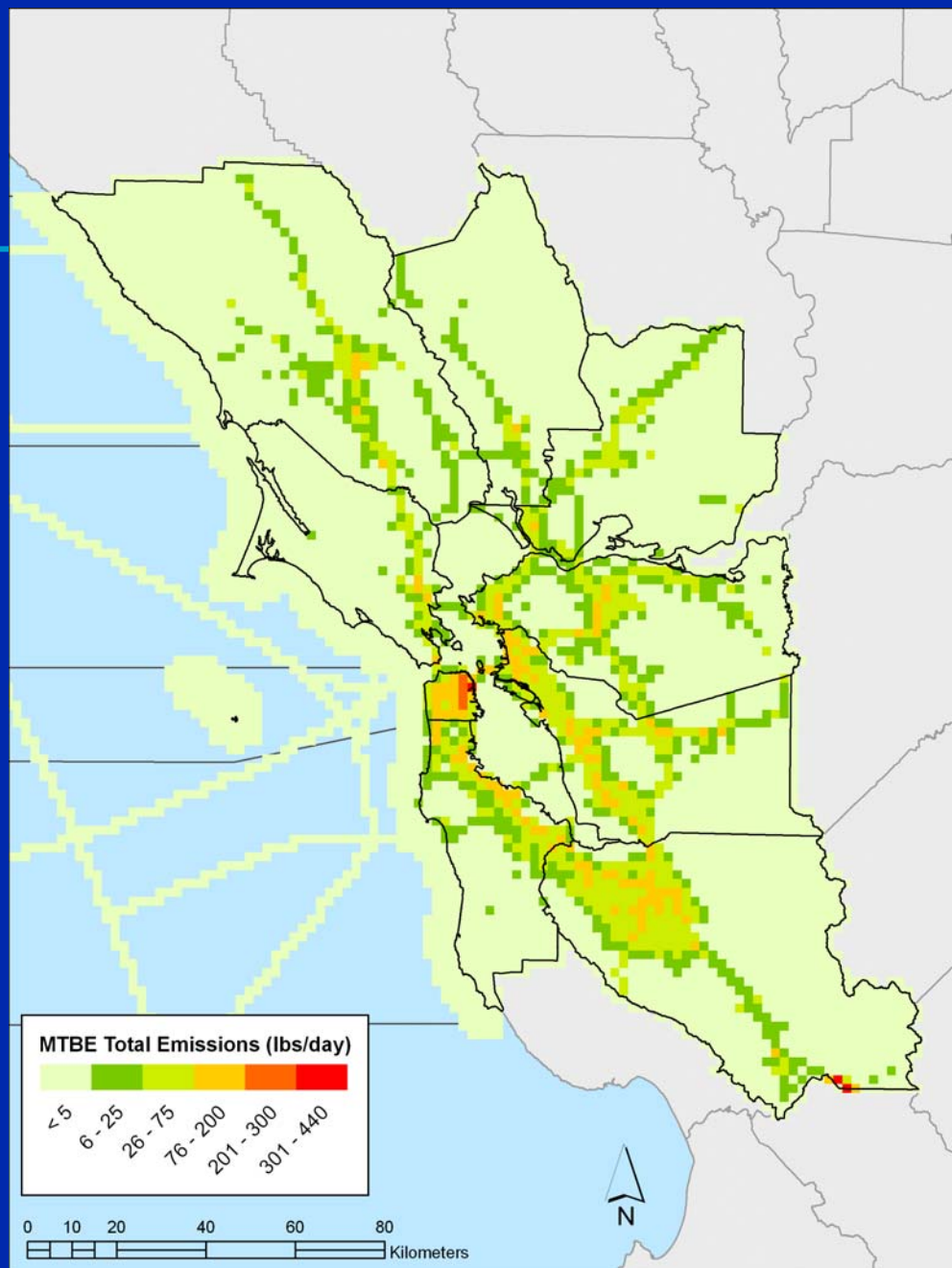
2.8 equivalent tons/day



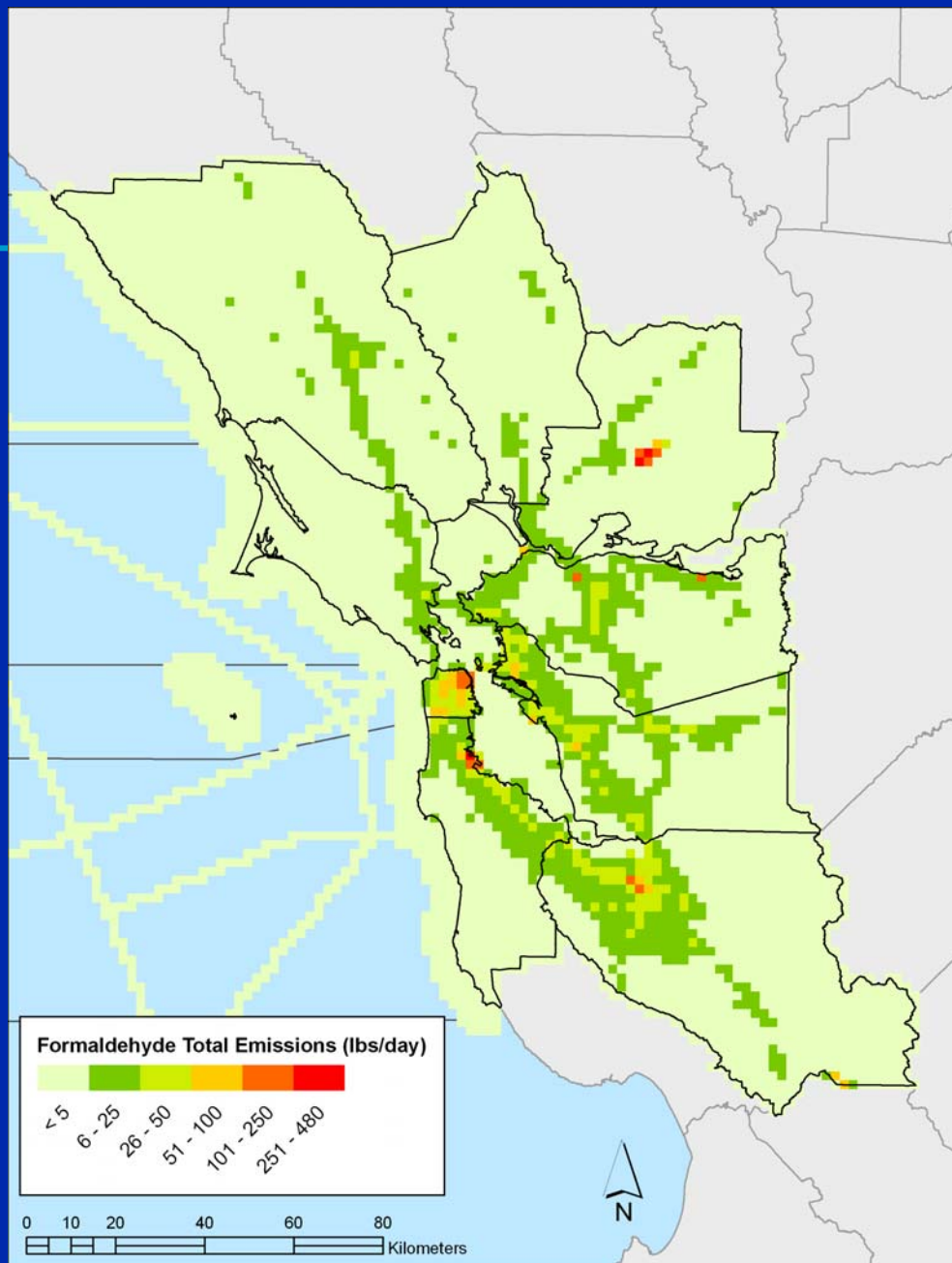
Geographic Distribution of Total Acrolein Emissions



Geographic Distribution of Total MTBE Emissions



Geographic Distribution of Formaldehyde Emissions



Strengths of the TAC Inventories

- Emissions for on-road mobile sources were estimated with EPA-recommended methods and California-specific chemical speciations.
- Emissions for point sources were directly reported.
- Spatial allocations are accurate and well-resolved for the modeling grid (2 km × 2 km).

Weaknesses of the TAC Inventories

- Some TACs are likely omitted (not components of TOG or PM_{10} or infrequently measured for chemical speciation profiles).
 - Quinoline and hydrazine are among the EPA's designated urban air toxics.
 - Others include radionuclides, titanium tetrachloride, hydrochloric acid, hydrofluoric acid, and sulfuric acid.
- Chemical speciation profiles contain significant uncertainties.
- Conservative assumptions were applied. For example, chromium compounds are emitted as chromium (VI) rather than chromium (III).

Primary Recommendations

- Prioritize TACs and continue inventory development and revisions using bottom-up methods.
- Add emissions for omitted TACs.
- Investigate emissions of chromium (VI) from sources of fugitive dust.

Emission Inventory-Based Site Evaluation

