

HAZARDOUS AIR POLLUTANTS

In Maine:

Emissions Inventory

Ranking System

AND

Prepared by:

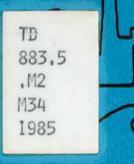
Bureau of Air Quality Control

Division of Technical Services

Bureau of Health

Environmental Health Unit

Environmental Toxicology Program



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Department of Environmental Protection

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JOSEPH E. BRENNAN GOVERNOR

March 25, 1985

COMMISSIONER -

Senator Ronald E. Usher, Senate Chair Representative Michael H. Michaud, House Chair Joint Standing Committee on Energy and Natural Resources State House, Station 2 Augusta, Maine Ø4333

Dear Senator Usher and Representative Michaud:

The Bureau of Air Quality Control of the Maine Department of Environmental Protection has completed the toxic air pollution inventory mandated in Chapter 835 of the Public Laws of 1984. With this report we respectfully submit the results of the inventory to the Committee on Energy and Natural Resources.

The report explains the process we followed in gathering data, what the collected information tells us and suggestions as to how the State of Maine should proceed to better control existing and future toxic emissions into the air. It is broken into three sections; the inventory, the air toxics ranking system and an appendix.

Throughout the inventory process the Maine Bureau of Health has been very helpful. In particular I want to thank Dr. Terry Shahata and Norman Anderson. In addition, credit should be given to David Dixon of the Bureau of Air Quality who organized and managed the inventory and data resolution processes. Finally, special thanks goes to the hundreds of businesses and industries that completed the questionnaires that form the basis of the data from which the inventory is drawn. Without their wholehearted help nothing could have been done.

In closing, as you use this report, questions and comments should be directed to Mr. Dixon. He will be most able and happy to help you. He can be reached at 289-2437.

Sincerely,

Henry E. Warren Commissioner

HEW/JLB/glk

cc: Richard Davies, Governor's Office

Portland •

EXECUTIVE SUMMARY

This report has been compiled by the Bureau of Air Quality Control and the Bureau of Health's Environmental Toxicology Program. Its purpose is to determine whether or not emissions of Hazardous Air Pollutants in Maine presents a serious environmental concern. The evaluation was carried out simultaneously with the Air Bureau collecting information on types and quantities of emissions while the Bureau of Health studied the relative toxicity or potential health hazard.

The report is divided into two sections: (1) the emission inventory and (2) the priority ranking system. Both sections are interdependent and must be evaluated collectively in order to determine the extent of environmental problems created by hazardous air pollutants. The hazard created is dependent on both factors, the toxicity or potency of the substance and the exposure.

The emissions inventory section includes a description of the procedures used to estimate emissions from point and area sources and discusses the findings. Based upon a list of 199 potential hazardous air pollutants selected to be inventoried and selected cutoffs for reporting, emissions of 6,143 tons per year were identified.

Of the 199 potentially hazardous air pollutants 58 were found to be emitted in Maine.

The Bureau of Health's toxicity ranking score is based upon four considerations: mutagenicity, carcinogenicity, reproductive effects and acute toxicity. The score in each of the four categories is a function of the evidence for each effect reported in the literature.

The toxicity score is combined with the exposure score to provide a numerical ranking of the inventoried hazardous air pollutants. The priority ranking system will be used by both Bureaus as a guide to systematically continue work on the hazardous air pollutant program.

The hazardous air pollutants which have been determined to undergo initial review are TOLUENE (present in gasoline and used as a solvent in paints and coatings), TETRACHLOROETHYLENE (used as a solvent principally for cleaning and extraction processes), FORMALDEHYDE (used in the manufacture of resins; as a perservative and hardening and reducing agent; in embalming fluids; as a corrosion inhibitor and sterilizing agent), BENZENE (present in gasoline and used as a raw material for organic chemicals), EPOXYPROPANE (used as a fumigant, and in the manufacture of urethane foams, surfactants and detergents, and synthetic lubricants), and CHLORINE (used as a bleaching and oxidizing agent).

Based upon the findings of the report, the following recommendations were developed. The recommendations presume that no single source of hazardous air pollutants is of such great public health concern that immediate pollutant abatement action is needed. The quantities emitted and potential health consequences of exposure to high concentrations or repeated exposures to lower concentrations, however, clearly demand continuing attention to sources emitting these potentially hazardous air pollutants. The entire process is systematic and procedures established by both Bureaus represent an appropriate level of effort.

The following recommendations are made based on the above conclusions:

Authorize the Commissioner to take immediate action when he finds imminent public health risk attributable to existing emissions of hazardous air pollutants.

Provide the Board of Environmental Protection the clear authority to require proper control of hazardous air pollutant emissions from any new source or new process of an existing source whenever the proposed emissions may pose an unacceptable health risk.

Require that existing air emission licensed sources undergo licensing review for hazardous air pollutants.

Authorize the Commissioner to collect additional source information for hazardous air pollutants.

SECTION 1. EMISSIONS INVENTORY

1. INTRODUCTION

Chapter 835 of Public Laws of 1984 required that the Department of Environmental Protection carry out and maintain an inventory of sources in the State emitting any substance that may be a potential hazardous air pollutant. The inventory was specifically required to include the following:

- (1) the number of sources;
 - (2) the location of each source or category of source;
 - (3) the quantity emitted by each source or category of source;
- (4) the total emissions; and

(5) the percentage of total emissions generated by sources with existing air licenses.

This section of the report describes the process utilized for data collection, emission calculations and quality assurance. Then it presents the findings of the inventory and our recommendations to the Committee.

2. PROCEDURE

The Department developed a comprehensive questionnaire based upon EPA guidelines and examples of data collected in other states. We solicited and received input from the Air Quality Research Coalition. The coalition is a paper tanneries, group representing industries, electronics pulp & manufacturers and other industries which showed a common interest in an air toxics program. As a result of discussions, we concluded that the short term objective of the inventory should be to develop a prioritization scheme for addressing specific air toxics. Based on that objective we developed a general questionnaire which asked only for the total use of a substance and total plant wide emissions, rather than production line by production line data. The scope of the inventory was further limited:

1. By establishing a list of 199 substances which were considered to be potential hazardous air pollutants by the Bureau of Air Quality Control and the Bureau of Health of the Department of Human Services, and;

2. By establishing a minimum use rate of 2000 pounds per year, below which no reporting was required.

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The inventory questionnaire was designed to obtain information from 4 categories of sources:

(1) process sources

(2) incineration sources

(3) storage facilities

(4) loading, unloading, transfer operations.

In late July questionnaires were sent to approximately 700 sources based upon three factors:

(1) facilities with existing air emission licenses,

(2) Standard Industrial Classification (SIC) and employment level, using the Maine Marketing Guide,

(3) hazardous waste manifest data of the Bureau of Oil and Hazardous Materials.

Questionnaries were due back October 31, 1985. At that time the response rate was between 60-65 percent. In looking through the early response data it was clear that 7 categories in the original sample would have little if any impact on the inventory. Because they would have such little impact we removed them from the sample. They were:

(1) schools

(2) hospitals

(3) veterinarians

- (4) hotels
- (5) nursing homes
- (6) oil distribution companies
- (7) sewage treatment plants

Eliminating these categories resulted in a 75 percent response rate. All the remaining sources which had not reported were contacted individually via telephone during the first two weeks of November. Each was reminded of the inventory and its importance.

During the first week in November, all questionnaires were preliminarily reviewed. Based upon this evaluation 3 categories of sources were selected for on-site follow-up, as a verifying technique to assure that the data was accurate. The categories selected were:

(1) Sources that reported significant emissions. Several were selected for an on-site inspection for quality assurance purposes and to verify methods used to calculate emissions,

(2) Sources who responded and reported no emissions. We selected several for on-site inspections based upon their SIC code and employment level or similarity with sources that reported emissions,

(3) Sources who did not respond at all. We went to all those whom we expected may be significant sources based upon their SIC code and level of employment or similarity with sources that reported emissions.

During December and early January 1985, a total of 54 sources received on-site follow-up inspections to verify data. This represents a quality control check of 29 percent. Significant additional emissions were found during this process. By this point the inventory response rate was 518 out of 619 or 84 percent. Based upon the follow-up inspections, we believe that the sources inventoried in this report include the majority of significant emitters in Maine.

In order to deal with the volume of data reported during the inventory we developed an on-line computer data handling system. This system utilized the Department's Honeywell Level 6 mini-computer. Simply put, data could be entered directly to the computer and reports written using a generalized information retrieval system. Inherent in the design of the system was a method to maintain the confidentiality of the information reported to us.

The quality of the responses to the questionnaire varied significantly. In some cases, it was evident that a great deal of time and effort had gone into the preparation of the reports. Documentation of calculations, detailed process descriptions, and control/emission assumptions were provided. The majority of submittals, however, simply reported only the amounts of chemicals used and left estimates of the emissions to the Department. In those cases, staff relied on talking with the particular plant operator, comparing between similar industrial sources and making the best estimate based on the chemical and its use. Because of the nature of this estimation process, we purposely made our assumptions conservative which means that the emissions may be somewhat overestimated.

More uncertainty is probably introduced by the methods used to calculate emissions. It was recognized that stack tests or engineering studies should not be required for collecting emission estimates for the purpose of this inventory. Two methods were commonly employed; emission factors and material balance calculations. Emission factors are a typical emission rate that would The factor is then multiplied by the be expected from a type of source. process rate to estimate emissions. Material balance calculations use the amount of product purchased minus what is sent out as a waste and what is tied up in the product. The quantity unaccounted for is assumed to be lost to the Much may initially be lost to the sewer system but even that will air. eventually evaporate. This method may, therefore, overestimate emissions at the source but is quite accurate for accounting for all that eventually reaches the ambient air.

It needs to be pointed out that time did not permit a thorough review of all questionnaires so in many cases estimates are based upon the source's

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response. Those that were checked during on-site visits indicate the data are very reliable. Nevertheless, there were instances of misunderstanding and cases of unreasonable assumptions, so it is conceivable that reported values may in some cases inaccurately represent real emission rates.

On the other hand the minimum cut-off exclusion and the select nature of the list tend to make the process underestimate the total statewide emissions of hazardous air pollutants. The use of emission factors, and material balance calculations with conservative assumptions probably overstates emissions in many instances. The estimates therefore should be treated as estimates, not as absolute values. The data is valid for drawing comparisons and developing a prioritization ranking scheme which was the major goal of the project.

The inventory does not address two other situations which may be of concern. It neither provides qualitative and quantitative estimates of reaction products resulting from emissions of individual precursors, nor estimates hazardous air emissions from trace constituents of oil or coal burning. The latter would increase the total reported emissions of hazardous air pollutants so that more attention may be directed at the program; the regulatory significance is not great because control is, and should be, dictated by emissions of the major constituents.

The data collected is not intended to be used to assess local ambient air impacts because it did not include flow rates, densities, elevations and other similar characteristics of the emission that are necessary to quantify impact.

3. AREA SOURCES

Area source emission estimates are included for several categories of sources whose emissions are too small to catalogue individually but whose collective emissions are significant. Area sources are reported by category with a geographical breakdown by county and major urban areas. In most cases, the distribution is based upon population distributions or data on vehicle miles traveled in an area. Emissions for the area sources are based exclusively on emission factors. The source of the data used, emission factors and assumptions are all listed for each catagory in Appendix 3.

Table 1 contains each of the categories of sources included in the area source inventory:

AREA SOURCE CATEGORIES

Dry Cleaners Degreasers Open Burning Dumps Agricultural Open Burning Forest fires Architectural Surface Coating - (Painting) Highway Markings Printing Residential Wood Burning Waste Oil Burning Automotive Emissions - (Lead) Gasoline Stations Operation - (filling underground tanks, vehicular refueling and spillage)

Gasoline Terminals

4. FINDINGS

1. The study found 57 of the 199 inventoried hazardous air pollutants were emitted in Maine. Since this inventory was limited to sources using in excess of 2,000 pounds per year, it is not all inclusive. Some sources may use less than 2,000 pounds of a pollutant which the inventory would not detect.

2. The study indicated 187 sources of the reporting 619 had some reportable emissions. Given enough time and resources to follow-up on all inventoried sources, there is little doubt more emissions would have been found.

3. Of the 178 sources which reported emissions of potential hazardous air pollutants, 78 sources or approximately 44 percent have existing air emission licenses.

4. Area sources were found to be a significant source of hazardous air pollutants in a few cases. They are dry cleaners, gasoline marketing, residential wood burning, and waste oil burning.

5. Table 2 presents a statewide summary of the top 25 hazardous air pollutants for both point and area sources.

6. The total emissions of hazardous air pollutants reported is 6,143 tons (12,286,000 Lb.). Compared to a total emissions inventory of 106,000 tons (212,000,000 Lb.) of sulfur dioxide, the generic category of hazardous air

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TOP 25 HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE

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DOLLUTANT		EMISSIONS (POUNDS)	TOTAL
POLLUTANT	POINT SOURCE	AREA SOURCE	TOTAL
1. Toluene	2,459,231	52,360	2,511,591
2. 1,1,1, Trichloroethane	1,496,539	0,	1,496,539
3. Chlorine	1,337,003	0	1,337,003
4. Xylene	1,091,274	0	1,091,274
5. Acetone	804,521	0	804,521
6. Tetrachlorethylene	144,450	563,600	708,050
7. Methyl Cellosolve	632,248	0	632,248
8. Methyl Ethyl Ketone	512,611	0	512 , 611
9. Methyi Mercaptan	428,940	0	428,940
10. Chlorine Dioxide	296,787	· 0	296,787
11. Methylene Chloride	288,569	0	288,569
12. Ethylene Glycol Ethyl E	ther 277,118	0	277,118
13. Hydrogen Sulfide	269,912	0	269,912
14. N-Butyl Acetate	253,563	0	253,563
15. Lead	630	202,800	203,430
16. Methyl Metharcylate	184,550	0	184,550
17. Butanol	149,490	0	149,490
18. Diphenylmethane 4,4-di-			
isocyanate	146,000	0	146,000
19. Hydrogen Chloride	44,731	96,000	140,731
20. 1,2 – Epoxypropane	124,600	0	124,600
21. Trichloroethylene	110,600	0	110,600
22. Formaldehyde	79,401	0	79,401
23. Turpentine	77,130	0	77,130
24. Ethyl Acetate	66,537	0	66,537
25. Benzene	0	52,392	52,392

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pollutants is quite small. However, this is not a fair comparison due to the varying public health risk which hazardous air pollutants pose. This comparison is made to put the volume of hazardous air pollutants in perspective with a more familiar criteria air pollutant.

7. Of the 6,143 tons of inventoried hazardous air pollutants only 167 tons (2.7%) would have sufficient evidence to be classified as carcinogenic. This is not to suggest that carcinogens should receive special treatment but rather to confirm that most of what were identified as hazardous air pollutants were not carcinogenic.

5. DISCUSSION OF FINDINGS

The problem with a discussion of the results of the inventory is the age old one of comparing apples and oranges. The list of hazardous air pollutants is composed of chemicals with such varying public health risks that they should not be compared to each other. It is the role of the Bureau of Health and the Scientific Advisory Committee to perform this risk assessment.

The findings represent broad generalizations of the data reported. In many cases it is not the generalizations that are very useful but rather the specific data. That data is reported in a series of appendices:

1. Appendix 1 is the List of Hazardous Air Pollutants for which the Department inventoried. This list should be referred to in the following appendices as it identifies a six letter code for each pollutant. The code was used as part of the computer data handling system.

2. Appendix 2 reports the total process emissions for all inventoried point sources.

3. Appendix 3 reports the individual area source documentation sheets. The initial sheet describes the category, data collection, data assumptions, emission factor(s), emission factor source, calculation method, and annual emissions summary. The second sheet details the emissions by county and for a few major cities.

4. Appendix 4 reports a breakdown of hazardous air pollutants by county.

5. Appendix 5 reports a breakdown of hazardous air pollutants by Standard Industrial Classification (SIC).

6. RECOMMENDATIONS

following recommendations legislation are made with the The for understanding that while significant quantities of hazardous air pollutants are emitted into the air at present there is no single hazardous air pollutant or source that represents a clear threat to public health from these emissions. Nevertheless, it is prudent to limit or possibly regulate future sources of hazardous air pollutants that may pose a direct threat to public health and require the same level of control generally required for the criteria air pollutants. At the same time the Department should be authorized to gather the source specific information necessary to conduct ambient air quality impact assessments.

With this in mind, the Department's recommendations are:

A. LEGISLATIVE

1. Emergency Provision to Protect Public Health

The Commissioner should be given the clear authority to take immediate action to stop emissions of hazardous air pollutants when they pose an imminent threat to public health. The language should be similar to the existing emergency provisions found in Hazardous Waste Management law, Title 38, Section 1310.

2. New Emission Sources

Legislation is needed to clarify and improve the capacity of the Board of Environmental Protection to regulate hazardous air pollutant emissions from any new source or new process or operation at an existing source when it finds that the proposed emissions pose an unacceptable public health risk.

3. Existing Emission Sources

Existing law should be clarified to clearly indicate that hazardous air pollutants are treated similar to the criteria pollutants in the licensing of any existing air emission source. This may require existing sources of hazardous air pollutants to use Best Practical Treatment.

4. Source Information

The Commissioner should be given the authority to solicit additional source information such as:

a. process schedules and design specifications;

b. stack or discharge vent operating parameters;c. control equipment specifications and design operating conditions; d. emission testing either periodically or using continuous emissions monitors.

B. DEPARTMENTAL

1. Emissions Inventory

The emissions inventory should be upgraded to include source specific emission data. Testing and documentation of emission estimates should be expanded to cover sources emitting less than 2000 lbs/year.

SECTION II. RANKING HAZARDOUS AIR POLLUTANTS FOR FORMAL EVALUATION

1. INTRODUCTION

Section 585-C of 38 M.R.S.A. mandated the Maine Department of Environmental Protection (DEP) to establish an air emissions inventory for hazardous air pollutants. With the assistance of the Maine Bureau of Health and other state agencies, the DEP developed an initial list of two hundred air pollutants to be classified as "hazardous." It is the responsibility of the Bureau of Health (22 M.R.S.A. Section 1696) to assess the public health risk associated with these hazardous air pollutants. In conducting its assessments, the Bureau relies on the advice and constructive criticism of its Scientific Advisory Panel. This Panel, established concurrently with the hazardous air pollutant legislation, consists of expert health professionals from academia, industry, consulting firms, and private practice.

2. APPROACH

In order to know the sequence in which pollutants will be assessed, it is first necessary to develop a ranking system based on a preliminary evaluation of each pollutant's toxicity and the quantity that is emitted into the ambient Two components are assessed in the ranking system: toxicity and air. exposure. Combining these two components is necessary for a balanced perception of actual public health risk. The ranking measures all the pollutants against a standardized set of criteria, and assigns numerical scores based on these criteria. This method has inherent problems as it does not consider the particular toxicologicial or emission characteristics of a chemical. On the other hand, consideration of specific characteristics may undermine the purpose of the ranking system. Since it is only a scheme for setting priorities, its effectiveness can be diluted by introducing criteria which are not easily comparable among a wide variety of pollutants. Yet, while not directly translatable into a measure of public health risk, this ranking system should provide a relative index of the pollutants' potential health threats.

3. TOXICITY COMPONENT

The toxicity component of the ranking system provides a preliminary score based on the studies cited in the National Institute for Occupational Safety and Health's <u>Registry of Toxic Effects of Chemical Substances</u> (RTECS). The effects categories are based on the following health criteria: carcinogenicity, mutagenicity, reproductive effects, and acute effects. Values for each category range from zero, except acute toxicity, which ranges from one to four. The toxicity factor for a pollutant is determined first by adding the individual category scores, and then adding to this sum the standard deviation of the four values. The reason for adding the standard deviation will be explained below.

The mechanism by which the health effects are scored is outlined in Table 1. Though each category is evaluated by different parameters, the basic ordering reflects a hierarchy of concern regarding a pollutant's human health risk. A pollutant is scored as a carcinogen, mutagen, or a reproductive toxin according to its likelihood of acting as such in human beings. The acute toxicity category is scored according to potency criteria, since any substance can be toxic to human beings if administered in large enough doses. With the exception of the acute toxicity category, no consideration is placed on route of exposure.

