# RESEARCH TRIANGLE INSTITUTE

## **JULY 1993**

# TOXIC RELEASE INVENTORY DATA VERIFICATION: COMPARISON OF TOXIC RELEASE INVENTORY (TRI) DATA TO OTHER ENVIRONMENTAL DATA BASES

## FINAL REPORT

Prepared for:

North Carolina Department of Health, Environment, and Natural Resources Pollution Prevention Program Raleigh, North Carolina

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#### **FINAL REPORT**

## TOXIC RELEASE INVENTORY DATA VERIFICATION: COMPARISON OF TOXIC RELEASE INVENTORY (TRI) DATA TO OTHER ENVIRONMENTAL DATA BASES

#### **1.0 INTRODUCTION**

In 1986, the Environmental Protection Agency (EPA) introduced a Federal law that required certain facilities to report their annual releases of toxic materials to the environment. This law, the Emergency Planning and Community Right-to-Know Act, is described in Section 313 of Title III of the Superfund Amendments and Reauthorization Act (SARA). It states that those facilities meeting certain criteria are required to complete and file the Toxic Release Inventory (TRI) Reporting Form annually. The criteria are as follows and all must apply:

- The facility has 10 or more full-time employees.
- The facility conducts manufacturing operations within the Standard Industrial Classification (SIC) codes 20 through 39.
- The facility manufactures, processes, or in any other way uses any of the listed toxic chemicals in amounts greater than the threshold quantities.

The threshold quantities for manufacturers and processors are:

- 75,000 lb during the 1987 calendar year,
- 50,000 lb during the 1988 calendar year, and
- 25,000 lb during the 1989 calendar year and in subsequent years.

The threshold quantities for users are as follows:

• 10,000 lb during the 1987 calendar year and in subsequent years.

The Section 313 toxic chemical list contains over 300 specific chemicals and about 20 chemical categories. However, these numbers may fluctuate yearly due to the addition or removal (delisting) of chemicals. Facilities are required to complete one TRI form for each listed chemical that is manufactured, processed, or used in quantities exceeding the threshold levels. Specifically, the forms contain the quantities of each chemical that are being released to air, water, and land. These forms cover the activity of the preceding calendar year and

are filed by July of each year. Facilities are required to submit TRI forms to EPA and their State governments.

In this way, TRI provides a complete, useful inventory of the toxic chemicals that are being released to the environment. Consequently, TRI data are used in a wide variety of efforts including those with significant impacts on environmental priorities and policy at both the Federal and State levels. It is crucial, therefore, that the TRI data be as accurate as possible.

The two primary objectives of this project are: (1) to determine the accuracy of and identify errors in TRI reporting data in the State of North Carolina for the years 1987 through 1990, and (2) to develop computer methodologies that can be used to compare TRI data with data in other environmental data bases. This report contains the results obtained from implementing the second objective. A separate report documents the results from the first objective.

To support North Carolina in its pollution prevention efforts, RTI has developed a multimedia waste reduction management system (WRMS) for use in identifying pollution prevention opportunities, targeting State resources, and tracking pollution prevention progress in the State. WRMS is a relational data base that integrates existing multimedia environmental release and waste generation data from the TRI and several single-media data bases. The purpose of this report is to describe how WRMS was used in an effort to verify TRI data by comparing TRI data in WRMS to toxics data reported in single-media data bases and by comparing TRI data across reporting years.

The goal of this effort was to develop computerized methodologies for implementing comparisons between TRI data in WRMS and toxics data reported to single-media data bases and to identify significant discrepancies within TRI and between the TRI and single-media data bases. Five specific comparisons were identified by the North Carolina Pollution Prevention Program (NCPPP):

- 1) Compare TRI data to release data in single-media data bases.
- 2) Identify facilities reporting releases above the TRI reporting thresholds to singlemedia data bases but not to TRI.
- 3) Compare TRI data across reporting years.
- 4) Compare facilities reporting under TRI to the State of North Carolina with the facilities reporting to EPA.
- 5) Link and compare TRI data to SARA Title III Section 312 data.

The methodology and results from the first three of these comparisons are described in the remainder of this report. The fourth comparison could not be completed because the necessary data on facilities reporting under TRI to the State of North Carolina were not available in computerized format, and neither project resources nor the Scope of Work allowed for computerization of these data from hard copy. The fifth comparison also could not be completed because the Section 312 data were not available in computerized form for the data years covered by WRMS (1987 through 1990). These data were available in computerized form for 1991 and 1992, and they were examined to assess the possibilities of such a link and comparison in future years. It was determined that the Section 312 data contained no facility identifiers in common with TRI data that would allow a computerized link between the two data bases.

This report is divided into two major sections. The Methodology section briefly describes WRMS and the single-media data bases and details the methods used to make each of the three comparisons. The Results section presents results of the comparisons. The computer programs used to implement these comparisons are included in Appendices A through D. The results of each comparison are presented in Appendices E through G. A detailed description of WRMS and its development may be found in A Multimedia Waste Reduction Management System for the State of North Carolina, prepared by RTI for NCPPP in April 1993.

#### 2.0 METHODOLOGY

WRMS is a dBASE IV data base that contains data from TRI on releases of toxics to air, water, and land and transfers to publicly owned treatment works (POTWs) or offsite hazardous waste treatment, storage, or disposal facilities. WRMS also contains data on hazardous wastes, conventional water pollutants, and criteria air pollutants reported by TRI facilities from the following single-media sources:

- North Carolina Annual Report (NCAR) on Hazardous Waste Data Base,
- National Pollutant Discharge Elimination System (NPDES) Data Base,
- Municipal/Industrial Pretreatment Monitoring Data Base, and
- Air Quality National Emissions Data System (NEDS).

The hazardous waste, conventional water pollutant, and criteria air pollutant data included in WRMS from the above sources cannot be compared to the TRI data in WRMS because they cover different categories of pollutants. For the comparisons between TRI and the single-media data bases, toxics data from the single-media sources (which were not included in WRMS so as not to create overlaps in coverage) were required. Not all of these data sources contain toxics data that could be compared with TRI data; only the NPDES and Pretreatment data bases were found to contain toxics data that could be compared directly to TRI data. NCAR contains data only on hazardous wastes. Although these were not toxics data, per se, in many cases, the toxics reported to TRI were probably contained in the hazardous wastes reported in NCAR; therefore, the NCAR data were also compared with the TRI data for constituents that might be found in each waste. NEDS currently contains only data on criteria air pollutants, so no comparison to TRI data could be made. However, the State of North Carolina is beginning to collect air toxics data, and these data may be available in the future for comparisons with TRI data. In summary, toxics data from the NPDES and Pretreatment data bases and hazardous waste data from NCAR were compared to TRI.

The toxics data from the NPDES and Pretreatment data bases had not been incorporated in WRMS, so the first step in the comparison was to convert them to the WRMS format. In order not to include them in WRMS itself, the converted data were stored in a separate dBASE file. The conversion process had two main components: matching facilities in the single-media data to those in TRI and converting facility, chemical, and emissions data to the WRMS format. All potential TRI facilities (i.e., manufacturing facilities) were included from the single-media data bases, whether they matched a facility already in WRMS or not. Nonmanufacturing facilities were excluded. Similarly, only toxic chemicals from the TRI list were included from the single-media data bases. For NPDES and Pretreatment data, concentration and flow data had to be converted to emissions data. The hazardous waste data were already in WRMS, so no additional conversion was needed.

The program used to convert these data, VCONVERT.PRG, is a modified version of the conversion program used to convert NPDES and Pretreatment data for WRMS. That conversion program (CONVERT.PRG) is described in detail in *A Multimedia Waste Reduction Management System for the State of North Carolina* (prepared for NCPPP by RTI, 1993). The program code for the verification conversion program is listed in Appendix A.

The rest of this section describes the methodology used for each comparison. Each methodology was implemented as a dBASE program. The program listings can be found in Appendices B through D.

#### 2.1 Comparison of TRI Data to Single-Media Data

The single-media data bases available for this comparison (as discussed above) were the NPDES and Pretreatment data bases, which cover direct releases to water and transfers to POTWs, respectively, and the NCAR hazardous waste data. NPDES data were compared to TRI water release data, Pretreatment data were compared to TRI POTW transfer data, and hazardous waste data were compared to TRI offsite transfers. Because air toxics data may be

available in the near future, the methodology was designed to include comparisons for air. Air toxics data would be compared to TRI stack air release data.

Due to the different nature of the comparison of hazardous waste data to TRI data, this comparison was done differently than that for NPDES and Pretreatment. In addition, no formal computer methodology was developed for the hazardous waste comparison. The following discussion applies only to the NPDES and Pretreatment comparisons.

Records with missing data or data based on a reported range were excluded from the comparison. For example, NPDES and Pretreatment data are both reported as concentration in a waste stream. If the flow data for a facility was missing, these concentration data could not be converted to emissions data. These records were retained in the converted data, but were not used in the comparison. Similarly, TRI data may be reported as a range; these data were retained in WRMS but were not used in the comparison.

For each year of data, the facilities and chemicals reported in each single-media data base were matched with those reported to TRI and the quantities reported compared. If a chemical was reported by a facility to either the single-media data base or TRI but not to both, the quantity for the data base to which it was not reported was treated as zero. Not reported is distinguished from a reported quantity of zero, however. It should be noted that nonzero releases might not be reported due to threshold reporting requirements. However, because the thresholds for TRI are not release quantities but manufactured quantities, there is no clear minimum release value that can be used for unreported releases found in other data bases.

The quantities reported to each data base were compared and the differences between them calculated. All differences were calculated as positive values, regardless of which data base reported the higher value; in this way, differences could be summed across chemicals for a facility without equal but opposite discrepancies cancelling each other out.

To determine what level of discrepancy was significant, the data for each single-media data base were examined and the differences compared to typical quantities reported to each single-media data base. Based on this examination, the significance level for the NPDES data base was set at 1,000 lb (i.e., a difference in the quantity reported to TRI versus NPDES of 1,000 lb or more would be considered significant), and the significance level for the Pretreatment data base was set at 10,000 lb. The level for the Pretreatment data base was set at higher because the quantities reported to the Pretreatment data base are typically much higher than those reported to NPDES, and using a significance level of 1,000 lb would have resulted in an enormous output for the Pretreatment comparison. Because no data were available for air toxics, the significance level for air was set to 1,000 lb. This may need to be modified when the air toxics data become available.

For each facility with one or more chemicals showing a discrepancy over the significance level, the significant discrepancies were summed to the facility level and the facilities sorted by total discrepancy. The data for those facilities and the chemicals with significant discrepancies were then output in a report, with the largest total discrepancies first. The report shows facility identification information, chemical name, and the quantities reported to TRI and the single-media data base.

This methodology was computerized in a dBASE program called TASK6.PRG. The program prompts the user to choose a single-media data base and a year for comparison, performs all calculations, and outputs the results to a text file named T6xyy.TXT where x is a one letter code indicating the single-media data base selected (N = NPDES, P = Pretreatment, A = air) and yy is the year selected (e.g., 87, 88, etc.). For example, the output file for a comparison of TRI and Pretreatment data for 1989 would be T6P89.TXT. The TASK6 program is listed in Appendix B.