<u>Carcinogenicity</u> For carcinogenicity assessment, RTECS reflects the weight of evidence criteria used by the International Agency for Research on Cancer (IARC) and the National Toxicology Program. The evidence is considered with regard to whether it represents the results of human or animal studies, and whether the resulting associatons are positive, suspected, indefinite, negative, or lacking data. Using the assumption that the compounds with the closest associations with human carcinogenesis warrant the most immediate attention, the following scoring scheme was developed.

Human: Positive/Suspected	4
Animal Positive	3
Animal Suspected	2
All Other Non-negative data	1
Adequate Negative data	0

<u>Mutagenicity</u> Though actual guidelines still need to be developed for interpreting mutagenicity data, there is the general recognition that those tests which more closely reflect human physiological responses to the suspected mutagen also provide the greatest weight of evidence. The following scheme was thus developed for interpreting RTECS mutagenicity data in light of this general premise.

In vitro or in vivo human, in vivo mammalian	4
In vitro mammalian	3
Non-mammalian	2
No Data	1
Adequate Negative data	0

The study (or studies) which have the highest ranking determine the pollutant's score for this health effect.

<u>Reproductive Effects</u> Scoring for reproductive effects follows basically the weight of evidence ranking used by IARC in its evaluation of potential carcinogens. This approach differs from IARC's in that the triggering studies have not necessarily been peer reviewed. Because of the generally limited data base for this category, however, less demanding criteria have been employed.

Human data, or	4
2 species and 2 routes in 1 species	
2 species or 2 routes in 1 species	3
1 species tests	2
No data	1
Adequate Negative data	0

<u>Acute Toxicity</u> As shown in the Table 2, numerous regulatory agencies categorize acute toxicity into roughly four levels. This categorization forms the basis of the scoring assignments for this effect. To reflect the importance of actual human data, pollutant scoring gives priority to any human lethal dose studies cited. For those pollutants which lack human data, priority is given to the LC50* (4-hour exposure period) in any animal species, since this is a measure of toxicity via inhalation. Pollutants with neither human lethal dose data nor lethal dose animal inhalation data are scored according to the LD50. For the sake of consistency, the rat has been used as the reference species for the LD50. When rat data are not available for a pollutant, toxicity data from any rodent is used.

> Level 4 LC (human) or LC50 (animal): less than 0.2 mg/liter, less than 200 ppm; or LD (human) or LD50 (animal): less than 50 mg/kg

4

3

2

1

- Level 3 LC (human) or LC50 (animal): 0.2 - 2.0 mg/liter, 200 - 2000 ppm; or LD (human) or LD50 (animal) 50 - 500 mg/kg
- Level 2 LC (human) or LC50 (animal): 2 - 20 mg/liter, 2000 - 20,000 ppm; or LD (human) or LD50 (animal): 500 - 5000 mg/kg
- Level 1 LC (human) or LC50 (animal): greater than 20 mg/liter, greater than 20,000 ppm; or LD (human) or LD50 (animal): greater than 5,000 mg/kg; or No Data

Deviation. There are several difficulties involved with the derivation of a general toxicity value for a wide variety of pollutants. In this scheme,

there is the concern that pollutants with wide deviations in their health effect scores (particularly in the carcinogenicity, mutagenicity, and reproductive effects scores) may be underestimated in their rankings because of inadequate data. To compensate for this, the standard deviation of the four health effects scores was added to the sum.

4. EXPOSURE COMPONENT

Emissions estimates for all hazardous air pollutants have been provided by the Department of Environmental Protection's Bureau of Air Quality Control. These estimates are in the form of statewide summations in pounds per year from industrial, commercial, residential, and mobile sources.

* For explanation of terms, see GLOSSARY

5. RANKING ALGORITHM

The DEP inventory, because of the wide range in the emissions estimates, is not easily comparable to the toxicity scale. A variety of methods can be used, each given varying degrees of weight to each component. It is unknown at this time how different modifications of the exposure component affect its weight in the final scoring. Given the uncertainties involved with the weighting process, a non-parametric approach is used. In this approach, the toxicity and exposure components for each pollutant are ranked in order of decreasing toxicity an decreasing emissions. The toxicity and exposure ranks are then added to produce the total score. Pollutants decrease in priority, therefore as their total scores increase. The results of this algorithm are presented in Table 3.

6. DISCUSSION

Despite the efforts of the National Institute for Occupational Safety and Health to list all journal sources, it is entirely possible that some important studies have not been included. Also, with the exception of carcinogenicity, the health effects scoring is drawn from sources that have not been peer reviewed. Finally, "negative data" cannot be distinguished in RTECS from "no data." A provision has been made in the ranking system to score a health category as zero if findings of a negative effect have been demonstrated.

The health effects categories were given equivalent scoring ranges to minimize the biases involved in weighing the different health effects. This procedure, however, does not eliminate the biases inherent in the criteria themselves. Both carcinogenicity and reproductive toxicity are highly correlated with In other words, a compound having a high score for either mutagenicity. carcinogenicity or reproductive effects most likely has a high score for mutagenicity as well. Ranking, therefore, may be disproportionately weighted against pollutants which may be very toxic, but not mutagenic. On the other hand, this unequal emphasis may also more closely reflect the public health Compounds present in sufficient concentrations to produce acute threat. effects are generally more easily detected and controlled than compounds which can produce effects through long-term, low level exposures. Nevertheless. this concern can of unequal weighting be mitigated to a certain extent by prior consideration of "supertoxics," or compounds whose acute toxicity doses are less than the most toxic dose levels described in Table 2. These compounds are listed in Table 4. Before the formal risk assessments are conducted, assessments will be done on these pollutants to determine whether a potential public health problem exists due to acute exposures.

Important facets of a compound's potential toxicity, such as chronic toxicity and synergistic effects, have not been explicitly considered in this process. People are rarely exposed in the ambient environment to only one pollutant. There are several examples to illustrate the possibility that exposures to mixtures of pollutants can create unforeseen effects. There are also many ways in which pollutants can promote disease development through interactions with lifestyle or genetic factors. This shortcoming may be alleviated to a certain extent through the unequal weighting. As a class, mutagenic compounds may be more likely to produce interactive effects than non-mutagenic compounds. By generally scoring mutagenic compounds higher than non-mutagenic compounds, this system may indirectly consider these effects. Unfortunately, there is currently no systematic method to assess these concerns. They will be assessed in the formal risk assessments, however, and attempts will be made to include these criteria in future ranking systems.

In many instances, the basis for a pollutant's statewide emissions estimate has questionable credibility. Actual emissions test for hazardous air pollutants have rarely been done in Maine. Much information has been obtained through emission factors, material balances, or purchase records. Estimates of hazardous air pollutants emitted from combustion processes are, for the most part, unavailable. Pollutant exposures resulting from consumer or small commercial usage are also difficult to estimate. These limitations, however, have been recognized at the outset. Also recognized was the fact that the hazardous air pollutant program could not wait for these limitations to be overcome. Rather, the intention is to more rigorously address exposure criteria as the public health concerns related to hazardous air pollutants become more clearly understood.

For all of these reasons, the ranking system should be kept flexible so that it can easily respond to additional information or more critical review of the literature. It would be counterproductive for the Bureau of Health to verify the data used in the system, since the purpose of the ranking process is to order pollutants without becoming involved in specifics. On the other hand, assistance from outside sources is appreciated and is encouraged during this and all ensuing steps of the pollutant assessment process.

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GLOSSARY

- LC: lethal concentration
- LD: lethal dose
- LC50: lethal concentration for 50% of the animals in the study
- LD50: lethal dose for 50% of the animals in the study
- mg/liter: milligrams of compound per liter of air
- ppm: parts per million (volume)
- mg/kg: milligrams of compound administered per kilogram body weight

TOXICITY RANKING SYSTEM

		Score
	A. Carcinogenicity (C)	
	Human: Positive/Suspected Animal Positive Animal Suspected All Other Non-negative data Adequate Negative data	4 3 2 1 0
	B. Mutagenicity (M)	
	In vitro or in vivo human, in vivo mammalian In vitro mammalian Non-mammalian No Data Adequate Negative data	4 3 2 1 0
	C. Reproductive Effects (R)	
	Human data, or 2 species and 2 routes in 1 species 2 species or 2 tests in 1 species 1 species No data Adequate Negative data	4 3 2 1 0
	D. Acute Toxicity (A)	
	Level 4 LC (human) or LC50 (animal): less than 0.2 mg/liter, less than 200 ppm; or LD (human) or LD50 (animal): less than 50 mg/kg	4
	Level 3 LC (human) or LC50 (animal): 0.2 - 2.0 mg/liter, 200 - 2000 ppm; or LD (human) or LD50 (animal) 50 - 500 mg/kg	3
	Level 2 LC (human) or LC50 (animal): 2 - 20 mg/liter, 2000 - 20,000 ppm; or LD (human) or LD50 (anmial): 500 - 5000 mg/kg	2
	Level 1 LC (human) or LC50 (anima1): greater than 20 mg/liter, greater than 20,000 ppm; or LD (human) or LD50 (anima1): greater than 5,000 mg/kg; or No Data	1
Toxicity	Sum = Sum of Individual Health Effects Scores S = Standard Deviation of a pollutant's effects scores Score = Sum + S -16-	

ACUTE TOXICITY TESTS FOR SELECTED FEDERAL REGULATORY PROGRAMS

CATEGORY	A: Mists, Dusts, and Fumes
OSHA	Highly Toxic Toxic
HMTA	Poison A or B
FHSA	Highly Toxic Toxic
FIFRA	I II III IV
RCRA	Acutely Hazardous
	.02 .2 2 20 200 LC ₅₀ (mg/L)
CATEGORY	
un zoonn	
OSHA	Highly ToxicToxic
FHSA	Highly Toxic Toxic
CWA	Hazardous
CATEGORY	20 200 2000 20,000 200,000 LC ₅₀ (ppm) C: Oral Toxicity
OSHA	Highly ToxicToxic
HMTA	Poison A or B
FIFRA	I II III IV
CWA	Hazardous
RCRA	Acutely Hazardous
	5 50 500 5000 50,000 LD ₅₀ (mg/kg)
OSHA HMTA FHSA FIFRA CWA RCRA	Occupational Safety and Health Act Hazardous Materials Transportation Act Federal Hazardous Sustances Act Federal Insecticide, Fungicide, and Rodenticide Act Clean Water Act Resource Conservation and Recovery Act
Source:	U.S. Environmental Protection Agency, <u>Chemical Substances</u> <u>Designation, Vol. I:</u> <u>Overview and Analysis</u> , Washington, D.C., 1981. -17-

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RANKINGS FOR HAZARDOUS AIR POLLUTANTS

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Pollutant	Toxicity Score	Emissions (1bs/yr)	Toxicity Rank	Emissions Rank	Total
Toluene Tetrachloroethylene Formaldehyde Benzene Epoxypropane Chlorine Methylene Chloride 1,1,1-Trichloroethane Lead Styrene Trichloroethylene Benzo-a-Pyrene Xylene Methyl Mercaptan 1,2 Dichloroethane Methyl Cellosolve	12.5 13.2 15.5 15.5 13.5 11.7 12.5 11.0 12.5 15.5 13.2 14.6 10.3 10.5 13.2 9.4	2,511,623708,05079,40152,392124,6001,433,003288,5691,496,539178,63013,001110,86515,1801,091,274428,94032,400632,248	18 15 2 2 10 27 18 30 18 2 15 7 34 31 15 36	$ \begin{array}{r} 1 \\ 6 \\ 21 \\ 24 \\ 19 \\ 3 \\ 13 \\ 2 \\ 16 \\ 32 \\ 20 \\ 30 \\ 4 \\ 9 \\ 28 \\ 7 \\ \end{array} $	19 21 23 26 29 30 31 32 34 34 34 35 37 38 40 43 43
Methyl Methacrylate Hydrogen Chloride Bis 2-ethylhexyl phthalate Chlorine Dioxide Ethylene Glycol Ethyl Ether Napthalene Acetone Methyl Ethyl Ketone Arsenic Hydrogen Sulfide Hydrazine Ethylene Oxide Formic Acid	11.3 12.5 13.4 9.4 9.4 12.5 7.5 8.0 16.0 8.5 14.6 13.4 10.5 15.5	184,550 44,731 10,190 296,787 277,118 11,799 804,521 512,611 430 269,912 740 1,535 50,433 184	28 18 11 36 36 18 46 43 1 41 7 11 31 2	15 26 34 10 11 33 5 8 50 12 49 45 25 54	43 44 45 46 47 51 51 51 51 53 56 56 56 56
Chromium Methyl Chloride Zinc Cadmium Epichlorhydrin Phenol n-Butyl Acetate Diethyl Sulfate Butanol Copper Diphenyl Methyl 4,4-Diisocyanate Manganese Turpentine Nitric Acid Ethyl Acetate	11.3 13.4 14.6 15.5 12.3 6.6 12.0 6.6 13.4 5.5 .10.0 5.5 7.0 5.5	16,800 946 238 12 5,180 253,563 5,424 149,490 107 146,000 9,320 77,130 41,450 66,537	28 11 7 2 24 49 26 49 11 51 35 51 47 51	29 47 52 58 37 14 38 17 55 18 35 22 27 23	57 58 59 60 61 63 64 66 66 69 70 73 74 74

RANKING FOR HAZARDOUS AIR POLLUTANTS (

(cont.)

Pollutant	Toxicity Score	Emissions (lbs/yr)	Toxicity Rank	Emissions Rank	Tota
Ethyl Benzene	12.5	80	18	56	74 76
Furfural	9.2	7,110	40	36	80
Barium	9.4	2,156	36	44	
Mercury	12.3	16	24	57	81
Bipheny1	8.0	4,510	43	· 40	83
Cyanide	8.5	2,955	41	42	83
p-Nitrophenol	10.5	228	31	53	84
Methyl Isobutyl Ketone		14,045	57	31	88
Ethanolamine	5.5	3,318	51	41	92
1,2 Dichlorobenzene	5.5	2,400	51	43	94
Tetrahydrofuran	8.0	400	43	51	94
Oxalic Acid	7.0	868	47	48	95
Titanium Oxide	4.0	5,240	57	39	96
Acetic Anhydride	5.5	1,403	51	46	97

SUPERTOXICS

Acrolein Beryllium Arsine Benzotrichloride Aldicarb Cadmium Bis (chloromethyl) ether Arsenic Trioxide Cyanide Hydrogen Cyanide Dioxin Methyl Isocyanate Nitrogen Mustard Phosphorus Tetrachlordibenzofuran Thallium ** APPENDICES A **

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APPENDIX A HAZARDOUS AIR POLLUTANT CODES

MNEUMONIC	SUBSTANCE	CAS #
ACETAL	Acetaldehyde	75 - Ø7-Ø
ACETAM	Acetamide	60-35-5
ACETIC	Acetic anhydride	108-24-7
ACETON	Acetone	67-64-1
ACETYL	2-Acetylaminofluorene	53-96-3
ACROLE	Acrolein	107-02-8
ACRYLA	Acrylamide	79-06-1
ACACID	Acrylic acid	79-10-7
AETHYL	Acrylic acid, Ethyl ester	140-88-5
ACRYLO	Acrylonitrile	107-13-1
ALDICA	Aldicarb	116-06-3
ALLYLC	Allyl chloride	107-05-1
ALPHAB	Alpha benzene hexachloride	319-84-6
AMINOD	p-Aminodiphenyl	92-67-1
ANILIN	Aniline and salts	62-53-3
ANISID	p-Anisidine	104-94-9
ANTIMO	Antimony (dust and salts) as Sb	7440-36-0
ARSINE	Arsine	7784-42-1
ARSINE	Arsenic (dust and salts) as As	7440-38-2
ASBEST	Asbestos	1332-21-4
AURAMI	Auramine (technical grade)	2465-27-2
BARIUM	Barium (dust and salts) as Ba	7440-39-3
BENZEN	Benzene	71-43-2
BENZID	Benzidine	92-87-5
BENZOA	Benzo(a) pyrene	50-32-8
BENZOT	Benzotrichloride	98-07-07
BENZYL	Benzyl chloride	100-44-7
BERYLL	Beryllium (dust and salts) as Be	7440-41-7
BETAPR	Beta-Propiolactone	57-57-8
BIPHEN	Biphenyl	92-52-4
BETHER	Bis(chloromethyl) ether	542-88-1
BPHTHA	Bis(2-ethylhexyl) phthalate	117-81-7
BROMIN	Bromine	7726-95-6
BUTADI	1,3,-Butadiene	106-99-0
BUTANE	Butanethiol	109-79-5
BUTANO	Butanol (n-Butyl Alcohol)	71-36-3
BACETA	n-Butyl acetate	123-86-4
BACEIA BAMINE	n-Butylamine	109-73-9
CADMIU	Cadmium (dust and salts) as Cd	7440-43-9
CTETRA	Carbon Tetrachloride	56-23-5
CDISUL	Carbon disulfide	75 - 15-Ø
CHLORI	Chlorine	7782-50-5
CDIOXI	Chlorine dioxide	10049-04-4
CACETO	2-Chloroacetophenone (Phenacylchloride)	532-27-4
CANILI	p-Chloroaniline	106-47-8
CHFORM	Chloroform	67-66-3
CMETHY	Chloromethyl methyl ether	107-30-2
CNITRO	p-Chloronitrobenzene	100-00-5
CPRENE	Chloroprene	126-99-8
CHROMI	Chromium (VI) insoluble compounds	7440-47-3
CHRYSE	Chrysene	218-01-9
COBALT	Cobalt (dust and salts) as Co	7440-48-4
COPPER	Copper (fumes, dusts & mists) as Cu	7440-50-8
CRESOL	Cresol (all isomers)	1319-77-3
CYANIM	Cyanimide	420-04-2
CTURTLI	CT CLITITITAC	-120 01-2

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CYANIK	Cyanic acid (K salt)	59Ø-28-3
CYANIN	Cyanic acid (Na salt)	917-61-3
CYANID	Cyanides (as Cn)	57-12-5
CYANOA	Cyanoacetamide	107-91-5
		460-19-5
CYANOG	Cyanogen	
CYCLOH	Cyclohexane	110-82-7
DIAMIN	2,5-Diaminotoluene	95-70-5
DIAZOM	Diazomethane	334-88-3
DETHAN	1,2-Dichlorethane	107-06-2
DBENZE	1,2-Dichlorobenzene	95-50-1
	1,2-Dichloropropane	78-87-5
DPROPA		91-94-1
- DBENZI	3,3-Dichlorobenzidine	
DPHTHA	Diethyl phthalate	84-66-2
DSULFA	Diethyl sulfate	64–67–5
DIISOC	Diisoctyl phthalate	27554-26-3
DIISOD	Diisodecyl phthalate	26761-40-0
DBENDI	3,3-Dimethoxybenzidine (o-dianisidine)	119-90-4
DIMHYD	1,1-Dimethyl hydrazine	57-14-7
		77-78-1
DSULFT	Dimethyl sulfate	
DAMINO	Dimethylaminoazobenzene	60-11-7
DCARBA	Dimethylcarbamyl chloride	- 79–44–7
DINITR	m-Dinitrobenzene	99-65-Ø
DIOXAN	1,4-Dioxane	123-91-1
TDIOXI	Total Dioxins	1746-01-6
DIPHYD	Diphenylhydrazine	122-66-7
DMETHA	Diphenylmethane 4,4-di-isocyanate(MDI)	101-68-8
	Epichlorohydrin	106-89-8
EPICHL		75-56-9
EPOXYP	Epoxypropane (Propylene oxide)	
ETHANE	Ethanethiol	75-08-1
ETHANO	Ethanolamine	141-43-5
EACETA	Ethyl acetate	141-78-6
EBENZE	Ethyl benzene	100-41-4
ECHLOR	Ethyl chloride	75-00-3
EETHER	Ethyl ether	6Ø-29-7
ETHYLE	Ethylene	74-85-1
	Ethylene glycol ethyl ether	110-80-5
EGLYCO	Ethylene oxide	75-21-8
EOXIDE		151-56-4
EIMINE	Ethyleneimine (Aziridine)	
FLUORI	Fluorine	7782-41-4
FORMAL	Formaldehyde (gas)	5Ø-ØØ-Ø
FORMAM	Formamide	75-12-7
FORMIC	Formic acid	64-18-6
FURFUR	Furfural	98-Ø1-1
FALCOH	Furfuryl alcohol	98-ØØ-Ø
GLYCID	Glycidaldehyde	765-34-4
	Hexachlorobutadiene	87-68-3
HBUTAD		77-47-4
HPENTA	Hexachlorocyclopentadiene	
HNAPHT	Hexachloronaphthalene	1335-87-1
HPHOSP	Hexamethylphosphoramide	68Ø-31-9
HYDRAZ	Hydrazine (and acid salts)	302-01-2
HBROMI	Hydrogen bromide	10035-10-6
HCHLOR	Hydrogen chloride	7647-01-0
HCYANI	Hydrogen cyanide	74-90-8
HSULFI	Hydrogen sulfide	7783-06-4
		123-31-9
HYDROQ	Hydroquinone (dihydroxy benzene)	111-42-2
IMINOD	2,2-Iminodiethanol	
IODINE	Iodine	7553-56-2
IACETA	Isoamyl acetate	123-92-2
IALCOH	Isoamyl alcohol	123-51-3