#### 2.2 Comparison of Single-Media Data to TRI Reporting Thresholds

The only single-media data bases available for this comparison (as discussed above) were the NPDES and Pretreatment data bases. Because air toxics data may be available in the near future, the methodology was designed to include comparisons for air. Release data for each single-media data base were compared to the reporting thresholds for TRI. Hazardous waste data were not compared to TRI reporting thresholds, since hazardous wastes may contain multiple constituents.

TRI has two reporting thresholds: one for chemicals "manufactured or processed," and one for chemicals "otherwise used." The manufactured or processed threshold started at 75,000 lb per year (lb/yr) for the 1987 reporting year, and decreased to 50,000 lb/yr in the 1988 reporting year and 25,000 lb/yr in the 1989 reporting year and thereafter. The otherwise used threshold is 10,000 lb for all reporting years. Note that these thresholds are for a quantity of chemical manufactured, processed, or otherwise used, not for quantities released. In fact, a facility could manufacture a chemical in quantities exceeding the threshold and release none of it. Similarly, another facility might manufacture the same chemical in quantities below the threshold but release most of it. The first facility would be required to report to TRI and the second would not. If a facility releases a quantity of a chemical that exceeds these reporting thresholds, then presumably that facility is manufacturing, processing, or otherwise using that chemical in quantities that exceed the threshold and should, therefore, be reporting to TRI. It is also possible that a facility reporting releases below these thresholds to single-media data bases is in fact manufacturing, processing, or otherwise using the chemical in quantities exceeding the TRI reporting thresholds, but since the single-media data bases do not report quantity manufactured, processed, or otherwise used, that more direct comparison cannot be made.

Records with missing data were excluded from the comparison. For example, NPDES and Pretreatment data are both reported as concentration in a waste stream. If the flow data for a facility were missing, then concentration data could not be converted to emissions data. These records were retained in the converted data but were not used in the comparison.

For each year of data, facilities and chemicals were compared to those reported to TRI. If a chemical was reported by a facility to the single-media data base but not to TRI, the quantity reported was compared to both the manufactured or processed threshold for that reporting year and the otherwise used threshold. Any release exceeding one of the TRI reporting thresholds and not reported to TRI was considered significant. Any releases either reported to TRI or not exceeding either threshold were excluded from further analysis. For the purposes of this comparison, a chemical was considered reported to TRI by a facility if there was a record in TRI for that facility and chemical, even if the reported quantity was zero. As was mentioned earlier, it is theoretically possible to exceed the TRI reporting threshold without actually releasing any of the chemical. Although a quantity exceeding the threshold reported to TRI may represent a significant discrepancy, reporting was the issue in this comparison and not discrepancies in quantity.

Facilities with significant releases were divided into two groups: facilities reporting one or more chemicals over the manufactured or processed threshold, and facilities reporting no chemicals over the manufactured or processed threshold but one or more over the otherwise used threshold. For each facility with one or more chemicals showing a discrepancy, all releases in excess of the otherwise used threshold (which is the lower of the thresholds) were summed to the facility level, and the facilities sorted by total quantity of these releases. The data for those facilities and the chemicals exceeding the otherwise used threshold were then output in a report, with the largest total releases first. The report shows facility identification information, chemical name, and the quantity reported to the single-media data base.

This methodology was computerized in a dBASE program called TASK8.PRG. The program prompts the user to choose a single-media data base and a year for comparison, performs all calculations, and outputs the results to a text file named T8xyy.TXT where x is a one letter code indicating the single-media data base selected (N = NPDES, P = Pretreatment, A = air) and yy is the year selected (e.g., 87, 88). For example, the output file for a comparison of TRI thresholds and Pretreatment data for 1989 would be T8P89.TXT. The TASK8 program is provided in Appendix C.

#### 2.3 Comparison of TRI Data Across Reporting Years

Three steps were used to compare the TRI data across the reporting years 1987 through 1990: reporting patterns, stopped reporting, and quantity discrepancies. These analyses were applied sequentially to each facility/chemical combination (corresponding to one Form R, the

form used to report to TRI); if a facility/chemical combination was flagged as a discrepancy in any step, it was not considered in later steps. Each step is described in the following paragraphs.

In the reporting pattern step, the data were examined for irregularities in a facility's pattern of reporting for a particular chemical. There are a variety of reasons that facilities may change their reporting status from one year to the next, including changes in reporting thresholds, chemicals listed and delisted, and changes in production practices at a facility. Irregularities were changes in reporting status that did not obviously fit into one of these scenarios. One common theme of the above reasons for change in reporting status is that changes resulting from them would likely be permanent or long term. For example, if a chemical is delisted, it is not likely to be relisted the next year. Similarly, facilities are not expected to change production practices for one year then change back. They could but it would not be typical. If a facility stopped reporting for any of these reasons, it would not be expected to start reporting again. Therefore, irregularities were defined as facilities that stopped reporting then resumed reporting after a period of not reporting (e.g., they report in one or more years, then do not report in one or more years, then report again in one or more years). In other words, reporting was considered irregular if all reporting years were not contiguous--if there was a gap in reporting. So a facility that does report, then does not report, then does report would be considered irregular because it has reported in two noncontiguous years.

In looking at reporting patterns, the quantity reported was not considered. Not reporting was also distinguished from reporting a release of zero. Facilities that reported a release of zero for a chemical were considered to have reported.

Facility/chemical combinations showing pattern discrepancies were output in a report showing the facility and chemical information and the reporting status in each year of the analysis. Facilities are shown in alphabetical order, as there was no way to rank a particular reporting pattern as "worse" than another. These facility/chemical combinations were then omitted from further consideration.