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ISOPHO		Isophorone	78-59-1
ISOPRO		Isopropylamine	75-31-Ø
KETENE		Ketene (unsaturated ketone)	463-51-4
LEADPB	•	Lead (dust and salts) as Pb	7439-92-1
			108-31-6
MALEIC		Maleic anhydride	
MANGAN		Manganese	7439-96-5
MELAMI		Melamine	108-78-1
MERCUR		Mercury (metal and salts) as Hg	7439-97-6
MCELLO		Methyl cellosolve	109-86-4
MCHLOR		Methyl chloride	74-87-3
MEKETO		Methyl ethyl ketone (MEK)	78-93-3
MIODIN		Methyl iodine	74-88-4
MISOCY		Methyl isocyanate	624-83-9
		Methyl mercaptan	74-93-1
MMERCA			80-62-6
MMETHA		Methyl methacrylate	
MISOBU		Methyl-iso-butylketone	108-10-1
MMETHY		Methylchloromethylether	107-30-2
MECHLO		Methylene Chloride	75-09-2
MEDIAN		4,4-Methylene-dianiline	101-77-9
MHYDRA		Methylhydrazine	60-34-4
MONOCH		Monochlorobenzene (chlorobenzene)	108-90-7
NNAPHT		n-phenyl-beta-naphthylamine	135-88-6
NALINE		Napthalene	91-20-3
		Napthylamine(alpha)	134-32-7
NAMINA			91-59-8
NAMINB		Napthylamine (beta)	744Ø-Ø2-Ø
NICKEL		Nickel (dust and salts) as Ni	
NITRIC		Nitric acid	7697-37-2
NITROA		p-Nitroaniline	100-01-6
NBENZE		Nitrobenzene	98-95-3
NBIPHE		4-Nitrobiphenyl	92-93-3
NMUSTA		Nitrogen mustard	51-75-2
NGLYCE		Nitroglycerine	55-63-Ø
NPHENO		p-Nitrophenol	100-02-7
NPROPA		1-Nitropropane	108-03-2
NSOMET		Nitroso-n-methylurea	684-93-5
			62-75-9
NSODIM		n-Nitrosodimethylamine	
NSOMOR		n-Nitrosomorpholine	59-89-2
NSOPHE		p-Nitrosophenol	104-91-6
NTOLUM	•	m-Nitrotoluene	99-08-1
NTOLUP		p-Nitrotoluene	99-99-ø
OCTACH		Octachloronaphthalene	2234-13-1
OXALIC		Oxalic acid	144-62-7
PCPHEN		Pentachlorophenol (PCP)	87-86-5
PHENOL		Phenol	108-95-2
PDIAMI		p-Phenylenediamine	106-50-3
		Phenylhydrazine	100-63-0
PHYDRA			75-44-5
PHOSGE		Phosgene	7723 - 14 - Ø
PHOSPH		Phosphorus	
PICRIC		Picric acid	88-89-1
PCBIPH		Polychlorinated byphenyls (PCBs)	11097-69-1
PROPAN		1,3-Propane sultone	1120-71-4
PROPYL		Propyleneimine	75-55-8
PYRIDI		Pyridine	110-86-1
QUINOL		Quinoline	91-22-5
QUINON		Quinone	106-51-4
RESORC		Resorcinol	108-46-3
		Rotenone	83-79-4
ROTENO		Selenium (dust and salts) as Se	7782-49-2
SELENI SOXIDE			96-Ø9-3
		Styrene oxide	ラローロラーン

STYREN TEREPH TETHYL TDIBEN TETHAN TFURAN TFURAN THALLI TITANI TOLUEN TDIAMI TDIISO TOLUID TBENZE TETHA1 TETHA2 TRICHL TURPEN URETHA VBROMI VCHLOR VCYCLO VFLUOR VIDENE XYLENE XYLIDI ZINCZN	Styrene, monomer Terephthalic acid Tetrachlorethylene (perchlorethylene) 2,3,7,8-Tetrachlorodibenzofuran 1,1,2,2-Tetrachloroethane Tetrahydrofuran Thallium (dust and salts) as Tl Titanium oxide Toluene 2,4-Toluene-diamine 2,4-Toluene-di-isocyanate o-Toluidine 1,2,4-Trichlorobenzene 1,1,1-Trichloroethane, methyl chloroform 1,1,2-Trichloroethane Trichloroethylene Turpentine Urethane Vinyl Bromide Vinyl Bromide Vinyl chloride Vinyl chloride Vinyl fluoride Vinyl fluoride Vinyl fluoride (1-1-Dichloroethene) Xylene (all isomers) Xylidine Zinc (dust and salts) as Zn	100-42-5 100-21-0 127-18-4 51207-31-9 79-34-5 109-99-9 7440-28-0 13463-67-7 108-88-3 95-80-7 584-84-9 95-53-4 120-82-1 71-55-6 79-00-5 79-01-6 8006-64-2 51-79-6 593-60-2 75-01-4 106-87-6 75-02-5 75-35-4 1330-20-7 1300-73-8 7440-66-6
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** APPENDIX B **

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APPENDIX B

SUMMARY OF HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE

		EMISSIONS (POUNDS	S)
POLLUTANT	POINT SOURCE	AREA SOURCE	TOTAL
1 Toleran	0 450 071	F0 760	0 511 501
1. Toluene	2,459,231	52,360	2,511,591
2. 1,1,1, Trichloroethane	1,496,539	0	1,496,539
3. Chlorine	1,337,003	0	1,337,003
4. Xylene	1,091,274	0	1,091,274
5. Acetone	804,521	0	804,521
6. Tetrachlorethylene	144,450	563,600	708,050
7. Methyl Cellosolve	632,248	0	632,248
8. Methyl Ethyl Ketone	512,611	0	512,611
9. Methyl Mercaptan	428,940	0	428,940
10. Chlorine Dioxide	296 , 787	0	296 , 787
11. Methylene Chloride	288,569	0	288,569
12. Ethylene Glycol Ethyl Ethe	er 277,118	0	277,118
13. Hydrogen Sulfide	269,912	0	269,912
14. N-Butyl Acetate	253,563	0	253,563
15. Lead	630	202,800	203,430
16. Methyl Metharcylate	184,550	0	184,550
17. Butanol	149,490	0	149,490
18. Diphenyl methane 4,4-di-	•		
isocyanate	146,000	0	146,000
19. Hydrogen Chloride	44,731	96,000	1,40,731
20. 1,2 - Epoxypropane	124,600	0	124,600
21. Trichloroethylene	110,600	Ō	110,600
22. Formaldehyde	79,401	0	79,401
23. Turpentine	77,130	Õ	77,130
24. Ethyl Acetate	66,537	Õ	66,537
25. Benzene	0	52, 392	52,392
26. Formic Acid	50,433	0	50,433
27. Nitric Acid	41,450	Õ	41,450
28. 1, 2 - Dichloroethane	32,400	Õ	32,400
29. Methyl Chloride	16,800	0	16,800
30. Benzo (a) Pyrene	0	15,000	15,000
Jo. Denzo (a) ryrene	v	12,000	19,000

APPENDIX B (CONTINUED)

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SUMMARY OF HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE

		EMISSIONS (POUNDS)	
POLLUTANT	POINT SOURCE	AREA SOURCE	TOTAL
31 Mathyl Isabutyl Katana	14 045	0	14 045
31. Methyl Isobutyl Ketone	14,045 13,001	0	14,045
32. Styrene	11,799		13,001
33. Napthalene 34. Bis Phthalate		0	11,799
	10,190	0	10,190
35. Manganese	9,320	0	9,320
36. Furfural	7,110	0	7,110
37. Phenol	5,810	0	5,810
38. Diethyl Sulfate	5,424	0	5,424
39. Titanium Oxide	5,240	0	5,240
40. Biphenyl	4,510	0	4,510
41. Ethanol Amine	3,318	0	3,318
42. Cyanimide	2,955	0	2,955
43. 1,2 Dichlorobenzene	2,400	0	2,400
44. Barium	2,156	0	2,156
45. Ethylene Oxide	1,535	0	1,535
46. Acetic Anhydride	1,403	0	1,403
47. Chromium	184	820	1,004
48. Zinc	946	0	946
49. Oxalic Acid	868	0	868
50. Hydrazine	740	0	740
51. Tetrahydrofuran	400	0	400
52. P-nitrophenol	228	0	228
53. Cadmium	0	200	200
54. Copper	107	0	107
55. Ethyl Benzene	80	0	80
56. Mercury	16	· 0	16
57. Epichlorohydrin	12	0	12

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** APPENDIX C **

APPENDIX C

AREA SOURCE EMISSIONS DATA

CATEGORY: DRY CLEANERS

DATA COLLECTION: Maine DEP statewide emissions inventory survey of dry cleaning establishments. 72% of respondents used a cleaning solvent of which 82% was perc.	CALCULATIONS:	
16.9% was stoddard, and .5% was freon 113. Each respondent on the avg. used 423.7 gal of perc., 86.5 gal of stoddard, and .5 gal freon 113.	Estimated use Perc. and stod.	<pre>#non-resp. x 72.6% x avg.use of responds</pre>
	Emissions VOC	#gal x wt/gal /2000 x EMF
DATA ASSUMPTIONS:		
1. Respondents and non-respondents of survey total 100% of ME dry		
cleaners.		
 Amount of solvent purchased equals amount of solvent consumed. Amounts used by units not responding corresponds to state avg of responding units. 		
 Perchlorethylene weighs 13.6 lb/gal; stoddard ~6.1 lb/gal. 		·

5. Use of freon 113 assumed to be normal.

EMISSION FACTOR(S):

Perchloroethylene----Stoddard solvent-----

2000 lb/ton consumed 2000 lb/ton consumed

TOTAL EMISSIONS:	
Perchloroethylene VOC's	281.8 tons
Stoddard solvent VOC's	25.84 tons

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EMISSION FACTOR SOURCE:

NEDS Source Classification Codes and Emission Factor Listing, EPA, Dec. 1984.

COUNTY	USE OF R UNITS PERC.	ESPONDING (GALS) STOD.	EMISSIC RESPOND VOC PERC.	DENTS (TONS)	#UNITS IN COUNTY NOT RESPONDING OR UNKNOWN	ESTIMATA USE OF N RESPONDS PERC.	ON-	EMISSIC NON-RES VOC PERC.			TOTAL STOD.VOC EMITTED (TONS
ANDROSCÓGGIN	1189	100 ing dia 200 ing dia 200 ing dia 200	8.0		7	2153.3		14.6		. 22.7	
AROOSTOOK	1525		10.3			ø	440.7	.ø	1.3	10.3	1.3
CUMBERLAND	8007	1760	54.4	5.3	11	3383.8	.0	23.0	.0	77.4	5.3
FRANKLIN	1976		13.4			ø	692.6	.ø	2.1	13.4	2.1
HANCOCK	585	702	3.9	2.1	3	922.8	.Ø	6.2	.0	10.2	2.1
KENNEBEC	2453		16.6		8	2461.0	188.8	16.7	۰5		.5
				×	_		503.7		1.5		1.5
KNOX	450		3.0		4	1230.5	251.8	8.3	.7	11.4	.7
LINCOLN			.0		1	307.6	62.9		.1		.1
OXFORD	8Ø5	972	5.4	2.9	1	307.6	62.9	2.0	.1	7.5	3.1
PENOBSCOT	2063	1325	14.0	4.0	7	2153.3	440.7	14.6	1.3	28.6	5.3
PISQUATAQUIS	245		1.6		1	307.6	62.9	2.0	.1	3.7	
SAGADAHOC			.ø	•	. 2	615.2	125.9	4.1	.3	4.1	
SOMERSET	600		4.0		2	615.2		4.1		8.2	
WALDO	600		4.0			ø	125.9	.ø	.3	4.0	
WASHINGTON	650		4.4		3	922.8	.0	6.2		10.6	
YORK	2157		14.6		9	2768.6	188.8	18.8	•2	33.4	•5 ,
							566.6		1.7		1.7
TOTALS	233Ø5	4759	158.4	14.5	59.0	18149.9	3714.8	123.4	11.3	281.8	25.8

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AREA SOURCE EMISSIONS DATA

DRY CLEANERS

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Degreasers - Area source

DATA COLLECTION:

Population data obtained from the Maine Dept. of Human Services bulletin: "Population projections of Main counties and minor civil divisions for total population - July 1 (1982-1991)."

DATA ASSUMPTIONS:

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Population projections for 1983 are representative of 1984.
 All area source degreasers are of the cold cleaning variety.

Note: A major source of cold cleaning degreasing chemicals in ME is the Safety Kleen Corp. who supplies and recycles these chemicals,of which petroleum naptha represents nearly 100%. Rich Baker of the Oil and Hazardous Waste Materials Bureau, confirms this information.

EMISSION FACTOR(S):

All VOC emissions	4 lb/cap-yr
All reactive VOC emissions	3 lb/cap-vr

CALCULATIONS

POLLUTANT	FÖRMULA				
Total VOC Total reactive VOC	Pop. x emf / 2000 Pop. x emf / 2000				

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TOTAL EMISSIONS

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VOC emissions	2297.4 tons
Reactive VOC emiss.	1723.0 tons

EMISSION FACTOR SOURCE:

EPA-450/2-77-028, Sec.4.3.2.2, Cold Cleaning Degreasing

AREA SOURCE EMISSIONS

DEGREASERS

COUNTIES AND SELECTED CITIES	POPULATION	VOC	REACTIVE VOC							
		tons	tons							
NDROSCOGGIN	100900	201.8	151.3							
Auburn	22300	44.6	33.4							
Lewiston	40650	81.3	60.9							
AROOSTOOK	90400	180.8	135.6							
Presque Isle	1Ø85Ø	21.7	16.2							
UMBERLAND	220100	440.2	330.1							
Brunswick	1755Ø	35.1	26.3							
Portland	61100	122.2	91.6							
S.Portland	23000	46.0	34.5							
Westbrook	1525Ø	30.5	22.8							
FRANKLIN	28500	57.0	42.7							
IANCOCK	42700	85.4	64.0	·						
(ENNEBEC	11285Ø	225.7	169.2							
Augusta	22050	44.1	33.0							
Waterville	17050	34.1	25.5							
NOX	34050	68.1	51.0							
LINCOLN	26150	52.3	39.2							
DXFORD	48800	97.6	73.2							
PENOBSCOT	141150	282.3	211.7							
Bangor	33350	66.7	50.0					,		
Brewer	8600	17.2	12.9							
PISQUATAQUIS	1785Ø	35.7	26.7							
SAGADAHOC	28050	56.1	42.0							
Bath	995ø	19.9	14.9							
SOMERSET	47050	94.1	70.5							
VALDO	28500	57.0	42.7							
WASHINGTON	3355Ø	67.1	50.3							
YORK	148100	296.2	222.1							
Biddeford	20400	40.8	30.6							
Saco	1265Ø	25.3	18.9							
Sanford	19450	38.9	29.1							
TOTALS	1148700	2297.4	1723.0	.ø						

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Residential Open Burning (Dumps)

DATA COLLECTION:

1

Data concerning which towns employ open burning of solid waste and and the town's population obtaine from the Maine Dept. of Envt'l-Protection's Bureau of Land Quality Control (based on 1984 inspection reports on all Maine dumps).

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CALCULATIONS

	متبالة مشال جيسد بالثان فبلب بجيب ملتك فيشد جيجه نشلك وونك جويه فيجه إلجه ويجه بالحد فيده ويحد ويده
Pollutants	Pop. x 6.66t/1000 per wk x 52 wk/yr x 60% x EMF ton/20001b
B(a)P	Amt. Part. x .00012

DATA ASSUMPTIONS:

1 -31 -

- 1. Assumed solid waste disposal of 6.66 tons/thous people per week.
- 2. Assumed 60% burning of solid waste disposed.

- 3. Assumed brush burning is insignificant at Maine dumps.
- 4. Assumed 1984 dump conditions are representative for 1983.

EMISSION FACTOR(S):

Particulates	16	lb/ton
Sulfur oxides	1	lb/ton
Carbon monoxide	85	lb/ton
VOC methane	13	lb/ton
VOC non-methane	3Ø	lb/ton
Nitrogen oxides	6	lb/ton
Benzo(a)pyrene	.00012	lb/ton

TOTAL EMISSIONS

Particulates	210.2	tons
Sulfur oxides	13.1	tons
Carbon monoxide	1117.0	tons
VOC methane	170.8	tons
VOC non-methane	394.2	tons
Nitrogen oxides	78.8	tons
Benzo(a)pyrene	.0252	tons

EMISSION FACTOR SOURCE:

AP-42, Sec.2.4: Open Burning, Table 2.4-1. Maine DEP, Bureau of Land Quality Control and conversation with DHS toxicologist Norm Anderson Jan.30, 1985.

AREA SOURCE EMISSIONS

RESIDENTIAL OPEN BURNING

(DUMPS)

COUNTIES AND SELECTED CITIES		POP. AREAS OPEN BURN SOL. WASTE	PARTICULAT tons	SULFUR OXIDES tons	CARBON MONOXIDE tons	VOC METHANE tons	VOC NON-METHAN tons	NITROGEN OXIDES tons	B(a)P tons	tons
ANDROSCOGGIN	100900		12.3	.77	65.5	10.0		4.6	.0014	
Auburn	22300		.ø	-00	.0	.ø		.0	.0000	
Lewiston	40650		.ø	.00	.0	.0		.0	.0000	
AROOSTOOK	90400			1.42		18.5		8.5	.0027	
Presque Isle	10850		.ø	.00	.ø	.ø		.0	.0000	
CUMBERLAND	220100		8.2	.51		6.7		3.0	.0009	
Brunswick	17550		.ø	.00	.0	.0		.0	.0000	
Portland	61100		.0	.00		.0		-0	.0000	
S.Portland	23000		.ø	.00	.ø	.ø		.0	.0000	
Westbrook	15250		.ø	.00		.0		.Ø~		
FRANKLIN	28500			.63		8.2		3.8	.0012	
HANCOCK	42700			.62		8.0		3.7	.0011	
KENNEBEC	112850		9.2	. 57		7.4		3.4	.0011	
Augusta	22050		.ø	.00		.0		.0	.0000	
Waterville	17050		.Ø	.00	.ø	.ø		.0	.0000	
KNOX	34050			.10		1.3		6	.0001	
LINCOLN	26150			.63				3.7	.0012	
OXFORD	48800	15560	25.8	1.61	137.4	21.0	48.4	9.6	.0031	
PENOBSCOT	141150	19155	31.8	1.99	169.1	25.8	59.7	11.9	.0038	
Bangor	33350		.ø	.00	.ø	.0	.0	.0	.0000	
Brewer	8600		.Ø	.00	.ø	.Ø	.0	.0	.0000	
PISQUATAQUIS	17850	9324	15.4	.96	82.3	12.5	29.0	5.8	.0018	
SAGADAHOC	28050	735	1.2	.07	6.4	.9	2.2	.4	.0001	
Bath	9950		.ø	.00	.0	.0	.0	.0	.0000	
SOMERSET	47050	8399	13.9	.87	74.1	11.3	26.1	5.2	.0016	
WALDO	28500	7787	12.9	.80	. 68.7	10.5	24.2	4.8	.0015	
WASHINGTON	33550	5434	9.0	- 56	47.9	7.3	16.9	3.3	.0010	
YORK	148100	9213	15.3	.95	81.3	12.4	28.7	5.7	.0018	
Biddeford	20400		.0	.00	.0	.0	.ø	.ø	.0000	
Saco	12650		.ø	.00	.ø	.0	.0	.Ø	.0000	
Sanford	19450		.0	.00	.0	.0	0	.0	.0000	
TOTALS	1148700	126487.0	210.2	13.14	1117.0	170.8	394.2	78.84	.0252	.ø
		available.				Total VOC	565.0			

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AREA SOURCE EMISSIONS DATA

CATEGORY:

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Agricultural burning - Blueberries

DATA COLLECTION:

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Contact was made with Ed McLaughlin and David Yarborough, both asst. scientists with the University of Maine Plant and Soil Dept.