In the stopped reporting step, facilities that stopped reporting during the time period analyzed were examined for the total quantity reported in the last year they reported. The total quantity released was the sum of stack, fugitive, water, land, POTW, and offsite releases. If this quantity was greater than the manufactured or processed reporting threshold for the next reporting year, the facility/chemical combination was flagged as showing a discrepancy. The manufactured or processed reporting threshold was 50,000 lb in 1988 and 25,000 lb in 1989 and thereafter. So if a facility reported a chemical only in 1987, the quantity would be compared to the threshold for 1988, the next year in which they might have reported. Facility/chemical combinations showing stopped reporting discrepancies were output in a report showing the facility and chemical information and the quantity reported or nonreporting status in each year of the analysis. This report is ordered by the last quantity reported, with the largest first. Each facility/chemical combination is treated separately, so a facility may have additional chemicals elsewhere in the report. The facility/chemical combinations in the stopped reporting discrepancies report were omitted from further consideration.

Finally, the quantities reported over time were compared for significant discrepancies. Only total quantity released was compared; the total quantity released and transferred was the sum of stack, fugitive, water, land, POTW, and offsite releases. Each year's quantity was compared to the next year's quantity. Two basic approaches to these comparisons were considered. The first was an absolute approach, in which the absolute difference in quantity between two reporting years was calculated and compared to a fixed cutoff. The second was a relative approach, in which the relative change in quantity between two reporting years was calculated and compared to a cutoff. The data spanned so many orders of magnitude that a log scale was used for this measure. The relative difference was calculated as:

Difference = 
$$\log_{10}\left(\frac{\text{Larger Quantity}}{\text{Smaller Quantity}}\right)$$

This results in a number representing the number of orders of magnitude difference between the two quantities (e.g., a value of 2 would indicate a difference of two orders of magnitude).

Neither of these approaches by themselves seemed to give adequate results; therefore, they were combined. A facility/chemical combination was considered to show a significant discrepancy if, over the years they reported, any two years showed a relative difference in quantity of more than one order of magnitude and an absolute difference in quantity of 3,000 lb or more. These cutoffs were determined by examination of the data.

Facility/chemical combinations showing quantity discrepancies were output in a report showing the facility and chemical information and the quantity reported or nonreporting status in each year of the analysis. The actual relative and absolute measures of discrepancies for each pair of years was not output, nor are the facilities ordered (except by alphabetical order), as there was no one measure of the magnitude of the discrepancy that could be used to order the data.

This methodology was computerized in a dBASE program called TASK7.PRG. The program can perform the analysis on up to five years of data. It analyzes the number of years available; if there are fewer than five, all years are analyzed. If there are more than five, the program prompts the user to choose a starting year for the analysis. Five years

starting with that year are analyzed (or all years available after the starting year, if that is fewer than five). Reporting pattern analysis is only performed if three or more years of data are available or selected for analysis. The program performs all calculations and outputs the results to three text files named T71.TXT, T72.TXT, and T73.TXT. These files contain the results of each step in the analysis: T71.TXT contains the results of reporting pattern analysis; T72.TXT contains the results of the stopped reporting analysis; and T73 contains the results of the quantity discrepancy analysis. The TASK7 program is listed in Appendix D.

#### 3.0 RESULTS

The results of the comparisons described in Section 2 are shown in Appendices E through G. Appendix E contains the results for the comparison of TRI data to single-media data for NPDES, Pretreatment, and NCAR for reporting years 1987 through 1990. Appendix F presents the results for the comparison of single-media data to TRI thresholds for NPDES and Pretreatment for reporting years 1987 through 1990. Appendix G shows the results for the comparison of TRI data across reporting years over the period 1987 to 1990. For confidentiality reasons, all facility identifying information in the Appendices has been replaced with the facility's SIC code and the results sorted by SIC code.

For all of these results, it should be noted that a significant discrepancy identified in the results does not necessarily indicate incorrect data or the failure of a facility to report correctly. The different data bases compared have different reporting requirements and procedures that could result in legitimate discrepancies among data reported to different data bases. Similarly, for the comparison of TRI data over time, legitimate changes in operations and production levels or reporting requirements could result in significant discrepancies. In addition, data entry errors could result in discrepancies where, in fact, there are none. Finally, NPDES and Pretreatment data had to be converted from average concentration and flow data to emissions, which may have introduced errors. The results are not intended to be a definitive identification of reporting problems, but a screening tool for targeting NCPPP's future efforts.

#### 3.1 Comparison of TRI Data to Single-Media Data

Detailed results of the comparison of TRI data to single-media data bases are presented in Appendix E. By far the most compelling result of this comparison for NPDES and Pretreatment was the rarity of finding the same chemical reported by the same facility to both TRI and NPDES or TRI and Pretreatment. Virtually all of the significant discrepancies took the form of a significant quantity of a chemical reported released by a facility to one of the data bases and not reported at all to the other. Not reported is distinguished from a reported quantity of zero in the comparison. Given that most of the significant discrepancies stem from differences in reporting (rather than actual quantities reported), it should be restated that there are a variety of legitimate reasons possible for these discrepancies, including the different reporting requirements of TRI versus the NPDES or Pretreatment data bases.

Similarly, the hazardous waste comparison to TRI constituents reflects many discrepancies. There are many reasons why these discrepancies might appear. Hazardous wastes many contain additional constituents not usually associated with a particular waste, constituents not on the TRI list, TRI constituents falling below the TRI reporting threshold, or nontoxic components, all of which would contribute to the overall waste quantity but not to the sum of TRI constituents. Also, specific constituents may be associated with several different waste codes and so may not have been a part of the waste to which they are being compared. To fully identify the reasons for discrepancies at any particular facility, a detailed waste characterization of that facility's hazardous wastes would be needed.