DATA ASSUMPTIONS:

 It was determined that the only significant agricultural field burning in Maine during 1983 occurred from blueberry field prep.
 Assumed that in any year 1/2 of the blueberry acreage is burned. The USAGE column represents this 1/2.

3. Assumed that of Washington County's burnable acreage, 25% was pruned.

4. The ratio of BaP/particulates of woodsmoke emissions is .00012.

CALCULATIONS

	······································
Pollutants	Ac x flf x emf / 2000
B(a)P	Part.emiss. x .00012

EMISSION FACTOR(S):

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1110101(0)	-
]	Particulates
(Carbon monoxide
7	VOC methane
7	VOC non-methane
1	Benzo(a)pyrene
	ing Factor

21 lb/ton 117 lb/ton 5.4 lb/ton 18 lb/ton .00012 x part. emiss. 2 tons/acre

TOTAL EMISSIONS

1. Sec. 2. Sec. 1. Sec

Particulates	523.9 tor	าร
Carbon monoxide	2919.Ø tor	ıs
VOC methane	134.7 tor	າຣ
VOC non-methane	449.Ø tor	າຣ
Benzo(a)pyrene	.0628 tor	าร

EMISSION FACTOR SOURCE:

AP-42: Open Burning of Agricultural materials: field crops unspecified (Sec.2.4) and conservation with DHS toxicologist Norm Anderson Jan.30, 1985. Acres of land in blueberry production supplied by David Yarborough. AREA SOURCE EMISSIONS

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AGRICULTURAL BURNING -BLUEBERRY

COUNTIES AND SELECTED CITIES	POPULATION	USAGE 1/2 tot ac burned/yr	PARTICULAT tons	CARBON MONOXIDE tons	VOC METHANE tons	VOC NON-METHAN tons	B(a)P tons			
ANDROSCOGGIN	100900	N/A								
Auburn	22300						•			
Lewiston	40650							,		
AROOSTOOK	90400		.33	1.87	.Ø8	.28	.00004			
Presque Isle	10850									
CUMBERLAND	220100		10.12	56.39	2.60	8.67	.00121			
Brunswick	17550									
Portland	61100				•					
S.Portland	23000	11								
Westbrook	15250					•				
FRANKLIN	28500	11								
HANCOCK	42700	5021	105.44	587.45	27.11	90.37	.01265			
KENNEBEC	112850	180	3.78	21.06	.97	3.24	.00045			
Augusta	22050	· N/A								
Waterville	17050									
KNOX	34050		53.65	298.93	13.79	45.99	.00643			
LINCOLN	26150			64.81	2,99		.00139			
OXFORD	48800									
PENOBSCOT	141150		1.36	7.60	.35	1,17	.00016			
Bangor	33350		1.50							
Brewer	8600									
PISQUATAQUIS	17850									
SAGADAHOC	28050									
Bath	20050 995Ø				•					•
SOMERSET	47050									
WALDO	28500		39.75	221.48	10.22	34.07	.00477			
WASHINGTON	33550			1645.37	75.94		.03543			
YORK	148100			14.04	.64					
Biddeford		-	2.52	14.04	.04	2.10	.00030			
	20400	,								
Saco	12650									
Sanford	19450									
TOTALS	1148700	24949.0	.523.9	2919.0	134.7	449.0	.ø628	.ø	.ø	.ø
Individual city					Total VOC	583.8				

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Forest Wildfire

DATA COLLECTION:

.

Forest acres burned by county in 1983 obtained from Maine Forest Service (Dept. of Conservation).

CALCULATIONS

POLLUTANT	FORMULA					
Particulates	Use x llt/ac x EMF/2000					
Carbon monoxide	Use x llt/ac x EMF/2000					
Total hydrocarbon	Use x llt/ac x EMF/2000					
Nitrogen oxides	Use x llt/ac x EMF/2000					
Benzo (a) pyrene	Amt part x EMF					

DATA ASSUMPTIONS:

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Fuel loading assumed to be 11 tons/acre (source: AP-42 sec.11.1). The ratio of BaP/part. of wood smoke emissions is .00012. This figure x the amt of particulates from forest wildfire emissions = the amt of BaP.

EMISSION FACTOR(S):

Particulates------Carbon monoxide------Total hydrocarbons----Nitrogen oxides------Benzo (a) pyrene----- 17 lbs/ton 140 lbs/ton 24 lbs/ton 4 lbs/ton .00012 x amt of part. lbs/ton

TOTAL EMISSIONS

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Particulates	83.4 tons
Carbon monoxide	687.4 tons
Total hydrocarbons	117.8 tons
Nitrogen oxides	19.6 tons
Benzo (a) pyrene	.Øl tons

EMISSION FACTOR SOURCE: -

AP-42 Forest Wildfires (Sec.11.1) and conversation with DHS toxicologist Norm Anderson, Jan.30, 1985, (for BaP).

COUNTIES AND SELECTED CITIES	POPULATION	USAGE (1983 acre burned)	PARTICULAT	CARBON MONOXIDE tons	TOTAL HYDROCARB. tons	NITROGEN OXIDES tons	B(a)P tons					
NDROSCOGGIN Auburn	100900 22300	21.5	2.0	16.5	2.8	.47	.0002			وی ہیں ہیں جب ختا بنا ہیں ہے۔		
Lewiston ROOSTOOK Presque Isle	40650 90400 10850	79.2	7.4	60.9	10.4	1.74	.0008					
UMBERLAND Brunswick Portland S.Portland	220100 17550 61100 23000	34.6	3.2	26.6	4. 5	.76	.0003					
Westbrook	1525Ø 285ØØ	9.9	٩	7.6	1.3	.21	.0001				•	
FRANKLIN HANCOCK	42700	53.2		40.9		1.17	.0005					
ENNEBEC	42700 112850	28.9		40.9		.63	.0003		•			
Augusta Waterville	22050 17050	20.9	21		2.0	-05	•0000					
NOX	34050	69.1	6.4	53.2	9.1	1.52	.0007					
LINCOLN	26150	21.0		16.1		.46	.0002					
DXFORD	48800	28.1		21.6		.61	.0003					
PENOBSCOT Bangor Brewer	14115Ø 3335Ø 86ØØ	108.9	10.1	83.8	14.3	2.39	.0012					
PISQUATAQUIS SAGADAHOC Bath	1785Ø 28Ø5Ø 995Ø	32.4 22.8		24.9 17.5		.71 .50	.0003 .0002					
SOMERSET	47050	28.1	2.6	21.6	3.7	.61	.0003			,		
VALDO	28500	16.6		12.7	2.1	.36	.0001					
VASHINGTON	3355Ø	290.0		223.3	38.2	6.38	.0032					
ORK Biddeford	148100 20400	48.5	4.5	37.3	6.4	1.06	.0005					
Saco	12650											
Sanford	19450											
TOTALS	1148700	892.8	83.4	687.4	117.8	 19.64	.0100	 Ø	 Ø		I	

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FOREST WILDFIRE

Individual city data not available.

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AREA SOURCE EMISSIONS

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Architectural Surface Coating

DATA COLLECTION:

Population figures obtained from "Population projection of Maine counties and minor civil divisions for total populations, 1982-1991" by the ME Dept. of Human Services.

Total VOC's	Population x EMF/2000

DATA ASSUMPTIONS:

37-

1. Population projections for 1983 were representative for the year 1983.

2. Usage of architectural surface coatings is proportional to the population.

EMISSION FACTOR(S):

VOC's-----

4.6 lb/cap-yr

TOTAL EMISSIONS:

CALCULATIONS:

Total VOC's	2642.01 tons

EMISSION FACTOR SOURCE:

AP-42, 4.2.1 Nonindustrial surface coating.

AREA SOURCE EMISSIONS DATA

COUNTY	POPULATION	VOC EMISSIONS (TONS/YR)
ANDROSCOGGIN	100900	232.0
AROOSTOOK	90400	207.9
CUMBERLAND	220100	506.2
FRANKLIN	28500	65.5
HANCOCK	42700	98.2
KENNEBEC	112850	259.5
KNOX	34050	78.3
LINCOLN	26150	60.1
OXFORD	48800	112.2
PENOBSCOT	141150	324.6
PISQUATAQUIS	17850	41.0
SAGADAHOC	28Ø5Ø	64.5
SOMERSET	47050	108.2
WALDO	28500	65.5
WASHINGTON	3355Ø	77.1
YORK	148100	340.6

TOTALS 1148700 2642.0

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AREA SOURCE EMISSIONS

DATA

CATEGORY:

Highway line and bridge painting

DATA COLLECTION:

All data obtained from the Maine Dept. of Transportation by personal communication. Line paint data was obtained from Dick Weeks of Division of Traffic Engineering, Jan.30,1985. Data for bridge painting obtained from John Butts and Ron Cyr of the Design Division and Bridge Maintenance Division respectively, Jan.31, 1985.

CALCULATIONS

POLLUTANT	FORMULA
VOC	Usage x 5.6 / 2000

DATA ASSUMPTIONS:

One gallon of solvent-based paint contains 5.6 lbs of VOC.
 100% of the solvent applied to the highway or bridge surface is emitted.

Note: Highway paint solvent primarily consist of <u>toluene</u> with small amounts of methylene chloride, xylene, hexane, vm+p naphtha, and 1,1 1,trichloroethane. Bridge paint primarily utilizes mineral spirits.

EMISSION FACTOR(S):

Total VOC

5.6 lb/gal

TOTAL EMISSIONS

VOCroad paint	422.3 tons	
VOCbridge paint	38.3 tons	
Total VOC	460.8 tons	

EMISSION FACTOR SOURCE:

EPA--NEDS Source Classification Codes and Emission Factor Listing--Surface Coating Operations, p96-97.

REA SOURCE E	MISSIONS			HIGHWAY LIN	E AND BRIDG	E PAINTING				
COUNTIES	POPULATION	LINE PAINT USAGE gals	LINE PAINT VOC tons	BRIDGE PNT USAGE gals	BRIDGE PNT VOC tons	TOTAL VOC tons				
NDROSCOGGIN	100900	4570	12.7	16	.04	12.8				
ROOSTOOK	90400	19806	55.4		.00	55.4				
UMBERLAND	220100	10665	29.8	836Ø	23.40	53.2				
			·							
RANKLIN	28500				.00	25.5				
IANCOCK IENNEBEC	42700 112850		34.1 25.5		.00 14.11	34.1 39.7				
NOX	34050	4570	12.7		.00	12.7				
INCOLN	26150				.00	17.0				
XEORD	48800				.00	34.1				
ENOBSCOT	141150	15236	42.6	24Ø	.67	43.3				
ISQUATAQUIS	17850				.06	12.8				
AGADAHOC	28050	3047	8.5		.00	8.5				
OMERSET	47050	12188	34.1		.00	34.1				
ALDO	28500				.00	17.1				
ASHINGTON	33550				.00	29.9				
ORK	148100	10665	29.8		.00	29.9				
COUNTY OF USE	_ ,			2965	8.30	.ø				
	- '			2905	0.10	• 10				
TOTALS	1148700	150828.0	422.3	13680.0	38.3	 460.8	.ø	.ø	.ø	 .ø

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Individual city data not available. *County breakdown of bridge paint used for maintenance is not available.

AREA SOURCE EMISSIONS DATA

CATEGORY: Printing Industry

DATA COLLECTION:

Population data obtained from the Maine Dept. of Human Services Bulletin, "Population projections of Maine counties and minor civil division for total population - July (1982-1991)."

CALCULATIONS

POLLUTANT	FORMULA					
VOC non-methane	EMF * Population					

DATA ASSUMPTIONS:

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1. Population projections for 1983 were representative for the year 1983.

EMISSION FACTOR(S):

Non-Methane VOC -----

.8 lb/yr/capita

TOTAL EMISSIONS

VOC non-methane 918960 lb/yr

EMISSION FACTOR SOURCE: AP-42, Sec. 4.9, table 4.9-2

AREA SOURCE EMISSIONS DATA

PRINTING INDUSTRY

.

COUNTIES AND SELECTED CITIES	POPULATION	FACTOR lb/yr/cap	TOTAL VOC non-meth. lbs		POLL	UTANTS				
ANDROSCOGGIN	100900	.8	80720							Na hay tan tan ang
Auburh	22300	.8								
Lewiston	40650	.8								
AROOSTOOK	90400	.8								
Presque Isle	10850	.8								
CUMBERLAND	220100	.8	176080							
Brunswick	17550	.8	14040							
Portland	61100	.8	4888 Ø							
S.Portland	23000	.8								
Westbrook	15250	8	12200							
RANKLIN	28500	.8								
HANCOCK	42700	.8	34160							
KENNEBEC	112850	.8	90280							
Augusta	22050	.8	1764Ø							
Waterville	17050	.8	1364Ø							
KNOX	34050	.8	27240							
LINCOLN	26150	-8	20920							
OXFORD .	48800	-8	39040							
PENOBSCOT	141150	.8	112920							
Bangor	3335Ø	.8	26680							
Brewer	8600	.8								
PISQUATAQUIS	1785Ø	-8	14280			,				
SAGADAHOC	28050	-8	22440					·		
Bath	995Ø	.8	7960			,				
SOMERSET	47050	.8								
WALDO	28500	.8								
WASHINGTON	33550	-8			•					•
YORK	148100	.8	118480							•
Biddeford	20400	.8								
Saco	12650	.8								
Sanford	19450	•8	15560							
TOTALS			918960	 Ø	 Ø		 Ø	ø	 Ø	 Ø

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Residential Wood Combustion

DATA COLLECTION:

State consumption of wood for Maine for 1982 was obtained from the Maine Office of Energy Resources 'Comprehensive Energy Plan (1983).' Population figures obtained from "Population Projections of Maine counties and minor civil divisions for total population, 1982-1991," by the Maine Dept. of Human Services.

DATA ASSUMPTIONS:

-43-

- 1. 2.5 tons/cord of wood.
- 2. State consumption of wood apportioned to the county level on the basis of county population as a % of the state population total.
- 3. 1982 wood usage data is representative of 1983 wood usage.
- 4. All wood burned in wood stoves.

EMISSION FACTOR(S):

•			· · ·	
	Particulates	42	lb/ton	
	SOx	-4	lb/ton	
	NOx	2.8	lb/ton	
	C0	260	lb/ton	
	P.O.M	.3	gm/kg	
	B(a)P	.0025	gm∕kg	
	VOC methane	1	lb/ton	
	VOC non-methane	100	lb/ton	

TOTAL EMISSIONS

Particulates	62632.4	tons
SOx ·	596.4	tons
NOx	4175.4	tons
CO .	387724.9	tons
P.O.M.	894.7	tons
B(a)P	7.4	tons
VOC methane	1491.2	tons
VOC non-meth	ane 149124.9	tons

EMISSION FACTOR SOURCE:

AP-42, Particulates, SOx, NOx, CO, VOC: meth. & non-meth (res.wood stoves) J.A.Cooper, JAPCA, Avg'80 "Environmental impact of Res. Wood Combustion and its implications- POM and BaP, Res. Wood Stoves.

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CALCULATIONS

Usage =	
State Usage *	1193000 cds
Tons per Cord	2.5 ton/cd
Pollutants	
(In lbs)	Usage * EMF / 2000 lb/t
(In kg)	Usage * EMF / 1000 g/kg

AREA SOURCE EMISSIONS

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RESIDENTIA L WOOD COMBUSTION

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COUNTIES AND SELECTED CITIES	POPULATION	USAGE	PARTICULAT	SOx	NOx	C0	P.O.M.	B(a) P	VOC methane	VOC non-meth.
		tons	tons	tons	tons	tons	tons	tons	tons	tons
ANDROSCOGGIN	100900	261978.1	5501.5	52.3	366.7	34057.1	78.5	.65	130.9	13098.9
Auburn	22300	57900.0		11.5	81.0	7527.0	17.3	.14	28.9	2895.0
Lewiston	40650	105544.2	2216.4	21.1	147.7	13720.7	31.6	.26	52.7	5277.2
AROOSTOOK	90400	234715.7	4929.0	46.9	328.6	30513.0	70.4	.58	117.3	11735.7
Presque Isle	10850.	28171.0		5.6	39.4	3662.2	8.4	.07	14.0	1408.5
CUMBERLAND	220100	571470.5	12000.8	114.2	800.0	74291.1	171.4	1.42	285.7	28573.5
Brunswick	17550	45567.0		9.1	63.7	5923.7	13.6	.11	22.7	2278.3
Portland	61100	158640.8	3331.4	31.7	222.0	20623.3	47.5	.39	79.3	7932.0
S.Portland	23000	59717.5	1254.0	11.9	83.6	7763.2	17.9	.14	29.8	2985.8
Westbrook	15250	39595.3		7.9	55.4	5147.3		.09	19.7	1979.7
FRANKLIN	28500	73997.7	1553.9	14.7	103.5	9619.7	22.1	.18	36.9	3699.8
HANCOCK	42700	110866.8	2328.2	22.1	155.2	14412.6	33.2	.27	55.4	5543.3
KENNEBEC	112850	293005.2	6153.1	58.6	410.2	38090.6	87.9	.73	146.5	14650.2
Augusta	22050	57250.9		11.4	80.1	7442.6	17.1	.14	28.6	2862.5
Waterville	17050	44268.8	929.6	8.8	61.9	5754.9	13.2	.11	22.1	2213.4
KNOX	34050	88407.8		17.6	123.7	11493.0	26,5	.22	44.2	4420.3
LINCOLN	26150	67896.2	1425.8	13.5	95.0	8826.5	20.3	.16	33.9	3394.8
OXFORD	48800	126704.9		25.3	177.3	16471.6	38.0	.31	63.3	6335.2
PENOBSCOT	141150	366483.7		73.2	513.0	47642.8	109.9	.91	183.2	18324.1
Bangor	33350	86590.3		17.3	121.2	11256.7	25.9	.21	43.2	4329.5
Brewer	8600	22329.1		4.4	31.2	2902.7	6.6	.05	11.1	1116.4
PISOUATAOUIS	17850	46345.9		9.2	64.8	6024.9	13.9	.11	23.1	2317.2
SAGADAHOC	28050	72829.3		14.5	101.9	9467.8	21.8	.18	36.4	3641.4
Bath	9950	25834.3		5.1	36.1	3358.4	7.7	.06	12.9	1291.7
SOMERSET	47050	122161.2		24.4	171.0	15880.9	36.6	.30	61.0	6108.0
WALDO	28500	73997.7		14.7	103.5	9619.7	22.1	.18	36.9	3699.8
WASHINGTON	33550	87109.6		17.4	121.9	11324.2	26.1	.21	43.5	4355.4
YORK	148100	384528.8		76.9	538.3	49988.7	115.3	.96		19226.4
Biddeford	20400	52966.8		10.5	74.1	6885.6	15.8	.13	26.4	2648.3
Saco	12650	32844.6		6.5	45.9	4269.8	9.8	.08	16.4	
Sanford	. 19450	50500.2		10.1	70.7	6565.0	15.1	.12		
TOTALS	1148700	2401086.2	62632.4	596.4	4175.4	387724.9	894.7	 7 . 45	1491.2	149124.9

AREA SOURCE EMISSIONS DATA

CATEGORY: Waste oil

DATA COLLECTION:

CALCULATIONS:

TOTAL EMISSIONS:

Chlorine

Note: PCB 5-7 ppm in fuel resulted in no detectable

Tot.pop/cty.pop x use x 8 x EMF x .01/2000

All other pollutants Tot.pop/cty.pop x use x 8 x EMF / 2000

DATA ASSUMPTIONS: 1. Assume waste oil weighs 81b/gal. 2. Assume even distribution by population.

EMISSION FACTOR(S):

_							
	Lead	.001000	% Waste	oil		····	
	Chlorine	.004000	% Waste	oil	Lead	11.99	tons
	Arsenic	.000018	% Waste	oil	Chlorine	47.99	tons
	Cadmium	.000010	% Waste	oil	Arsenic	.21	tons
	Chromium	.000035	% Waste	oil	Cadimium	.11	tons
					Chromium	.41	tons

EMISSION FACTOR SOURCE: A.D.Little, Waste Oil Combustion at a Bituminous Concrete Batching Plant, August 1984.

COUNTY	POPULATION	USAGE gal	LEAD tons	CHLORINE tons	ARSENIC	CADMIUM tons	CHROMIUM tons
ANDROSCOGGIN	100900	263515.2	1.05	4.21	.Ø18	.010	.Ø36
AROOSTOOK	90400	236092.9	.94	3.77	.016	.009	.Ø33
CUMBERLAND	220100	574823.7	2.29	9.19	.041	.022	.080
FRANKLIN	28500	74431.9	. 29	1.19	.005	.002	.010
HANCOCK	42700	111517.3	.44	1.78	.008	.004	.015
KENNEBDC	112850	294724.4	1.17	4.71	.021	.011	.041
KNOX	, 34050	88925.6	.35	1.42	.006	.003	.012
LINCOLN	26150	68294.5	.27	1.09	.004	.002	.009
OXFORD	48800	127448.4	.50	2.03	.009	.005	.017
PENOBSCOT	141150	368634.1	1.47	5.89	.026	.014	.051
PISQUATAQUIS	17850	46617.9	.18	.74	.003	.001	.006
SAGADAHOC	28050	73256.7	.29	1.17	.005	.002	.010
SOMERSET	47050	122878.0	.49	1.96	.008	.004	.017
WALDO	28500	74431.9	.29	1.19	.005	.002	.010
WASHINGTON	33550	87620.7	.35	1.40	.006	.003	.Ø12
YORK	148100	386785.0	1.54	6.18	.027	.015	.054
TOTALS	1148700	3000000	11.99	47.99	.215	.119	.419

WASTE OIL

AREA SOURCE EMISSIONS DATA

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Automotive lead emissions

DATA COLLECTION:

Lead emissions by country from Maine State Implementation Plan data Table 2.5.1 (3) "Area source automotive lead emissions by counties."

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CALCULATIONS:

DATA ASSUMPTIONS:

EMISSION FACTOR(S):

TOTAL EMISSIONS:

Lead emissions

89.4 tons

EMISSION FACTOR SOURCE:

AUTOMOTIVE LEAD EMISSIONS

AREA SOURCE EMISSIONS DATA

COUNTY	POPULATION	LEAD EMISSIONS (tons/yr)
ANDROSCOGGIN	· 100900	6.5
AROOSTOOK	90400	6.4
CUMBERLAND	220100	18.3
FRANKLIN	28500	2.1
HANCOCK	42700	4.4
KENNEBEC	1 1285Ø	8.6
KNOX	34050	2.2
LINCOLN	26150	2.4
OXFORD	48800	3.6
PENOBSCOT	141150	10.6
PISQUATAQUIS	178 5 Ø	1.3
SAGADAHOC	28050	2.3
SOMERSET	47050	4.3
WALDO	285ØØ	2.3
WASHINGTON	3355ø	2.8
YORK	148100	11.3

TOTALS 1148700

89.4

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AREA SOURCE EMISSIONS DATA

CATEGORY:

Gasoline Service Station Operations

DATA COLLECTION:

Gasoline usage obtained from Maine gasoline sales information, provided by the Maine Office of Energy Resources. Total gas sold was apportioned to the county level by % vehicle miles traveled (VMT). VMT data was obtained from the Maine Dept. of Transportation.

DATA ASSUMPTIONS:

1. VMT is proportional to amount of gasoline used.

2. 1982 VMT data is representative of 1983 conditions.

3. Gasoline applies only to leaded, unleaded, and regular gasoline.

4. All service stations operations utilize no emission controls.

5. % benzene and toluene emissions from fuel tanks similar to all other gasoline station operations.

EMISSION FACTOR(S):

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Splash filling Underground tank	11.5 lb/thous gal
breathing + emptying	1.0 lb/thous gal
Refueling loss	9.0 lb/thous gal
Spillage	.7 lb/thous gal
Benzene	.004 x VOC(tons)
Toluene	.004 x VOC(tons)

EMISSION FACTOR SOURCE:

AP-42, (Table 4.4-4): Hydrocarbons and Vapor Phase Organic Pollutants Table 2-6, page 13, "Benzene and toluene emissions from gas tanks," Jackson and Everett.

CALCULATIONS:

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TOTAL EMISSIONS:

VOC	5955.63 tons
Benzene	23.82 tons
Toluene	23.82 tons
Underground tanks Splash filling*	3109.80 tons
Breathing + emptying*	270.41 tons
Refueling loss*	2433.76 tons
Spillage*	189.29 tons

*All pollutants

AREA SOURCE EMISSIONS

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GASOLINE SERVICE STATION OPERATIONS

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		POLLUTANTS EMITTED	SPLASH FILLING	UND TANKS BREATHING +EMPYTING tons	REFUELING LOSS tons		TOTAL VOC tons	TOTAL BENZENE tons	
ANDROSCOGGIN	41667.0	VOC	239.58	20.83	187.50	14.58	462.50		
		Benzene	.95	.08	.75			1.85	
		Toluene	.95	. Ø8	.75	.05			1.85
AROOSTOOK	41096.2	VOC	236.30	20.54	184.93	14.38	456.16		
		Benzene	.94	.08	.73	.05		1.82	
		Toluene	.94	.08	.73				1.82
CUMBERLAND	114727.0	VOC			516.27		1273.46		
		Benzene	2.63					5.09	
		Toluene		.22					5.09
FRANKLIN	14840.0		85.33				164.72		
		Benzene		.02				.65	
		Toluene		.02					•65
HANCOCK	28539.0	VOC					316.78		
		Benzene	•65					1.26	
		Toluene	•65		.51				1.26
KENNEBEC	55365.8	VOC		27.68	249.14		614.56		
		Benzene	1.27	_	99	.07		2.45	
		Toluene	1.27						2.45
KNOX	14269.5		82.04	7.13	64.21	4.99	158.39		
		Benzene	.32		•25			.63	
		Toluene	.32			.01			.63
LINCOLN	15411.0		88.61				171.06		
		Benzene	•35					.68	
		Toluene		.03					•68
OXFORD	23402.0		134.56				259.76		
		Benzene	•53					1.03	
		Toluene	•53		.42				1.03
PENOBSCOT	34246.9			17.12		11.98	380.14		
		Benzene	.78					1.52	
		Toluene	.78	.06	.61	.04			1.52

TOTALS ON NEXT PAGE

AREA SOURCE EMISSIONS (CONT.)

GASOLINE SERVICE STATION OPERATIONS

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1999 - Albert A.

	USAGE THRU-PUT THOUS GALS	POLLUTANTS EMITTED	UNDERGRO SPLASH FILLING tons	UND TANKS BREATHING +EMPYTING tons	REFUELING LOSS tons	SPILLAGE	TOTAL VOC tons	TOTAL BENZENE tons	TOTAL TOLUENE tons
PISQUATAQUIS	8561.7	VOC	49.22	4.28	38.52	2.99	95.03		
		Benzene	.19	.01	.15	.01		.38	
		Toluene	.19	.01	.15	.01			38
SAGADAHOC	13698.7	VOC	78.76			4.79	152.05	•	
		Benzene	.31			.01		.60	
		Toluene	.31			.01			.60
SOMERSET	25695.1	VOC	147.74			8.99	285.21		
		Benzene	•59			.03		1.14	
		Toluene	.59			.03			1.14
WALDO	15411.1	VOC	88.61			5.39	171.06		
		Benzene	.35			.02		.68	
		Toluene	.35			.02			.68
WASHINGTON	18265.0		105.02			6.39	202.74		
		Benzene	.42			.02		.81	
		Toluene	.42			.02			.81
YORK	71347.6		410.24			24.97	791.95		
1014	/101/00	Benzene	1.64			.09	,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	3.16	
		Toluene	1.64			.09		5010	3.16
TOTALS	536543.6		3109.8	270.4	2433.7	189.2	5955.63	23.82	23.82

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	BENEZENE	TOLENE	VOC
Техасо	.Ø4 Tons	.04 Tons	10.6 Tons
Gulf	.15	.15	38.7
Mobil	•23	.23	57.5
Chevron	.23	.23	57.1
Koch	.004	.004	1.0
Getty	• Ø9	.Ø9	22.4
B.P.	. 29	.29	71.3
Exxon	.09	.09	21.9
Mobil (Bangor)	. 6Ø	. 6Ø	150.3
Texaco (Bangor)	•65	.65	161.9
Mobil (Hall)	.002	.002	.38
TOTAL	2.376	2.376	630.7

AREA SOURCE EMISSIONS FROM GASOLINE TERMINAL

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** APPENDIX D **

APPENDIX D

COUNTY CODES

CODE COUNTY Øl ANDROSCOGGIN ØЗ AROOSTOCK Ø5 CUMBERLAND Ø7 FRANKLIN Ø9 HANCOCK 11 KENNEBEC 13 KNOX 15 LINCOLN 17 OXFORD 19 PENOBSCOT 21 PISCATAQUIS SAGADAHOC 23 25 SOMERSET 27 WALDO 29 WASHINGTON 31 YORK

PAGE NO. 1

MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION HAZARDOUS AIR POLLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

	POLLUTANT	AMOUNT USED	EMISSIONS
COUNTY	IDENT.	TONS/YEAR	L8S.7YEAR
01	ACETON	645.1	206,310.0
	ACRYLO	57.5	0.0
	BACETA	6.3	4,930.0
	BUTANO	2.2	0.0
	CHLORI	75.1	100.0
	COPPER	125.0	100.0
	CRESOL	2.5	0.0
	DBENZE	2.5	0.0
	DMETHA	440.5	56,000.0
	EACETA	12.5	22,231.0
	EOXIDE	3.6	966.0
	EPICHL	31.0	12.0
······································	ETHANO	21.4	0.0
	FALCOH	2.0	0.0
	FORMAL	3,375.0	25,003.0
	FORMIC	7.5	1.0
	HCHLOR	89.3	512.0
	HSULFI	0.0	0.0
	IMINOD	8.5	Ů.O
	LEADPB	2.0	0.0
	MALEIC	143.0	
	MCELLO	10.5	0.0
	MECHLO	218.6	190,145.0
	MEKETO	265.0	103,174.8
	MELAMI	2,321.0	0.0
	NITRIC	52.1	2,500,0
	OXALIC	6.0	0.0
	PHENOL	3,524.0	3,247.0
	STYREN	86.0	21.0
	TDIOXI	1.6	0.0
······································	TETHAI	878.9	655,298.0
	TETHYL	218.7	19.0
,,,	TITANI	48.7	0.0
	TOLUEN	159.3	75,301.0
	TRICHL	8.8	0.0
	VCHLOR	501.4	0.0
	XYLENE	284.9	22,378.8
03	ACETIC	77.5	1,402.6
	COIOXI	6.0	0.0
	CHLORI	59.1	22,800.0
	EACETA	1.3	2,600.0
	EPOXYP	401.0	124,600.0
		36.5	600.0
	MEKETO	4.3	3,200.0
	PCBIPH	42.8	0.0
	TETHA1	7.3	. 14,600.0
······································	TETHYL	9.5	14,860.0
· ·	XYLENE	58.9	117,800.0
05	ACETON	. 144.9	125,311.0
	ANTIMU	6.7	0.0

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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 2 HAZARDOUS AIR POLLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

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2011154	POLLUTANT	AMOUNT USED	EMISSIONS LBS./YEAR
SOUNTY	IDENT.	TONS/YEAR	LBS./TEAR
·····	ASBEST	5.6	0.0
	BACETA	21.9	34,203.0
	BARIUM	41.7	374.0
	BUTANO	84.8	116,340.0
	CDIOXI	370.0	4,000.0
<u></u>	CHLORI	17,810.0	85,000.0
	CHROMI	1.2	0.0
	COPPER	2.7	6.6
	DBENZE	2.5	0.0
	DETHAN	68.1	32,400.0
	EACETA	3.1	6,208.0
	EGLYCO	32.0	7,461.0
	EOXIDE	0.5	569.0
	FORMAL	20.6	41,000.0
	FORMIC	166.9	4,892.0
	HCHLOR	82.5	6,160.0
	HSULFI	14.8	22,000.0
	LEADPB	20.2	550.0
	MANGAN	6.2	9,320.0
·	MCELLO	52.9	7,708.0
	MCHLOR	1.8	3,600.0
	MECHLO MEKETO	63.9 44.7	41,081.0 13,167.9
	MELAMI	50.0	0.0
	MISOBU	1.0	0.0
	MMERCA	34.9	30,700.0
	NITRIC	107.8	1,620.0
	OXALIC	1.3	0.0
	PCBIPH	40.5	0.0
	PHENOL	2.6	0.7
	STYREN	47.4	6,775.0
	TETHA1	194.1	168,290.0
	TETHYL	16.8	9,060.0
	TFURAN	0.2	400.0
	TITANI	2,669.2	1,800.0
	TOLUEN	2,656.9	606,635.0
	TRICHL	104.4	53,977.0
	TURPEN	186.0	12,000.0
	XYLENE	171.9	36,854.7
	ZINCZN	92.0	946.0
07	ACETON	24.6	42,363.0
·	BACETA	1.2	2,400.0
		3,364.0	34,000.0
	CHLORI	19,406.5	300,000.0
•	EACETA FORMIC	1.8	3,600.0
	HSULFI	530.0	100,000.0
	MECHLO	9.8	16,265.0
	MEKETO	40.7	76,958.0
	MERCA	374.0	180,000.0
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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION HAZARDOUS AIR POLLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

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	POLLUTANT	AMOUNT USED TONS/YEAR	EMISSIUNS LBS./YEAR
	IDENT.	TUNSZTEAR	LOGATIEAR
	NITRIC	4.5	9,000.0
	TITANI	1,000.0	0.0
	TOLUEN	23.0	43,676.0
	TRICHL	1.2	2,420.0
	TURPEN	9,000.0	30,000.0
09	ACETON	11.7	4,000.0
	CHLORI	53.0	0.0
	ETHANO	3.7	734.0
	MECHLO	1.0	2,000.0
	MEKETO	1.0	40.0
	STYREN	35.0	3,500.0
	TITANI	243.0	0.0
11	ACETON	13.5	15,000.0
	ASBEST	28.7	0.0
	BACETA	0.1	200.0
	CDIOXI	10.0	50.0
	CHFORM	0.0	0.0
	CHLORI	2,831.1	66,163.5
	CHROMI	1.2	48.0
	EACETA	0.8	1,600.0
	EGLYCO	2.5	4,964.0
	ETHANO	24.6	340.0
	FORMAL	92.7	4,118.0
	FORMIC	27.9	27,744.0
	HCHLOR	46.5	26,464.0
	MECHLO	3.0	0.0
	MEKETO	6.9	13,850.0
	OXALIC	14.0	560.0 252,675.0
		90.0	168,540.0
13	ACETON	12.4	24,750.0
	FORMIC	41.2	1,648.0
	HCHLOR	3.5	140.0
	MEKETO	0.8	1,608.0
	NITRIC	3.5	350.0
	TETHA1	1.8	36.0
	TETHYL	4.0	7,000.0
	TOLUEN	24.0	48,000.0
	XYLENE	5.5	8,316.0
15	ACETON	0.5	150.0
	HYDRAZ	3.5	140.0
•	MECHLO	1.5	360.0
	MMETHA	11.6	650.0
	PCBIPH	55.0	0.0
······································	STYREN	10.3	205.0
	TETHAL	. 3.9	4,257.0
17	ACETON	5.4	6,733.0
	BACETA	32.0	64,054.0
	BARIUM	3.7	0.0
	BUTANO	5.0	9,947.0
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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION HAZARDOUS AIR POLLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

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	POLLUTANT	AMOUNT USED	EMISSIONS
YTVUCC	IDENT.	TONS/YEAR	LBS./YEAR
	COLONI	7 190 0	67,895.0
	CDIOXI CHFORM	3,280.0	0.0
			250,000.0
·····	CHLORI DMETHA	14,676.0 45.0	90,000.0
		43.0	9,354.0
	EACETA	9.6	19,094.0
	EGLYCO	41.2	1,648.0
	FORMIC	1.9	72.0
	RUHLOR	220.0	41,000.0
	HSULFI		0.0
	LEADPB	0.6	42,092.0
	MCELLO	21.0	
	MECHLO	44.0	4,400.0
	MEKETO	15.4	30.578.0
	MMERCA	150.0	70,000.0
	VALINE	4.8	9,555.0
	NITRIC	23.1	920.0
	VPHENO	3.0	120.0
	OXALIC	8.1	0.0
	TETHA1	28.0	56,000.0
	TITANI	1,217.0	0.515
	TOLUEN	28.9	57,602.0
	TURPEN	4,000.0	12,000.0
	XYLENE	115.6	231,232.0
19	ACETON	140.7	224,700.8
	BACETA	1.2	0.0
	BARIUM	2.5	1,782.0
	3IPHEN	0.05	800.0
	COIOXI	2,800.0	87,900.0
	CHFORM	0.0	0.0
	CHLORI	76,502.0	300,200.0
	CTETRA	2.5	0.0
	CYANID	1.3	2,600.0
<u></u>	DJENZE	60.0	2,400.0
	EACETA	1.2	2,064.0
	ETHANO	6.4	0.0
· · · · · · · · · · · · · · · · · · ·	FORMIC	83.9	6,952.0
	HCHLOR	9,641.6	1,282.6
	HSULFI	377,394.8	45,110.0
	MECHLO	16.6	25,772.6
	MEKETO	87.1	143,730.6
	MERCUR	8.9	16.4
<u></u>	MMERCA	377,320.2	49,440.0
	VALINE	0.4	800.0
	NITRIC	35.3	0.005
	STYREN	25.0	2,500.0
	TETHA1	88.7	101,256.3
	TETHYL	3.1	6,200.0
	TITANI	726.0	1,780.0
	TOLUEN	734.4	1,433,143.2
	TRICHL	7.8	3,000.0

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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 5 HAZARDOUS AIR POLLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