Table I summarizes the number of facilities with significant discrepancies and the largest single discrepancy for each data base and year. For both NPDES and Pretreatment, the number of facilities with significant discrepancies decreased steadily from 1987 to 1990. Similarly, the largest discrepancy also decreased for both comparisons over that period. For both data bases, the chemical sodium sulfate (solution) appears frequently at the high end ofthe discrepancy list for 1987. In all cases, sodium sulfate (solution) was reported to TRI in large quantity and not reported to either the NPDES or Pretreatment data bases. This chemical was delisted from TRI in 1989, with the delisting taking effect retroactively for the 1988 reporting year, and it does not appear in later reporting years. In no instance was it seen reported to the NPDES or Pretreatment data bases, indicating that it likely was not required to be reported to those data bases.

The discrepancies between the NPDES and TRI data nearly all take the form of chemicals reported to TRI and not to NPDES. In no case was the same chemical reported by the same facility to both the NPDES and TRI data bases. The discrepancies between the TRI data and the Pretreatment data are more evenly divided between chemicals reported to TRI but not Pretreatment and chemicals reported to Pretreatment but not TRI. In a few cases (about four per year) the same chemical was reported by the same facility to both with significant discrepancies (greater than 10,000 lb difference between the two reported quantities).

#### 3.2 Comparison of Single-Media Data to TRI Reporting Thresholds

Detailed results of the comparison of single-media data to the TRI thresholds are presented in Appendix F. Few facilities reported releases exceeding the TRI reporting thresholds that did not report these releases to TRI. For NPDES, only one facility reported releases not reported to TRI that exceeded the TRI manufactured or processed threshold in 1987 and 1990; no facilities reported releases not reported to TRI that exceeded either the manufactured or processed or the otherwise used threshold for 1988 and 1989. Some

### TABLE I SUMMARY OF DISCREPANCIES BETWEEN TRI AND SINGLE-MEDIA DATA BASES

	1987	1988	1989	1990
	Total Number	of Facilities R	eporting	
TRI	779	742	863	618
NPDES	65	72	78	80
Pretreatment	539	544	761	791
Numb	er of Facilities	with Significar	nt Discrepancies	
• NPDES vs TRI	65	45	34	33
Pretreatment vs TRI	192	144	108	99
	Maximun	n Discrepancy	(lb)	
NPDES vs TRI	60 million	530,000	85,000	340,000
Pretreatment vs TRI	2.7 billion	1.3 billion	310 million	260 million

Notes:

- Maximum discrepancies are rounded.
- Because the number of facilities with significant discrepancies includes facilities that reported to only one of the data bases being compared, this number may exceed the total number of facilities reporting to either of the data bases individually.

Pretreatment facilities reported releases not reported to TRI that exceeded one or both of the TRI thresholds for each reporting year. It should be recalled that much larger quantities are typically reported to the Pretreatment data base than to the NPDES data base, so the difference in results between the two is not surprising.

Table II summarizes the number of facilities with releases exceeding each of the TRI thresholds that were not reported to TRI for each data base and year. For both NPDES and Pretreatment, this number of facilities did not change significantly between 1987 and 1990, despite the decreasing manufactured or processed threshold.

#### 3.3 Comparison of TRI Data Across Reporting Years

Detailed results of the comparison of TRI data across reporting years are presented in Appendix G. A large number of discrepancies were found: 278 facilities showed irregular reporting patterns, 268 facilities stopped reporting after reporting quantities over the reporting threshold the previous year, and 145 facilities had quantity discrepancies over time. While there may be overlap among these groups of facilities (a facility could be in one group for one chemical and another group for a different chemical), this is still a significant fraction of the approximately 1,200 facilities reporting to TRI in the period examined.

There are a variety of legitimate reasons that may account for some of the discrepancies found. Changes in the actual reporting requirements or in the interpretation of reporting requirements from year to year may have resulted in irregular reporting or stopped reporting. One major explanation for many of the stopped reporting discrepancies is the delisting of certain chemicals. Table III lists the chemicals delisted each year since TRI started. The delisting of sodium hydroxide in 1989 explains the failure of many facilities to report sodium hydroxide in 1989 after two years of reporting very large quantities. Sodium sulfate solution is a similar case (although shown as delisted in 1989, the delisting was retroactive to 1988, explaining the absence of reporting for sodium sulfate solution in 1988 after large quantities were reported in 1987). The interpretation of all these results should be made in light of such changes in reporting requirements.

## TABLE II SUMMARY OF FACILITIES EXCEEDING TRI REPORTING THRESHOLDS BUT NOT REPORTING TO TRI

	1987	1988	1989	1990
	Total Number	of Facilities Repo	rting	
NPDES	65	72	78	80
Pretreatment	539	544	761	791
Number of Facilit	ies Exceeding TI	<b>RI</b> Manufactured	or Processed Thres	hold
NPDES	1	0	0	1
· Pretreatment	28	17	31	23
Number of	Facilities Exceedi	ng TRI Otherwis	e Used Threshold*	
NPDES	0	0	0	0
Pretreatment	34	11	11	11

<sup>a</sup> Does not include facilities shown as exceeding the manufactured or processed threshold.