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· · ·	POLLUTANT	AMOUNT USED	EMISSIONS
COUNTY	IDENT.	TONS/YEAR	LBS./YEAR
	TURPEN	5,000.0	17,000.0
	XYLENE	112.3	209,900.0
21	ACETON	28.0	56,000.0
	JACETA	3.4	4,635.5
<u></u>	BIPHEN	3.7	3.710.0
	BUTANO	0.1	76.0
·	EACETA	0.1	224.0
	EGLYCO	0.1	48.0
	ETHANO	0.1	244.0
	MCHLOR	3.4	6,800.0
<u></u>	MEKETO	13.8	27,280.0
	MISOBU	1.8	3,609.6
	NALINE	0.2	444.0
	TETHAL	7.6	15,262.0
	TOLUEN	29.8	59,662.8
	XYLENE	15.7	21,378.1
23	ACETON	5.0	10,000.0
	BACETA	2.0	0.085
	BUTANO	9.0	11,020.0
	EACETA	10.5	17,056.0
	EBENZE	0.5	80.0
	EGLYCO	15.9	2,451.0
	LEADPB	160.0	0.0
	MCELLO	- 4 . 7	5,608.0
	MEKETO	13.0	26,000.0
	MISOBU	5.5	2,816.0
	TETHAI	3.5	5,960.0
	TOLUEN	6.2	11,780.0
	XYLENE	57.3	71,688.0
25	ACETON	98.9	64,806.0
	BACETA	70.3	140,600.0
	ЗРНТНА	1.5	3,000.0
· · · · · · · · · · · · · · · · · · ·	BUTANO	2.7	5,400.0
	CDIOXI	4,520.0	16,800.0
	CHLORI	13,141.0	16,400.0
	EACETA	0.8	1,600.0
•	EGLYCU	12.8	25,600.0
	ETHANO	1.0	2,000.0
	FORMAL	1.4	0.0
	FORMIC	58.1	2,320.0
	HCHLOR	84.7	3,300.0
	HSULFI	93.9	4,802.0
	LEADPB	2.0	80.0
	MCELLO	51.6	103,200.0
	MCHLOR	3.2	6,400.0
· · ·	MECHLO	1.4	2,713.0
· · · · · · · · · · · · · · · · · · ·	MEKETO	14.8	29,600.0
	MISOBU	1.6	2,739.0
ê	MMERCA	236.0	1,800.0
	NALINE	0.5	1,000.0
· · · · · · · · · · · · · · · · · · ·			
	·	58-	······································

MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 6 HAZARDOUS AIR PULLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

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DENT. PHENO XALIC HENOL ETHA1 ITANI OLUEN URPEN YLENE SCETON ACETA URFUR IEKETO NISOBU OLUEN YLENE ACETA DIOXI HLORI ORMAL ICHLOR ISULFI MERCA ITRIC PHENOL ITANI	$\begin{array}{r} \hline \text{TONS/YEAR} \\ \hline 2.7 \\ 4.8 \\ \hline 2.8 \\ 16.8 \\ \hline 2.617.8 \\ \hline 36.8 \\ 1.262.0 \\ \hline 71.0 \\ \hline 3.5 \\ 1.7 \\ \hline 7.3 \\ \hline 2.6 \\ \hline 6.1 \\ 14.2 \\ \hline 11.2 \\ \hline 22.5 \\ \hline 2.589.0 \\ \hline 13.900.0 \\ \hline 4.6 \\ \hline 17.4 \\ \hline 300.0 \\ \hline 200.0 \\ \hline 28.8 \\ \hline 1.2 \\ \hline 1.0 \\ \hline 650.0 \\ \end{array}$	LBS./YEAR 108.0 192.0 112.0 31,109.0 1,448.0 71,423.0 4,130.0 1,423.0 4,130.0 1,423.0 1,423.0 4,130.0 2,800.0 1,360.0 2,080.0 4,880.0 11,360.0 8,960.0 900.0 85,142.0 9,280.0 700.0 57,000.0 97,000.0 2,450.0 2,000.0 0,000
XALIC HENOL ETHA1 ITANI OLUEN URPEN YLENE CETON ACETA URFUR IEKETO IISOBU OLUEN YLENE ACETA DIOXI CHLOR ISOBU OLUEN YLENE ACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL YETHA1	$ \begin{array}{r} 4.8\\ 2.8\\ 16.8\\ 2.617.8\\ 36.8\\ 1.262.0\\ 71.0\\ 3.5\\ 1.7\\ 7.3\\ 2.6\\ 6.1\\ 14.2\\ 11.2\\ 22.5\\ 2.589.0\\ 13.900.0\\ 4.6\\ 17.4\\ 300.0\\ 200.0\\ 28.8\\ 1.2\\ 1.0\end{array} $	192.0 112.0 112.0 112.0 112.0 1.448.0 1.448.0 71.423.0 4.130.0 142.000.0 2.800.0 1.360.0 5.840.0 2.080.0 4.880.0 11.360.0 8.960.0 900.0 86.142.0 287.139.0 9.280.0 700.0 57.000.0 1.200.0 2.450.0 2.000.0 0
XALIC HENOL ETHA1 ITANI OLUEN URPEN YLENE CETON ACETA URFUR IEKETO IISOBU OLUEN YLENE ACETA DIOXI CHLOR ISOBU OLUEN YLENE ACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL YETHA1	$ \begin{array}{r} 4.8\\ 2.8\\ 16.8\\ 2.617.8\\ 36.8\\ 1.262.0\\ 71.0\\ 3.5\\ 1.7\\ 7.3\\ 2.6\\ 6.1\\ 14.2\\ 11.2\\ 22.5\\ 2.589.0\\ 13.900.0\\ 4.6\\ 17.4\\ 300.0\\ 200.0\\ 28.8\\ 1.2\\ 1.0\end{array} $	192.0 112.0 112.0 112.0 112.0 1.448.0 1.448.0 71.423.0 4.130.0 142.000.0 2.800.0 1.360.0 5.840.0 2.080.0 4.880.0 11.360.0 8.960.0 900.0 86.142.0 287.139.0 9.280.0 700.0 57.000.0 1.200.0 2.450.0 2.000.0 0
HENOL ETHA1 ITANI OLUEN URPEN VIENE CETON ACETA URFUR IEKETO IISOBU OLUEN VYLENE ACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL IETHA1	$\begin{array}{r} 2.8 \\ 16.8 \\ 2.617.8 \\ 36.8 \\ 1.262.0 \\ 71.0 \\ 3.5 \\ 1.7 \\ 7.3 \\ 2.6 \\ 6.1 \\ 14.2 \\ 11.2 \\ 22.5 \\ 2.5$	$ \begin{array}{r} 112.0\\ 31,109.0\\ 1,448.0\\ 71,423.0\\ 4,130.0\\ 142.000.0\\ 2,800.0\\ 142.000.0\\ 2,800.0\\ 1,360.0\\ 5,840.0\\ 2,080.0\\ 4,880.0\\ 11,360.0\\ 8,960.0\\ 900.0\\ 85,142.0\\ 287.139.0\\ 9,280.0\\ 700.0\\ 57,000.0\\ 97,000.0\\ 2,450.0\\ 2,000.0\\ 0,000.0$
ETHA1 ITANI OLUEN URPEN YLENE CETON ACETA URFUR IEKETO IISOBU OLUEN YLENE BACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL YETHA1	$ \begin{array}{r} 16.8 \\ 2.617.8 \\ 36.8 \\ 1.262.0 \\ 71.0 \\ 3.5 \\ 1.7 \\ 7.3 \\ 2.6 \\ 6.1 \\ 14.2 \\ 11.2 \\ 22.5 \\ 2.5 \\ $	$ \begin{array}{c} 1,448.0\\ 71,423.0\\ 4,130.0\\ 4,130.0\\ 142,000.0\\ 2,800.0\\ 1,360.0\\ 5,840.0\\ 2,080.0\\ 4,880.0\\ 11,360.0\\ 8,960.0\\ 900.0\\ 86.142.0\\ 287.139.0\\ 9,280.0\\ 700.0\\ 57,000.0\\ 97,000.0\\ 2,450.0\\ 2,000.0\\ \end{array} $
ITANI OLUEN URPEN VLENE CETON ACETA URFUR IEKETO IISOBU OLUEN VLENE DACETA DIOXI HLORI ORMAL ICHLOR ISULFI MERCA VITRIC PHENOL IETHAI	2.617.8 36.8 $1.262.0$ 71.0 3.5 1.7 7.3 2.6 6.1 14.2 11.2 22.5 $2.589.0$ $13.900.0$ 4.6 17.4 300.0 200.0 28.8 1.2 1.0	71,423.0 4,130.0 142.000.0 2,800.0 1,360.0 5,840.0 2,080.0 4,880.0 11,360.0 8,960.0 900.0 86.142.0 287.139.0 9,280.0 700.0 57,000.0 97,000.0 2,450.0 2,000.0
URPEN YLENE CETON ACETA URFUR IEKETO IISOBU OLUEN YLENE ACETA DIOXI HLORI ORMAL ICHLOR ISULFI MERCA IITRIC PHENOL IETHAI	$ \begin{array}{r} 1,262.0 \\ 71.0 \\ 3.5 \\ 1.7 \\ 7.3 \\ 2.6 \\ 6.1 \\ 14.2 \\ 11.2 \\ 22.5 \\ 2,589.0 \\ 13,900.0 \\ 4.6 \\ 17.4 \\ 300.0 \\ 200.0 \\ 28.8 \\ 1.2 \\ 1.0 \\ \end{array} $	$\begin{array}{c} 4,130.0\\ 142,000.0\\ 2,800.0\\ 2,800.0\\ 1,360.0\\ 5,840.0\\ 2,080.0\\ 2,080.0\\ 4,880.0\\ 11,360.0\\ 8,960.0\\ 900.0\\ 85,142.0\\ 287.139.0\\ 9,280.0\\ 700.0\\ 57,000.0\\ 97,000.0\\ 1,200.0\\ 2,450.0\\ 2,000.0\\ \end{array}$
YLENE CETON ACETA URFUR IEKETO IISOBU OLUEN YLENE ACETA DIOXI CHLORI ISULFI MERCA IITRIC PHENOL IETHAI	71.0 3.5 1.7 7.3 2.6 6.1 14.2 11.2 22.5 $2,589.0$ $13,900.0$ 4.6 17.4 300.0 200.0 28.8 1.2 1.0	$ \begin{array}{c} 142,000.0\\ 2,800.0\\ 1,360.0\\ 5,840.0\\ 2,080.0\\ 4,880.0\\ 11,360.0\\ 8,960.0\\ 900.0\\ 8,960.0\\ 900.0\\ 85,142.0\\ 287.139.0\\ 9,280.0\\ 700.0\\ 57,000.0\\ 97,000.0\\ 2,450.0\\ 2,000.0\\ \end{array} $
ACETON ACETA URFUR IEKETO IISOBU OLUEN YLENE ACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL IETHAI	3.5 1.7 7.3 2.6 6.1 14.2 11.2 22.5 $2,589.0$ $13,900.0$ 4.6 17.4 300.0 200.0 28.8 1.2 1.0	2,800.0 $1,360.0$ $5,840.0$ $2,080.0$ $4,880.0$ $11,360.0$ $8,960.0$ 900.0 $85,142.0$ $287,139.0$ $9,280.0$ 700.0 $57,000.0$ $1,200.0$ $2,450.0$ $2,000.0$
ACETA URFUR IEKETO IISOBU OLUEN CYLENE DACETA DIOXI CHLOR ISULFI IMERCA IITRIC PHENOL IETHAI	$ \begin{array}{r} 1.7 \\ 7.3 \\ 2.6 \\ 6.1 \\ 14.2 \\ 11.2 \\ 22.5 \\ 2.5 \\ $	$ \begin{array}{c} 1,360.0\\ 5,840.0\\ 2,080.0\\ 4,880.0\\ 11,360.0\\ 8,960.0\\ 900.0\\ 86,142.0\\ 287,139.0\\ 9,280.0\\ 700.0\\ 57,060.0\\ 97,000.0\\ 1,200.0\\ 2,450.0\\ 2,000.0\\ \end{array} $
URFUR IEKETO IISOBU OLUEN KYLENE BACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL IETHAI	7.3 2.6 6.1 14.2 11.2 22.5 $2.589.0$ $13,900.0$ 4.6 17.4 300.0 200.0 28.8 1.2 1.0	5,840.0 2,080.0 4,880.0 11,360.0 8,960.0 900.0 86,142.0 287,139.0 9,280.0 700.0 57,000.0 97,000.0 1,200.0 2,450.0
IEKETO IISOBU OLUEN KYLENE BACETA DIOXI CHLOR ISULFI MERCA IITRIC PHENOL IETHAI	2.6 6.1 14.2 11.2 22.5 $2,589.0$ $13,900.0$ 4.6 17.4 300.0 200.0 28.8 1.2 1.0	2,080.0 $4,880.0$ $11,360.0$ $8,960.0$ 900.0 $85,142.0$ $287,139.0$ $9,280.0$ 700.0 $57,000.0$ $97,000.0$ $1,200.0$ $2,450.0$ $2,000.0$
ISOBU OLUEN VLENE DACETA DIOXI CHLORI ISULFI MERCA VITRIC PHENOL IETHAI	$ \begin{array}{r} 6.1\\ 14.2\\ 11.2\\ 22.5\\ 2,589.0\\ 13,900.0\\ 4.6\\ 17.4\\ 300.0\\ 200.0\\ 200.0\\ 28.8\\ 1.2\\ 1.0\end{array} $	4,880.0 11,360.0 8,960.0 900.0 86,142.0 287,139.0 9,280.0 700.0 57,000.0 97,000.0 2,450.0 2,000.0
OLUEN YLENE DACETA DIOXI HLORI ORMAL ICHLOR ISULFI MERCA VITRIC PHENOL IETHAI	14.2 11.2 22.5 2,589.0 13,900.0 4.6 17.4 300.0 200.0 28.8 1.2 1.0	11,360.0 8,960.0 900.0 85,142.0 287,139.0 9,280.0 700.0 57,000.0 97,000.0 1,200.0 2,450.0 2,000.0
ACETA DACETA DIOXI CHLORI CORMAL ICHLOR ISULFI IMERCA NITRIC PHENOL IETHAI	11.2 22.5 2,589.0 13,900.0 4.6 17.4 300.0 200.0 28.8 1.2 1.0	8,960.0 900.0 85,142.0 287,139.0 9,280.0 700.0 57,000.0 97,000.0 1,200.0 2,450.0 2,000.0
DACETA DIOXI CHLORI ICHLOR ISULFI IMERCA VITRIC PHENOL IETHAI	22.5 2,589.0 13,900.0 4.6 17.4 300.0 200.0 28.8 1.2 1.0	900.0 86.142.0 287.139.0 9.280.0 700.0 57.000.0 97.000.0 1.200.0 2.450.0 2.000.0
EDIOXI CHLORI FORMAL ICHLOR ISULFI MERCA VITRIC PHENOL FETHAI	2,589.0 13,900.0 4.6 17.4 300.0 200.0 28.8 1.2 1.0	86.142.0 287.139.0 9,280.0 700.0 57.000.0 97,000.0 1,200.0 2,450.0 2,000.0
CHLORI ORMAL ICHLOR ISULFI MERCA IITRIC PHENOL IETHAI	$ \begin{array}{r} 13,900.0\\ 4.6\\ 17.4\\ 300.0\\ 200.0\\ 28.8\\ 1.2\\ 1.0\\ \end{array} $	287,139.0 9,280.0 700.0 57,000.0 97,000.0 1,200.0 2,450.0 2,000.0
ORMAL ICHLOR ISULFI MERCA IITRIC PHENOL IETHAI	4.6 17.4 300.0 200.0 28.8 1.2 1.0	9,280.0 700.0 57,000.0 97,000.0 1,200.0 2,450.0 2,000.0
ICHLOR ISULFI IMERCA IITRIC PHENOL IETHAI	17.4 300.0 200.0 28.8 1.2 1.0	700.0 57,000.0 97,000.0 1,200.0 2,450.0 2,000.0
ISULFI IMERCA IITRIC PHENOL IETHAI	300.0 200.0 28.8 1.2 1.0	57,000.0 97,000.0 1,200.0 2,450.0 2,000.0
MERCA NITRIC PHENOL TETHA1	200.0 28.8 1.2 1.0	97,000.0 1,200.0 2,450.0 2,000.0
HENOL MENOL METHA1	28.8 1.2 1.0	1,200.0 2,450.0 2,000.0
PHENOL TETHA1	1.2 1.0	2,450.0 2,000.0
TETHA1	1.0	2,000.0
	0.010	V • V
TURPEN	5,000.0	2,000.0
CETON	12.4	21,597.0
NTIMO	4.0	0.0
BACETA	0.1	Ú.C
3PHTHA	3.6	7,190.0
BUTANO	3.5	6,707.0
CHLORI	6.6	9,200.0
CHROMI	1.7	136.0
CYANID	1.6	355.0
		5,424.0
		0.0
		217,500.0
		4,656.0
······································		<u> </u>
		600.0
		473,640.0
		5,832.0
		41,344.0
and a second		183,900.0
		25,660.0
		116.0
		189,796.0
	SULFA ACETA GLYCO ORMIC URFUR CHLOR YDRAZ CELLO ECHLO ECHLO EKETO METHA ITRIC XALIC CBIPH ETHA1	SULFA 135.6 ACETA 0.1 GLYCO 108.8 ORMIC 116.4 URFUR 3.2 CHLOR 13.5 YDRAZ 0.3 CELLO 236.9 ECHLO 4.2 EKETU 32.3 METHA 25.896.0 ITRIC 130.9 XALIC 2.9 CBIPH 6.7

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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 7 HAZARDOUS AIR POLLUTANT SYSTEM PROCESS EMISSIONS BY COUNTY (HAP114)

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	POLLUTANT	AMOUNT USED	EMISSIONS
YTVUGC	IDENT.	TONS/YEAR	LBS./YEAR
•	TETHYL	84.0	107,311.0
	TITANI	2.0	0.0
	TOLUEN	27.3	40,648.0
	TRICHL XYLENE	117.5 28.1	51,468.0 52,226.0
	ATEENE		
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** APPENDIX E **

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APPENDIX E

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STANDARD INDUSTRIAL CLASSIFICATION

DIVISION	MAJOR GROUP	INDUSTRY
MANUFACTURING	2Ø 22 24 25 26 28 29	FOOD AND KINDRED PRODUCTS TEXTILE MILL PRODUCTS LUMBER AND WOOD PRODUCTS WOOD FURNITURE AND FIXTURES PAPER AND ALLIED PRODUCTS CHEMICALS AND ALLIED PRODUCTS PETROLEUM REFINING AND RELATED INDUSTRIES
	3Ø 31 32 33 34 35	RUBBER AND MISCELLANEOUS PLASTICS LEATHER AND LEATHER PRODUCTS STONE, CLAY, GLASS, AND CONCRETE PRODUCTS PRIMARY METAL PRODUCTS FABRICATED METAL PRODUCTS MACHINERY
	36 37 38 39	ELECTRICAL AND ELECTRONIC MACHINERY TRANSPORTATION EQUIPMENT MEASURING, ANALYZING, AND CONTROLLING INSTRUMENTS; PHOTOGRAPHIC, MEDICAL, AND OPTICAL GOODS; WATCHES AND CLOCKS MISCELLANEOUS MANUFACTURING INDUSTRIES
TRANSPORTATION, COMMUNICATIONS, ELECTRIC, GAS, AND SANITARY SERVICES		
	40 42 49	RAILROAD TRANSPORTATION MOTOR FREIGHT TRANSPORTATION AND WAREHOUSING ELECTRIC, GAS, AND SANITARY SERVICES
WHOLESALE TRADE	51	WHOLESALE TRADE- NONDURABLE GOODS
SERVICES	72 73 75 8Ø	PERSONAL SERVICES BUSINESS SERVICES AUTOMOTIVE REPAIR, SERVICE, AND GARAGES HEATLTH SERVICES
PUBLIC ADMINISTRATION	97	NATIONAL SECURITY AND INTERNATIONAL AFFAIRS

MAR 11, 1985DEPARTMENT OF ENVIRONMENTAL PROTECTIONPAGE NO. 1_____HAZARDOUS AIR POLLUTANT SYSTEMSTATENIDE SUMMARY OF PROCESS EMISSIONS BY INDUSTRIAL CLASSIFICATION (HAP112)