Manufactured or Processed Threshold:

75,000 lb in 1987 50,000 lb in 1988 25,000 lb in 1989 and thereafter

Otherwise Used Threshold:

10,000 lb for all reporting years

# TABLE IIICHEMICALS DELISTED FROM TRI BETWEEN 1988 AND 1991

Reporting Year <sup>a</sup>	Chemicals Delisted
1988	CI Acid Blue 9, diammonium salt
	CI Acid Blue 9, disodium salt
	Titanium dioxide
1989	Melamine
	Sodium sulfate (solution)
	Sodium hydroxide
1990	Aluminum oxide
	Terephthalic acid
1991	CI Pigment Blue 15 (from Copper Compounds)
	CI Pigment Green 7 (from Copper Compounds)
	CI Pigment Green 36 (from Copper Compounds)

<sup>\*</sup>This is the reporting year for which the Form R instructions first note the chemical as delisted, typically the same year as the *Federal Register* notice delisting the chemical. However, chemicals are sometimes delisted retroactively, and any unnecessary reports deleted from (or never entered to) the TRIS data base, from which TRI data for WRMS were obtained. For example, sodium sulfate (solution) was delisted in 1989, retroactive to 1988, and all reports for sodium sulfate (solution) for 1988 have been omitted from TRIS.

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APPENDIX A PROGRAM CODE FOR VCONVERT.PRG

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Filename: VCONVERT.PRG Anne E. Crook, Research Triangle Institute Author: 2/15/93 Last updated: 2/16/93 Created: Comments: Main WRMS Data Base Data Verification Conversion Utility program for converting data from NPDES and Pretreatment data base formats to WRMS format. This program uses a two-level menu: the first (menu\_1) lets the user choose a data source to convert from (e.g., NPDES); the second lets them select a conversion activity. There are 3 conversion activities: facility matching, chemical matching, and emissions conversion, which must be done in that order. In addition, the program creates a working file or files for use in the conversion program. The main VCONVERT program calls other programs to perform each of these activities (creating working files, facility matching, chemical matching, and emissions conversion). Each data source has its own program for each of these activities, since each data source is different and requires different operations to be converted. The VCONVERT program expects to find the following subdirectories on the default drive: \WRMS \WRMS\NPDES . \WRMS\PRETREAT It expects the main WRMS data files and the VCONVERT program itself to be in \WRMS, and the data source specific program files, original data files, working file(s), and any ancillary files needed by the conversion program to be in the appropriate subdirectory listed above. \* Call the procedure that initializes environment with SET commands, defines \* menus, and defines memvars DO initial \* Main program loop to display menus; the loop redisplays the main menu after \* lower level ones return control, until the user selects quit and the loop is exited. DO WHILE .T.  $mexit_1 = .F.$ && set menu\_1 exit flag to false source = space(10)&& initialize data source variable DO show\_1 && display menu\_1 DO eval\_1 && evaluate menu\_1 choice \* If user selected EXIT, exit menu\_1 loop, otherwise continue IF mexit\_1 EXIT ENDIF IF source = space(10)LOOP ENDIF \* Check for working data file, and create it from original data if it does \* not exist. If original data file not found, display error message and \* return to menu\_1. If working data exists, restore status information \* that tells where user left off. \* The working, original, and source variables are set when the user makes

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* a choice in menu_1.
  IF .NOT. file(working)
                                         && no working file
   IF .NOT. file(original)
                                           && no original data file either
     @ 1,0 CLEAR
                                             && Display error message
     @ 9,27 TO 13,52 DOUBLE
     @ 11,29 SAY "Source file not found."
     @ 22,50 SAY "Press any key to continue ..."
     WAIT ""
     LOOP
                                             && go back to menu 1
                                           && original data file found
   ELSE
     DO &source.vwork
                                          && call prog to create working file
     SAVE TO &source.vstatus ALL LIKE 1_*
                                            && create status variable file
   ENDIF
                                             && so user can resume later
  ELSE
                                         && working file found
   RESTORE FROM & source.vstatus ADDITIVE
                                           && restore status of where user
  ENDIF
                                           && left off
* Loop to display menu_2
 DO WHILE .T.
   mexit_2 = .F.
                                         && set menu_2 exit flag to false
   DO show_2
                                         && display menu_2
   DO eval 2
                                         && evaluate menu_2 choice
* If user selected EXIT or completed conversion, exit menu_2 loop; otherwise
* continue
    IF mexit_2 .OR. 1_emiss
     EXIT
   ENDIF
* Save status memory variables for use next time to determine where to pick up
* conversion.
   SAVE TO &source.vstatus ALL LIKE 1_*
 ENDDO
                                          && restart menu_2 loop
* User has exited from menu_2 or completed conversion for selected data source
ENDDO
                                          && restart menu_1 loop
* User has exited from menu_1 (main menu)
* Reset environment, close all files, quit to dBase
DO reset
RETURN && vconvert.prg
PROCEDURE initial
* Initialize program environment
  CLEAR ALL
                           && closes all files, clears screen and memory
  SET BELL OFF
                           && Turns off warning bell
 SET CLOCK OFF
                           && Turns off clock display
  SET DEVICE TO SCREEN
                           && Directs all output to the screen
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&& Disables exact matching SET EXACT OFF && Disables help SET HELP OFF SET SAFETY OFF && Suppresses overwrite warnings && Turns off delete, insert, caps, and numlock SET SCOREBOARD OFF && indicators SET STATUS OFF && Turns off status line SET TALK OFF && Suppresses echo of command progress \* Initialize variables to be shared by all programs called by VCONVERT These are status variables that reflect how much of the conversion \* process has been completed (and therefore, where the user left off) \* so that the user can resume in the same place later. Working and original \* hold the names of the working and original data files that the programs \* expect to find for each data source. PUBLIC 1\_fac\_done, 1\_chm\_done, 1\_emiss, working, original \* Call the procedure that defines the menus DO defmenus RETURN && initial PROCEDURE defmenus \* These are popup menus displayed in the middle of the screen \* The PROMPT text appears on the menu itself; the MESSAGE text appears below \* the menu and clarifies the menu prompt \* Define menu\_1 to choose data source SET BORDER TO DOUBLE DEFINE POPUP menu\_1 FROM 8,28 TO 14,50 DEFINE BAR 1 OF menu\_1 PROMPT "NPDES"; MESSAGE "Convert data from NPDES format" DEFINE BAR 2 OF menu\_1 PROMPT "Pretreatment"; MESSAGE "Convert data from Pretreatment format" DEFINE BAR 3 OF menu\_1 PROMPT "AIRS"; MESSAGE "Convert data from AIRS format" DEFINE BAR 5 OF menu\_1 PROMPT "Exit to dBase"; MESSAGE "Exit WRMS Conversion Utility to dBase" ON SELECTION POPUP menu\_1 DEACTIVATE POPUP \* Define menu\_2 to choose conversion activity DEFINE POPUP menu\_2 FROM 8,28 TO 14,50 DEFINE BAR 1 OF menu\_2 PROMPT "Match Facilities"; MESSAGE "Match facilities to facilities already in FACILITY file" DEFINE BAR 2 OF menu\_2 PROMPT "Match Chemicals"; MESSAGE "Match chemicals to chemicals already in CHEMICAL file" DEFINE BAR 3 OF menu\_2 PROMPT "Convert Emissions"; MESSAGE "Convert emissions data to WRMS format" DEFINE BAR 5 OF menu\_2 PROMPT "Exit to Source Menu"; MESSAGE "Return to Data Source Menu" ON SELECTION POPUP menu\_2 DEACTIVATE POPUP RETURN && defmenus PROCEDURE show\_1 \* Display menu\_1 with screen header and instructions CLEAR 00,1 SAY "WRMS DATA VERIFICATION CONVERSION UTIL