	3 () 1.T 1.T	AMOUNT USED	EMISSIONS
	POLLUTANT	TONS/YEAR	LUS./YEAR
20	ACETIC	77.5	1,402.6
	BACETA	22.5	900.0
	CDIOXI	6.0	0.0
	CHLORI	19.9	2,400.0
······	EPOXYP	401.0	124,600.0
	HCHLOR	· 36.5	600.0
22	ANTIMO	4.0	0.0
	BIPHEN	23.7	4,510.0
	CHLORI	1.3	0.0
	CHROMI	1.2	48.0
	DBENZE	60.0	2,400.0
	DSULFA	135.6	5,424.0
. <u></u>	EGLYCO	2.1	4,140.0
	FORMIC	72.4	22,296.0
	MEKETO	15.4	13,664.0
	OXALIC	22.1	560.0
	TETHAI	143.8	287,673.3
	TOLJEN	10.8	11,521.0
	XYLENE	174.6	337,390.0
2.4	BACETA	5.2	6,366.0
	BUTANO	0.7	1,238.0
	DMETHA	45.0	90,000.0
	EACETA	1.8	3,695.0
	EGLYCO	0.1	48.0
	ETHANO	0.1	244.0
	FORMAL	4.6	9,280.0
	MECHLO	44.0	4,400.0
	MEKETO	5.8	11,466.0
	NALINE	5.0	9,999.0
	PHENOL	1.2	2,450.0
	TETHAL	23.5	47,000.0
	TOLUEN	30.9	61,802.0
	XYLENE	14.4	. 28,799.0
25	BACETA	10.3	16,351.5
	EACETA	1.9	3,754.0
	FURFUR	7.3	5,840.0
	MEKETO	2.8	2,160.0
· ,	MISDBU	7.9	8,489.6
	TOLJEN	32.1	47,050.8
٦،	XYLENE	44.6	65,863.1
25	ACETON	95.1	51,506.0
		83.8	<u>116,300.0</u> 296,787.0
	CDIDXI Chform	16,933.0	and a second
	CHLORI	87,058.6	0.0 1,304,902.5
	DETHAN	68.1	32,400.0
	EGLYCO	3.8	1,200.0
	EGLIJE	3.6	966.0
	ETHANO	29.7	1,074.0
	FORMAL	113.3	45,118.0

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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 2 HAZARDOUS AIR POLLUTANT SYSTEM STATENIDE SUMMARY OF PROCESS EMISSIONS BY INDUSTRIAL CLASSIFICATION (HAP112)

		•	
	· · ·	AMOUNT USED	EMISSIONS
· · · · ·	POLLUTANT	TUNS/YEAR	LBS./YEAR
	FORMIC	55.9	13,200.0
	HCHLOR	186.4	31,766.0
	HSULFI	378,553.5	269,912.0
	MCELLO	17.9	6,700.0
	MECHLO	17.2	15,900.0
	MEKETO	12.6	21,700.0
	MMERCA	378,315.1	428,940.0
	NITRIC	61.8	11,320.0
	TETHA1	17.4	24,700.0
	TETHYL	8.6	14,700.0
	TITANI	9,039.6	3,580.0
	TOLUEN	3,278.3	1,930,000.0
	TRICHL	75.2	36,300.0
	TURPEN	24,448.0	77,130.0
	XYLENE	190.2	348,900.0
	ZINCZN	89.7	900.0
23	ACETON	88.2	4,804.0
	ANTIMO	6.7	0.0
	BACETA	3.8	0.0
	BARIUM	2.0	0.0
	BUTANO	3.2	40.0
	CHLORI	58,671.1	0.0
	CHROMI	1.2	0.0
	CTETRA	<u> </u>	0.0
	EGLYCO	8.0	320.0
	ETHANO	26.4	0.0
	FALCOH	20.4	0.0
	FURFUR	3.2	1.270.0
		9,607.0	52.6
····	HCHLOR		
		8.5	0.0
	MCELLO	10.5	0.0
	MECHLO	126.5	5,065.0
·····	MEKETO	176.8	134.8
	MERCUR	8.9	16.4
	MMETHA	25,896.0	183,900.0
	NITRIC	2.1	0.0
	PCBIPH	0.5	0.0
	TETHA1	444.4	2,751.0
	TETAYL	222.7	179.0
	TITANI	39.2	0.0
	TOLUEN	119.5	593.0
	TRICHL	8.8	0.0
	XYLENE	327.2	2,243.8
29	MECHLO	1.5	275.0
30	ACETON	535.0	143,920.0
	BACETA	0.1	0.0
	BUTANO	0.1	0.0
	OMETHA	412.5	0.0
	EACETA	1.0	1,860.0
	EPICHL	31.0	12.0

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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 3 HAZARDOUS AIR POLLUTANT SYSTEM STATENIDE SUMMARY OF PROCESS EMISSIONS BY INDUSTRIAL CLASSIFICATION (HAP112)

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	AMOUNT USED	EMISSIONS
POLLUTANT	TONS/YEAR	L8S./YEAR
FORMAL	3,375.0	25,003.0
FORMIC	7.5	1.0
MALEIC	143.0	0.0
MCELLO	0.1	0.0
MECHLO	20.9	35,740.0
MEKETO	63.9	62,031.0
MELAMI	2,321.0	0.0
ММЕТНА	11.6	650.0
PHENOL	3,524.0	3,247.0
STYREN	96.3	226.0
TETHA1	366.1	592,710.0
TOLJEN	17.8	31,725.0
TRICHL	1.2	2,420.0
XYLENE:	1.4	2,200.0
ACETON	250.3	471,927.8
BACETA	97.6	195,262.0
BPHTHA	5.1	10,190.0
BUTANO	10.5	20,892.0
CHLORI	4.6	9,200.0
DMETHA	28.0	56,000.0
EACETA	25.4	45,529.0
EGLYCO	110.1	220,054.0
ETHAND	1.0	2,000.0
FORMAL	1.4	0.0
FORMIC	251.1	10,044.0
HCHLOR	3.5	140.0
HSULFI	0.0	0.Ü
LEADPA	5.0	80.0
MCELLO	309.4	618,932.0
MCHLOR	6.6	13,200.0
MECHLO	13.3	26,582.6
MEKETO	194.3	353,229.0
NALINE	0.9	1,800.0
NPHENO	5.7	228.0
OXALIC	1.7	308.0
PHENOL	2.8	112.0
TETHA1	6.3	10,200.0
TETHYL	33.4	66,750.0
TITANI	93.0	1,660.0
TOLUEN	146.2	287,836.0
XYLENE	79.4	158.847.0
2 <u>ACRYLO</u> MELAMI	<u> </u>	0.0
TDIOXI		0.0
VCHLOR	501.4	0.0
3TETHA1	23.1	28,665.0
4 CYANID	1.6	355.0
EACETA	1.8	2,808.0
EGLYCO	2.5	4,964.0
MCELLO	0.5	1,008.0
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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 4 HAZARDOUS AIR POLLUTANT SYSTEM STATEWIDE SUMMARY OF PROCESS EMISSIONS BY INDUSTRIAL CLASSIFICATION (HAP112)

		AMOUNT USED	EMISSIONS
	POLLUTANT	TONS/YEAR	LBS./YEAR
	MECHLO	1.8	1,032.0
	MEKETO	9.0	18,045.9
	NITRIC	4.7	4,700.0
	PCBIPH	6.2	0.0
		25.8	34,558.0
	TETHA1		
	TOLUEN .	7.6	11,882.0 47,520.0
	TRICHL	39.6	17,876.7
	XYLENE	12.9	
-	ZINCZN	2.3	46.0
<u>5</u> .	BARIUM	2.5	1,782.0
	CYANID	1.3	2,600.0
	EACETA	0.1	35.0
	FORMIC	122.3	4,892.0
	MECHLO	1.4	2,713.0
	MEKETO	0.6	1,182.6
	TETHA1	68.2	101,303.0
	TOLUEN	26.6	53,203.2
	TRICHL	3.1	1,677.0
5	ACETON	32.6	39,899.0
	BACETA	20.7	34,203.0
	BARIUM	43.4	374.0
	CHLORI	48.0	100.0
	C OP > ER	127.7	106.6
	EACETA	2.5	5,000.0
	EGLYCO	5.9	5,941.0
	HCHLOR	170.5	6,672.0
	HSULFI	0.0	0.0
	LEADP8	8.22	550.0
	MECHLO	104.3	190,571.0
	NITRIC	108.5	4,470.0
	OXALIC	6.0	0.0
	PHENOL	2.6	0.7
	TETHA1	221.0	178,089.0
	TETHYL	4.0	7,000.0
	TITANI	1.9	0.0
	TRICHL	82.9	5,948.0
	XYLENE	57.5	19,469.0
1.	ACETON	51.8	59,974.0
	BACETA	0.2	280.0
	BUTAND	9.0	11,020.0
	CHLORI	2.0	0.0
	CHROMI	1.7	136.0
	EACETA	1.0	256.0
	EBENZE	0.5	80.0
	EGLYCO	34.9	40,451.0
	HCHLOR	13.5	5,500.0
	HYDRAZ	0.3	600.0
	LEADPB	160.0	0.0
	MCELLO	4.7	5,608.0
	MECHLO	7.8	6,290.0
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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 5 HAZARDOUS AIR POLLUTANT SYSTEM STATENIDE SUMMARY OF PROCESS EMISSIONS BY INDUSTRIAL CLASSIFICATION (HAP112)

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	POLLUTANT	TONS/YEAR	LBS./YEAR
	MEKETO	11.3	19,790.0
	MISOBU	2.2	2,816.0
			20,960.0
- <u>.</u>	NITRIC	122.7	12,775.0
	STYREN	107.4	150,581.0
	TETHA1	85.9	
·	TETHYL	50.6	40,561.0
	TOLUEN	0,9	
	TRICHL	0.5	1,000.0
	XYLENE	70.4	99,096.0
33	MISOBU	1.6	2,739.0
	TETHA1	18.0	32,709.0
	TOLJEN	8.2	14,223.0
39	ACETON	16.3	32,490.0
	MEKETO	0.8	1,608.0
	TOLUEN	2.9	5,915.0
	XYLENE	1.2	2,353.0
40	BACETA	0.1	200.0
	EACETA	0.8	1,600.0
42	EACETA	1.0	2,000.0
	TRICHL	8.0	16.000.0
49	ASBEST	34.3	0.0
	CHLORI	12,000.0	0.0
	HYDRAZ	3.5	140.0
	MANGAN	6.2	9,320.0
	PCBIPH	95.5	0.0
51	ACETON	75.3	0.0
	BACETA	2.4	0.0
	CHLORI	644.7	0.0
····	DBENZE	2.5	0.0
	EGLYCO	7.7	0.0
	FORMIC	48.2	0.0
	MCELLO	34.5	0.0
	MECHLO	8.22	0.0
	MEKETO	42.6	0.0
	MISOBU	1.0	0.0
	NITRIC	86.2	0.0
	OXALIC	1.3	0.0
	TETHAI	62.3	0.0
	TETHYL	7.1	0.Ŭ
	TOLJEN	56.4	0.0
	TRICHL	20.4	0.0
	XYLENE	42.6	0.0
7.2	TETHYL	0.0	0.0
73	CRESOL	2.5	0.0
<u> </u>	DBENZE	2.5	0.0
	MECHLO	2.5	0.0
75	EGLYCO	6.6	0.0
	TOLJEN	1.5	0.0
75	XYLENE	4.1	8,236.0
30	EOXIDE	2.0	569.0

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MAR 11, 1985 DEPARTMENT OF ENVIRONMENTAL PROTECTION PAGE NO. 6 HAZARDOUS AIR POLLUTANT SYSTEM STATEWIDE SUMMARY OF PROCESS EMISSIONS BY INDUSTRIAL CLASSIFICATION (HAP112)

		AMOUNT USED	EMISSIONS
	POLLUTANT	TONS/YEAR	LBS./YEAR
	XYLENE	1.9	0.0
97	CHLORI	10.2	20,400.0
	MCHLOR	1.8	3,600.0
	MEKETO	6.5	7,600.0
	MEKETO PCBIPH	42.8	0.0
	· TETHA1	. 2.8	5,600.0
<u></u>	TETHYL	9.7	15,260.0
	TETHYL: TFURAN TOLUEN	. 0.2 1.1	400.0 2,200.0
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** APPENDIX F **

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APPENDIX F

TOXICITY SCORING SUMMARY

Pollutant	CAS	RTECS	Categories					S	Tox'.
	Registry Number	Accession Number	С	М	R	A	Tot		Score
	· ·								
America Trievide	1327-53-3	CG3325000	4	4	4	4	16	0	16.0
Arsenic Trioxide	62-75-9	IQ0525000	4	4	4	4	16	Õ	16.0
N-Nitrosodimethylamine	107-13-1	AT5250000	4	4	4	3	15	0.50	15.5
Acrylonitrile Benzene	71-43-2	CY1400000	4	4	4	3	15	0.50	15.5
Chloroform	67-66-3	FS9100000	4	4	4	ž	15	0.50	15.5
Chromium (Potassium Chromate)	7789-00-6	GB2940000	4	4	3	4	15	0.50	15.5
Dioxin	1746-01-6	HP3500000	3	4	4	4	15	0.50	15.5
Epichlorohydrin	106-89-8	TX4900000	4	4	4	3	15	0.50	15.5
Ethyleneimine	151-56-4	KX5075000	3	4	4	4	15	0.50	15.5
Formaldehyde	50-00-0	LP8925000	3	4	4	4	15	0.50	15.5
Nitroso-n-Methylurea	684-93-5	YT7875000	4	4	4	3	15	0.50	15.5
•	100-42-5	WL3675000	4	4	4	3	15	0.50	15.5
Styrene Polychlorinated Biphenyls	11097-69-1	TQ1360000	4	4	4	2	14	1.00	15.0
Vinyl Chloride	75-01-4	KU9625000	4	4	4	2	14	1.00	15.0
Benzo(a)Pyrene	50-32-8	DJ3675000		4	4	3	14	0.58	14.6
Cadmium	7440-43-9	EU9800000	3 3	3	4	4	14	0.58	14.6
Cadmium chloride	10108-64-2	EV0175000	3	4	4	3	14	0.58	14.6
Carbon tetrachloride	56-23-5	FG4900000	4	4	3	3	14	0.58	14.6
Hydrazine	302-01-2	MU7175000	3	4	4	3	14	0.58	14.6
Beryllium	7440-41-7	DS1750000	4	4	1	4	13	1.50	14.5
Bis (chloromethyl) ether	542-88-1	KN1575000	4	4	1	4	13	1.50	14.5
	107-30-2	KN6650000	4	4	1	4	13	1.50	14.5
Chloromethyl methyl ether	107-30-2	KN6650000	4	4	1	4	13	1.50	14.5
Methylchloromethyl Ether	51-75-2	IA1750000	1	4	4	4	13	1.50	14.5
Nitrogen Mustard	53-96-3	AB9450000	3	4	4	2	13	0.96	14.0
Acetamide, N-flouren-2-yl	100-44-7	XS8925000	2	4	3	4	13	0.96	14.0
Benzyl chloride	77-78-1	WS8225000	3	4	2	4	13	0.96	14.0
Dimethyl Sulfate	51-79-6	FA8400000	z	4	4	2	13	0.96	14.0
Urethane	60-11-7	BX7350000	3 3	4	3	3	13	0.50	13.5
Dimethylaminoazobenzene	75-56-9	TZ2975000	3	4	3	3	13	1.00	13.5
Epoxypropane	75-07-0	AB1925000	ĩ	4	4	3	12	1.41	13.4
Acetaldehyde	1332-21-4	CI6475000	4	3	1	4	12	1.41	13.4
Asbestos	2465-27-2	BY3500000	4	4	1	3	12	1.41	13.4
Auramine	92-87-5	DC9625000	4	4	ī	3	12	1.41	13.4
Benzidine	57-57-8	RQ7350000	3	4	1	4	12	1.41	13.4
Beta-Propiolactone	117-81-7	TI0350000	3	4	4	1	12	1.41	13.4
Bis (2-ethylhexyl) phthalate	75-15-0	FF6650000	1	4	4	3	12	1.41	13.4
Carbon Disulfide	7758-98-7	GL8800000	1	4	4	3	12	1.41	13.4
Copper Sulfate	75-21-8	KX2450000	1	4	4	3	12	1.41	13.4
Ethylene Oxide	123-31-9	MX3500000	1	4	3	4	12	1.41	13.4
Hydroquinone	7553-56-2	NN1575000	1	3	4	4	12	1.41	13.4
Iodine	/353-50-2	MHT2/2000	т	5	-	7		T • 47	 ♥ ● T