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I T Y"
 @ 5,9 SAY "Choose a data source using the arrow keys and press <Enter>"
 @ 17,10 TO 20,67
 @ 18,11 SAY "This utility converts data from one of the above formats"
 @ 19,11 SAY "to the WRMS format for data verification."
 ACTIVATE POPUP menu_1
RETURN && show_1
PROCEDURE eval_1
• Evaluate menu_1 choice
* Under each data source choice, the source, working, and original variables
* are set to reflect the user's data source choice. The source variable will
* be appended to the beginning of subroutine calls to select the subdirectory
* and program specific to the chosen data source. The working and original
* variables tell VCONVERT what the names of the working and original data files
* are for the chosen data source, so it can see if they exist or (in the case
* of the working file) need to be created. Some data source choices also
* initialize some additional status variables for use by called programs.
 DO CASE
   CASE bar() = 5
                                 && User selected Exit
     mexit_1 = .T.
                                    && Set exit flag to true
                                 && User selected NPDES
   CASE bar() = 1
     source = "npdes\N"
     working = "npdes\nconcext.dbf"
     original = "npdes\npdes.dat"
     PUBLIC l_cauto
   CASE bar() = 2
                                 && User selected Pretreatment
     source = "pretreat\P"
     working = "pretreat\preext.dbf"
     original = "pretreat\pretreat.dat"
   CASE bar() = 3
                                 && User selected AIRS
     DO not_avail
      source = \operatorname{air}A
      working = "air\airsext.dbf"
*
      original = "air\airs.dat"
 ENDCASE
RETURN && eval_1
PROCEDURE show_2
Display menu_2 with instructions. Header should still be displayed from
* menu_1.
 @ 1,0 CLEAR
 § 5,7 SAY "Choose a conversion activity using the arrow keys and press <Enter>"
 @ 17,5 TO 19,75
 @ 18,6 SAY "Facility and Chemical Matching MUST be done before Convert Emissions."
 ACTIVATE POPUP menu_2
RETURN && show_2
PROCEDURE eval_2
* Evaluate menu_2 choice
* Conversion must be done in the correct order: Facility matching first,
* Chemical meatching second, and Emissions conversion last. If the user
* tries to select a choice out of this order, a message will be displayed
* telling them what segments they must do before the one they chose.
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\* Similarly, if the user selects a segment that has already been done, a \* message will be displayed indicating that. In both cases, the menu will

\* be redisplayed after the message is displayed.

DO CASE CASE bar() = 5&& User selected Return to main menu  $mexit_2 = .T.$ && Set exit flag to True CASE bar() = 1&& User selected Facility Matching && Facility matching already done IF l\_fac\_done DO already && Display message ELSE && Facility matching not already done && Call program to do facility matching DO &source.vfac ENDIF && User selected Chemical Matching CASE bar() = 2DO CASE && Chemical matching already done CASE l\_chm\_done && Display message DO already && Facility matching has been done CASE l\_fac\_done DO &source.vchem && Call program to do Chemical matching OTHERWISE && Facility matching not done yet && Direct user to do Facility matching •DO nofac && before selecting Chemical matching ENDCASE && User selected Emission conversion CASE bar() = 3DO CASE && Emission conversion already done CASE l\_emiss && Display message DO already && Chemical matching has been done (implies CASE 1\_chm\_done && Facility matching done also) DO &source.vemiss && Call program to do Emission conversion && Facility matching done, but not Chemical CASE l\_fac\_done && matching (otherwise, previous CASE would && have been true) && Direct user to do Chemical matching DO no\_chem && before selecting Emissions conversion && Neither Facility nor Chemical matching done OTHERWISE && (otherwise, a previous CASE would have been && true) && Direct user to do both Facility and DO no\_fac\_chm && Chemical matching before selecting && Emissions conversion ENDCASE ENDCASE RETURN && eval\_2 PROCEDURE reset \* Reset default environment; this reverses the SET commands in INITIAL CLEAR ALL SET ESCAPE ON SET EXACT OFF SET HELP ON SET SAFETY ON SET SCOREBOARD ON