Pollutant	CAS	RTECS	-		S	S	Tox.		
	Registry	Accession							Score
	Number	Number	С	М	R	Α	Tot		
Ve the landman in a	60-34-4	MV5600000	1	4	٦	4	12	1.41	13.4
Methylhydrazine	59-89-2	QE7525000	4	4	3 1	3	12	1.41	13.4
N-Nitrosomorpholine		CM8050000	3	4	1	4	12	1.41	13.4
Propyleneimine	75-55-8		1	4	4	3	12	1.41	13.4
Zinc Chloride	7646-85-7 [,]	ZHI 400000		4	2	2	12	1.15	13.4
1,2-Dichloroethane	107-06-2	KI0525000	4	4 4	4	2	12	1.15	13.2
Tetrachloroethylene	127-18-4	KX3850000	2		4	2	12	1.15	13.2
Trichloroethylene	79-01-6	KX4550000	2	4	4 2	23	12		12.8
1,3-Propanesultone	1120-71-4	RP5425000	3	4	2			0.82	12.8
Hexachlorobutadiene	87-68-3	ES0700000	2	4	3	3	12	0.82	
Hexamethylphosphamide	680-31-9	TD0875000	3	4	3	2	12	0.82	12.8
Nickel	7440-02-0	QR5950000	3	3	2	4	12	0.82	12.8
Acrolein	107-02-8	AS1050000	1	4	2	4	11.	1.50	12.5
Alpha-Napthalene	134-32-7	QM1400000	4	4	1	2	11	1.50	12.5
Arsenic	7440-38-2	CG0525000	4	4	2	1	11	1.50	12.5
Benzotrichloride	98-07-07	XT9275000	4	2	1	4	11	1.50	12.5
Beta-Napthylamine	91-59-8	QM2100000	4	4	1	2	11	1.50	12.5
Ethyl Benzene	100-41-4	DA0700000	1	4	4	2	11	1.50	12.5
Hydrogen Chloride	7647-01-0	MW4025000	. 1	4	2	4	11	1.50	12.5
Lead	7439-92-1	· 0F7525000	1	4	4	2	11	1.50	12.5
Methylene Chloride	75-09-2	PA8050000	· 1	4	4	2	11	1.50	12.5
Pentachlorophenol	87-87-5	SM6300000	1	2	4	4	11	1.50	12.5
Styrene Oxide	96-09-3	CZ9625000	1	4	4	2	11	1.50	12.5
Toluene	108-88-3	XS5250000	1	4	4	2 2	11	1.50	12.5
o-Toluidine	95-53-4	XU2975000	4	4	1	2	11	1.50	12.5
p-Aminodiphenyl	92-67-1	DU8925000	4	4	1	2	11	1.50	12.5
1,1-Dimethyl Hydrazine	57-14-7	MV2450000	3	4	1	3	11	1.26	12.3
2,4-Toluene Diamine	95-80-7	XS9625000	3	4	1	3	11	1.26	12.3
Acetamide	60-35-5	AB4025000	. 3	4	3	1	11	1.26	12.3
Acrylamide	79-06-1	AS3325000	1	4	3	3	11	1.26	12.3
Alpha Benzene Hexachloride	319-84-6	GV3500000	3	4	1	3	11	1.26	12.3
Dimethylcarbamyl Chloride	79-44-7	FD4200000	• 3	4	1	3	11	1.26	12.3
Diphenylhydrazine	122-66-7	MW2625000	3	4	1	3	11	1.26	12.3
Glycidaldehyde	765-34-4	MB3150000	3	4	1	3	11	1.26	12.3
Mercury	7439-97-6	OV4550000	1	4	3	3	11	1.26	12.3
Phenol	108-95-2	SJ3325000	1	4	3	3	11	1.26	12.3
Diethyl Sulfate	64-67-5	WS8750000	3	4	2	2	11	0.96	12.0
Arsine	7784-42-1	CG6475000	4	1	1	4	10	1.73	11.7
Chlorine	7782-50-5	F02100000	1	4	1	4	10	1.73	11.7
Chromium (Lead Chromate)	7758-97-6	GB2975000	4	4	1	1	10	1.73	11.7
Vinylidine Chloride	75-35-4	YZ8061000	1	4	4	1	10	1.73	11.7
Resorcinol	108-46-3	VG9625000	1	4	1	4	10	1.73	11.7
1,4-Dioxane	123-91-1	JG8225000	3	4	1	2	10	1.29	11.3
3,3-Dimethoxybenzidine	119-90-4	DD0875000	3	4	1	2	10	1.29	11.3
4-Nitrobiphenyl	92-93-3	DV5600000	3	4	1	2	10	1.29	11.3
Chloroprene	126-99-8	EI9625000	1	4	3	2	10	1.29	11.3
Melamine	108-78-1	OS0700000	3	4	1	2	10	1.29	11.3
Methyl Chloride	75-09-2	P A 6300000	1	3	4	2	10	1.29	11.3
Methyl Methacrylate	80-62-6	OZ5075000	1	4	3	2	10	1.29	11.3
Napthalene	91-20-3	QJ0525000	1	4	2	3	10	1.29	11.3
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Pollutant	CAS	RTECS		Cat	ego	rie	S	S	Tox.
	Registry	Accession	С	М	D	A	Tot		Score
	Number	Number	C	IAI	К	M	101		
									_
Phenylhydrazine	100-63-0	MV8925000	1	4	2	3	10	1.29	11.3
Vinyl Cyclohexene Dioxide	106-87-6	RN8640000	3	4	1	2	10	1.29	11.3
1,1,1-Trichloroethane	71-55-6	KJ2975000	1	3	3	3	10	1.00	11.0
Methyl Iodide	74-88-4	PA9450000	3	3	1	3	10	1.00	11.0
Rotenone	83-79-4	DJ2800000	1	3	3	3	10	1.00	11.0
2,5-Diaminotoluene	95-70-5	XS9700000	1	- 4	1	3	9	1.50	10.5
3,3-Dichlorobenzidine	91-94-1	DD0525000	3	4	1	1	9	1.50	10.5
4,4-Methylene Dianiline	101-77-9	BY5425000	1	4	1	3	9	1.50	10.5
Aldicarb	116-06-3	UE2275000	1	3	1	4	9	1.50	10.5
Aniline	62-53 - 3	BW6650000	1	4	1	3	9	1.50	10.5
Formic Acid	64-18-6	LQ4900000	1	4	1	3	9	1.50	10.5
Methyl Mercaptan	74-93-1	PB4375000	1	4	1	3	9	1.50	10.5
Picric Acid	88-89-1	TJ7875000	1	4	1	3	9	1.50	10.5
n-Butylamine	109-73-9	E02975000	1	4	1	3	9	1.50	10.5
p-Chloronitrobenzene	100-00-5	CZ1050000	1	4	1	3	9	1.50	10.5
p-Nitrophenol	100-02-7	SM2275000	1	4	1	3	9	1.50	10.5
p-Phenylenediamine	106-50-3	SS8050000	1	4	1	3	9	1.50	10.5
1,1,2-Trichloroethane	79-00-5	KJ3150000	2	4	1	2	9	1.26	10.3
Allyl chloride	107-05-1	UC7350000	1	4	2	2	9	1.26	10.3
Nitrobenzene	98-95-3	DA6475000	1	2	2	4	9	1.26	10.3
Xylene	1330-20-7	ZE2100000	1	2	4	2	9	1.26	10.3
Manganese Chloride	7773-01-5	000962500	1	3	3	2	9	0.96	10.0
Acrylic Acid, Ethyl Ester	140-88-5	AT0700000	1	4	1	2	8	1.41	9.4
Barium Chloride	10361-37-2	CQ8750000	1	1	2	4	8	1.41	9.4
Chlorine Dioxide	10049-04-4	F03000000	1	1	2	4	8	1.41	9.4
Chrysene	218-01-9	GC0700000	2	4	1	1	8	1.41	9.4
Ethylene Glycol Ethyl Ether	110-80-5	KK8050000	1	1	4	2	8	1.41	9.4
Formamide	75-12-7	LQ0525000	1	2	4	1	8	1.41	9.4
Methyl Cellosolve	109-86-4	KL5775000	1	1	4	2	8	1.41	9.4
N-Phenyl-Beta-Napthylamine	133-88-6	QM4550000	2	4	1	1	8	1.41	9.4
Nitroglycerine	55-63-0	OX2100000	1	2	1	4	8	1.41	9.4
Phosphorus	7723-14-0	TH3500000	1	1	2	4	8	1.41	9.4
Vanadium Pentoxide	1314-62-1	YW2450000	1	2	1	4	8	1.41	9.4
Vinyl Bromide	593-60-2	KU8400000	1	4	1	2	8	1.41	9.4
p-Anisidine	104-94-9	BZ5450000	1	4	1	2	8	1.41	9.4
Acrylic Acid	79-10-7	AS4375000	1	1	3	3	8	1.15	9.2
Barium Carbonate	513-77-9	CQ8600000	1	1	3	3	8	1.15	9.2
Ethyl Ether	60-29-7	KI5775000	1	3	1	3	8	1.15	9.2
Furfural	98-01-1	LT7000000	1	3	1	3	8	1.15	9.2
Furfuryl Alcohol	98-00-0	LU9100000	1	3	1	3	8	1.15	9.2
Maleic Anhydride	108-31-6	ON3675000	1	3	1	3	8	1.15	9.2
Quinoline	91-22-5	VA9275000	1	3	1	3	8	1.15	9.2
p-Chloroaniline	106-47-8	BX0700000	1	3	1	3	8	1.15	9.2
1,1,2,2-Tetrachloroethane	79-34-5	KI8575000	2	2	1	3	8. 7	0.82	8.8
Antimony	7440-36-0	CC4025000	1	1	1	4	7	1.50	8.5
Cyanide	57-12-5	GS7175000	1	1	1	4	7	1.50	8.5
Cyanogen	460-19-5	GT1925000	1	1	1		7	1.50	8.5
Fluorine	7782-41-4	LM6470000	1	1 1	1 1	4 4	7 7	1.50 1.50	8.5 8.5
Hexachloronapthalene	1335-87-1	QJ7350000	T	T	T	4	/	1.00	0.0

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Pollutant	CAS	RTECS	, Ca	tego	rie	S	S	Tox.
	Registry	Accession	a 14			m ,		Score
	Number	Number	СM	R	A	Tot		
				•				
Indregen Granida	74-90-8	MW6825000	1 1	1	4	7	1.50	8.5
Hydrogen Cyanide	7783-06-4	MX1225000	$\begin{array}{ccc} 1 & 1 \\ 1 & 1 \end{array}$	1	4	, 7	1.50	8.5
Hydrogen Sulfide	624-83-9	NQ9450000	1 1		4	, 7	1.50	8.5
Methyl Isocyanate	75-44-5	SY5600000	1 1		4	, 7	1.50	8.5
Phosgene Tetrachlorodibenzofuran	51207 - 31-9	HP5295000	1 1		4	, 7	1.50	8.5
	7440-28-0	XG3425000	1 1		4	, 7	1.50	8.5
Thallium Vizul Elugrido	75-02-5	YZ7351000	1 4		1	, 7	1.50	8.5
Vinyl Fluoride	78-87-5	TX9625000	1 2		3	, 7	0.96	8.0
1,2-Dichloropropane	584-84-9	CZ6300000	1 2 1 2	ī	3	, 7	0.96	8.0
2,4-Toluene Diisocyanate	92 - 52-4	DU8050000	1 3		2	7		8.0
Biphenyl Die zamethene	334-88-3	PA7000000	3 2		1	7	0.96	8.0
Diazomethane	77-47-4	GY1225000	1 1	2	3	7	0.96	8.0
Hexachlorocyclopentadiene	78-93-3	EL6475000	1 1	3	2	, 7	0.96	8.0
Methyl Ethyl Ketone	106-51-4	DK2625000	1 2		3	, 7	0.96	8.0
Quinone	109-99-9	LU5950000	$1 \frac{1}{3}$		2	7	0.96	8.0
Tetrahydrofuran ·	108-39-4	GO1250000	1 1		3	7	0.96	8.0
m-Cresol	99-65-0	CZ7350000	1 2		3	7	0.96	8.0
m-Dinitrobenzene	104-91-6	SM4725000	1 2		3	7	0.96	8.0
p-Nitrosophenol	67-64-1	AL3150000	1 2 1 2		2	, 7	0.50	7.5
Acetone	84-66-2	TI1050000	1 2		2	7	0.50	7.5
Diethyl phthalate	532-27-4	AM6300000	1 1	1	3	6	1.00	7.0
2-Chloroacetophenone	7726-95-6	EF9100000	1 1		3	6	1.00	7.0
Bromine	1307-96-6	GG2800000	1 1	ī	3	6	1.00	7.0
Cobalt Oxide	420-04-2	GS5950000	1 1		3	6	1.00	7.0
Cyanimide Hydrogen Bromide	10035-10-6	MW3850000	1 1	1	3	6	1.00	7.0
	75-31-0	NT8400000	1 1		3	6	1.00	7.0
Isopropylamine Nitric Acid	7697-37-2	QU5775000	1 1		3	6	1.00	7.0
Oxalic Acid	144-62-7	R02450000	1 1		3	6	1.00	7.0
1,2,4-Trichlorobenzene	120-82-1	DC2100000	1 1		2	6	0.58	6.6
1,3-Butadiene	106-99-0	EI9275000	1 2		2	6	0.58	6.6
Butanol	71-36-3	E01400000	1 2		2	6	0.58	6.6
Isoamyl Alcohol	123-51-3	EL5425000	1 2		2	6	0.58	6.6
Pyridine	110-86-1	UR8400000	1 2		· 2	6	0.58	6.6
n-Butyl Acetate	123-86-4	AF7350000	$\overline{1}$ $\overline{1}$	2	2	6	0.58	6.6
p-Nitroaniline	100-01-6	BY7000000	$\overline{1}$ $\overline{2}$		2	6	0.58	6.6
p-Nitrotoluene	99-99-0	XT3325000	1 2		2	6	0.58	6.6
1,2-Dichlorobenzene	95-50-1	CZ4500000	1 1	1	2	5.	0.50	5.5
1-Nitropropane	108-03-2	TZ5075000	1 1	1	2	5	0.50	5.5
2,2-Iminodiethanol	111-42-2	KL2975000	1 1	1	2	5	0.50	5.5
Acetic Anhydride	108-24-7	AK1925000	1 1	1	2	5	0.50	5.5
Butanethiol	109-79-5	EK6300000	1 1	1	2	5	0.50	5.5
Cyclohexane	110-82-7	GU6300000	1 2	1	1	5	0.50	5.5
Diphen. meth. 4,4-diisocyan.	101-68-8	NQ9350000	1 2	1	1	5	0.50	5.5
Ethanethiol	75-08-1	KI9625000	1 1	1	2	• 5	0.50	5.5
Ethanolamine	141-43-5	KJ5775000	1 1	1	2	5	0.50	5.5
Ethyl Acetate	141-78-6	AH5425000	1 1	1	2	5	0.50	5.5
Ethyl Chloride	75-00-3	KH7525000	1 1	1	2	5	0.50	5.5
Isophorone	78-59-1	GW7700000	1 1	1	2	5	0.50	5.5
Ketene	463-51-4	OA7700000	1 1	1	2	5	0.50	5.5

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Pollutant	CAS Registry	RTECS Accession		Cat	ego	orie	S	S	Tox. Score
Number				М	R	A	Tot		30016
Monochlorobenzene	108-90-7	CZ0175000	1	1	1	2	5	0.50	5.5
Potassium Cyanate Selenium	590-28-3 7782-49-2	GS6825000 VS7700000	1 1	1 1	1 2	2 1	5 5	0.50 0.50	5.5 5.5
Sodium Cyanate	917-61-3	GS7000000	1	ī	1	2	5	0.50	5.5
Turpentine	8006-64-2	Y08400000	1	1	1	2	5	0.50	5.5
Xylidine	1300-73-8	ZE8575000	1	1	1	2	5	0.50	5.5
m-Nitrotoluene	99-08-1	XT2975000	1	1	1	2	5	0.50	5.5
Cyanoacetamide	107-91-5	AB5950000	1	1	1	1	4	0	4.0
Diiosoctyl Phthalate	27554-26-3	TI1300000	1	1	1	1	4	0	4.0
Diisodecyl Phthalate	26761-40-0	TI1300000	1	1	1	1	4	0	4.0
Ethylene	74-85-1	KU5340000	1	1	1	1	4	0	4.0
Methyl Isobutyl Ketone	108-10-1	SA9275000	1	1	1	1	4	0	4.0
Octachloronapthalene	2234-13-1	QK0250000	1	1	1	1	4	0	4.0
Terephthalic Acid	100-21-0	WZ0875000	1	1	1	1	4	0	4.0
Titanium Oxide	13463-67-7	XR2275000	1	1	1	1	4	0	4.0

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** APPENDIX G **

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APPENDIX G

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TOXICITY RANKING FOR HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE

Pollutant	Toxicity Score	Emissions (lbs/yr)	Toxicity Rank	Emissions Rank	Tota
Arsenic	16.0	430	1	50	51
Chromium	15.5	184	2	54	56
Epichlorhydrin	15.5	12	2	58	60
Formaldehyde	15.5	79,401	. 2	21	23
Styrene	15.5	13,001	2	32	34
Benzene	15.5	52,392	2	24	26
Hydrazine	14.6	740	7	49	56
Cadmium	14.6	238	7	52	59
Benzo-a-Pyrene	14.6	15,180	7	30	37
Epoxypropane	13.5	124,600	10	19	29
Bis 2-ethylhexyl phthalate	13.4	10,190	11	34	45
Copper	13.4	107	11	55	66
Ethylene Oxide	13.4	1,535	11	45	56
Zinc	13.4	946	11	47	58
1,2 Dichloroethane	13.2	32,400	15	28	43
Tetrachloroethylene	13.2	708,050	15	6	21
Trichloroethylene	13.2	110,865	15	20	35
Hydrogen Chloride	12.5	44,731	18	26	44
Lead	12.5	178,630	18	16	34
Methylene Chloride	12.5	288,569	18	13	31
Napthalene	12.5	11,799	18	33	51
Ethyl Benzene	12.5	80	18	56	74
Toluene	12.5	2,511,623	18	1	19
Mercury	12.3	16	24	57	81
Phenol	12.3	5,180	24	37	61
Diethyl Sulfate	12.0	5,424	26	38	64
Chlorine	11.7	1,433,003	27	3	30
Methyl Chloride	11.3	16,8 0 0	28	29	57
Methyl Methacrylate	11.3	184,550	28	15	43
1,1,1-Trichloroethane	11.0	1,496,539	30	2	32
Méthyl Mercaptan	10.5	428,940	31	9	40
p-Nitrophenol	10.5	228	31	53	84
Formic Acid	10.5	50,433	31	25	56
Xylene	10.3	1,091,274	34	4	38
Manganese	10.0	9,320	35	35	70
Barium	9.4	2,156	36	44	80
Chlorine Dioxide	9.4	296,787	36	10	46
Methyl Cellosolve	9.4	632,248	36	7	43
Ethylene Glycol Ethyl Ether	9.4	277,118	36	11	47
Furfural	9.2	7,110	40	36	76
Hydrogen Sulfide	8.5	269,912	41	12	53
Cyanide	8.5	2,955	41	42	83

TOXICITY RANKING FOR HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE (con't

Pollutant	Toxicity Score	Emissions (1bs/yr)	Toxicity Rank	Emissions Rank	Tota:
Biphenyl	8.0	4,510	43	40	83
Methyl Ethyl Ketone	8.0	512,611	43	8	51
Tetrahydrofuran	8.0	400	43	51	94
Acetone	7.5	804,521	46	5	51
Nitric Acid	7.0	41,450	47	27	74
Oxalic Acid	7.0	868	47	48	95
n-Butyl Acetate	6.6	253,563	49	14	63
Butanol	6.6	149,490	49	17	66
Acetic Anhydride	5.5	1,403	51	46	97
1,2 Dichlorobenzene	5.5	2,400	51	43	-94
Diphenyl Methyl 4,4-Diisocyanate	5.5	146,000	51	18	69
Ethyl Acetate	5.5	66,537	51	23	74
Ethanolamine	5.5	3,318	51	41	92
Turpentine	5.5	77,130	51	22	73
Metĥyl Isobutyl Ketone	4.0	14,045	57	31	88
Titanium Oxide	4.0	5,240	57	39	96

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** APPENDIX H **

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APPENDIX H

EXPOSURE RANKING FOR HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE

Pollutant	Toxicity Score	Emissions (lbs/yr)	Toxicity Rank	Emissions Rank	Total
Toluene	12.5	2,511,623	18	1	19
1,1,1-Trichloroethane	11.0	1,496,539	30	2	32
Chlorine	11.7	1,433,003	27	3	30
Xylene	10.3	1,091,274	34	4	38
Acetone	7.5	804,521	46	5	51
Tetrachloroethylene	13.2	708,050	15	6	21
Methyl Cellosolve	9.4	632,248	36	. 7	43
Methyl Ethyl Ketone	8.0	512,611	43	· 8	51
Methyl Mercaptan	10.5	428,940	31	9	40
Chlorine Dioxide	9.4	296,787	36	10	46
Ethylene Glycol Ethyl Ether	9.4	277,118	36	11	47
Hydrogen Sulfide	8.5	269,912	41	12	53
Methylene Chloride	12.5	288,569	18	13	31
	6.6	253,563	49	14	63
n-Butyl Acetate	11.3	184,550	28	15	43
Methyl Methacrylate	12.5	178,630	18	16	34
Lead	6.6	149,490	49	17	66
Butanol	5.5	146,000	51	18	69
Diphenyl Methyl 4,4-Diisocyanate	13.5	124,600	10	19	29
Epoxypropane	13.2	110,865	15	20	35
Trichloroethylene	15.5	79,401	2	21	23
Formaldehyde	5.5	77,130	51	22	73
Turpentine	5.5	66,537	51	23	74
Ethyl Acetate	15.5	52,392	2	24	26
Benzene	10.5	50,433	31	25	56
Formic Acid	12.5	44,731	18	26	44 [°]
Hydrogen Chloride	7.0	44,751	47	27	74
Nitric Acid	13.2	32,400	15	28	43
1,2 Dichloroethane	11.3	16,800	28	29	57
Methyl Chloride		15,180	28	30	37
Benzo-a-Pyrene	14.6		57	31	88
Methyl Isobutyl Ketone	4.0 15.5	14,045	2	32	34
Styrene		13,001	18	33	51
Napthalene	12.5	11,799	18	34	45
Bis 2-ethylhexyl phthalate	13.4	10,190	35	35	70
Manganese	10.0	9,320	40	36	76
Furfural	9.2	7,110		. 37	61
Phenol	12.3	5,180	24 26	38	64
Diethyl Sulfate	12.0	5,424		39 ·	96
Titanium Oxide	4.0	5,240	57	40	90 83
Biphenyl	8.0	4,510	43		92
Ethanolamine	5.5	3,318	51	41	
Cyanide	8.5	2,955	41	42	83 94
1,2 Dichlorobenzene	5.5	2,400	51	43	
Barium	9.4	2,156	36	44	80

Pollutant	Toxicity	Emissions	Toxicity	Emissions	Total
	Score	(lbs/yr)	Rank	Rank	
Ethylene Oxide	13.4	1.535	11	45	56
Acetic Anhydride	5.5	1,403	51.	46	97
Zinc	13.4	946	11	47	58
Oxalic Acid	7.0	868	47	48	9 5
Hydrazine	14.6	740	7	49	56
Arsenic	16.0	430	1	50	51
Tetrahydrofuran	8.0	400	43	51	94
Cadmium	14.6	238	7	52	59
p-Nitrophenol	10.5	228	31	53	84
Chromium	15.5	184	2	54	56
Copper	13.4	107	11	55	66
Ethyl Benzene	12.5	80	18	56	74
Mercury	12.3	16	24	57	81
Epichlorhydrin	15.5	12	2	58	60
Epicnlornyarin	15.5	12	2	58	00

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EXPOSURE RANKING FOR HAZARDOUS AIR POLLUTANTS EMITTED IN MAINE (con't)

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** NOTES **

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