SET STATUS ON SET TALK ON RETURN && reset PROCEDURE already \* Display message telling user they have already done the selected task @ 1,0 CLEAR @ 9,18 TO 13,58 DOUBLE @ 11,20 SAY "You have already completed this task." @ 22,50 SAY "Press any key to continue ..." WAIT "" RETURN && already PROCEDURE no\_fac \* Display message telling user to match facilities before chemicals @ 1,0 CLEAR @ 9,21 TO 14,56 DOUBLE @ 11,23 SAY "You must Match Facilities before" @ 12,23 SAY "you can Match Chemicals." @ 22,50 SAY "Press any key to continue ..." WAIT "" RETURN && no\_fac PROCEDURE no\_chem \* Display message telling user to match chemicals before converting emissions @ 1,0 CLEAR @ 9,21 TO 14,55 DOUBLE @ 11,23 SAY "You must Match Chemicals before" @ 12,23 SAY "you can Convert Emissions." @ 22,50 SAY "Press any key to continue ..." WAIT "" RETURN && no\_chem PROCEDURE no\_fac\_chm \* Display message telling user to match facilities and chemicals before \* converting emissions @ 1,0 CLEAR @ 9,19 TO 14,61 DOUBLE @ 11,21 SAY "You must Match Facilities and Chemicals" @ 12,21 SAY "before you can Convert Emissions." @ 22,50 SAY "Press any key to continue ..." WAIT "" RETURN && no\_fac\_chm PROCEDURE not\_avail \* Display message telling user that the selected option is not available. @ 1,0 CLEAR @ 9,24 TO 13,55 DOUBLE

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@ 11,26 SAY "This option is not available."
@ 22,50 SAY "Press any key to continue..."
WAIT ""
RETURN && not_avail
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* EOF: vconvert.prg
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\* Filename: NVWORK.PRG \* \* Author: Anne E. Crook, Research Triangle Institute \* \* Created: 2/15/93 Last updated: 2/24/93 \* Called by:VCONVERT \* Comments: Creates working NPDES files for verification conversion program, with structure extended to include fields used in processing. Also cleans up some fields. -\*\*\*\*\* \* Initialize status variables. These variables indicate which steps of the \* conversion process have been completed.  $l_fac_done = .F.$ && Facility matching done  $l\_chm\_done = .F.$ && Chemical matching done  $l\_cauto = .F.$ && Auto chemical matching done  $1_{emiss} = .F.$ && Emissions conversion done \* Initialize other variable. myear = space(2)&& Year to be converted \* Ask for year being converted. @ 1,0.CLEAR SET CONFIRM ON @ 9,20 SAY "Enter year of data being converted: 19" @ 9,58 GET myear PICTURE \*99\* READ SET CONFIRM OFF \* Create intermediate NPDES data file - the original data file is in ASCII \* format; this just pulls it into dBase format. \* SET TALK ON echos the number of records processed to the screen so that the \* user can see the progress of the records being added. @ 1,0 CLEAR @ 8,11 SAY "Creating intermediate file ...." CREATE npdes\npdesint FROM npdes\nintstru SET TALK ON APPEND FROM npdes.dat TYPE SDF \* Convert averages from character to numeric format, and consolidate weighting \* variables (only one of cnt19 and cnt26 will be nonzero for a record). REPLACE ALL avg WITH val(cavg), weight WITH cnt19 + cnt26 SET TALK OFF @ 9,0 CLEAR CLOSE DATABASE @ 8,43 SAY "Done" \* The intermediate file is divided into two working files, one for flow data \* and one for pollutant concentration and loading data. @ 9,11 SAY "Creating Flow file ....." \* Create working Flow file. Pipe index is for linking to concentration file. \* Parameter code 50050 is flow; units code 03 is MGD. All flow values should \* be in MGD; any that aren't are not used. CREATE npdes\nflowext FROM npdes\flowstru INDEX ON npdes + pipe\_num TAG pipe SET TALK ON

APPEND FROM npdes/npdesint FOR param\_code = "50050" .AND. param\_units = "03"

SET TALK OFF @ 10,0 CLEAR CLOSE DATABASE @ 9,43 SAY "Done"

\* Create working Concentration file.

@ 11,11 SAY "Creating Concentration file ..." CREATE npdes\nconcext FROM npdes\concstru

\* Create index on NPDES number, conditional on it being nonstandard (i.e., not \* matching standard format of NC followed by numbers). The index will be used \* later in cleaning up nonstandard NPDES numbers.

INDEX ON npdes TAG npdes\_c FOR .NOT. like("NC\*", npdes) .OR. "O" \$ npdes

\* Create unique index on NPDES number. The index will be used later in \* creating new VFAC records for unmatched facilities.

INDEX ON npdes TAG npdes\_u UNIQUE

\* Create index on parameter code, conditional on no chemical match found. \* The index will be used in chemical matching.

INDEX ON param\_code TAG param\_c FOR chem\_mis

\* Create.index unique to pipe and parameter. The index will be used in \* emissions conversion.

INDEX ON npdes + pipe\_num + param\_code + param\_name TAG pipeparamu UNIQUE

\* Create index on pipe and parameter, but not unique. The index will be used \* in emissions conversion.

INDEX ON npdes + pipe\_num + param\_code + param\_name TAG pipeparam

\* Copy records with nonzero concentration values from intermediate data file. \* Parameter code 50050 is flow; all others are concentration or loading values. SET TALK ON APPEND FROM npdes\npdesint FOR param\_code <> "50050" .AND. avg > 0 SET TALK OFF @ 12,0 CLEAR CLOSE DATABASE @ 11,43 SAY "Done" \* Enter year in working Concentration file and bad flow and concentration files. @ 14,11 SAY "Entering year in working file ..." USE npdes\nconcext REPLACE ALL year WITH val("19" + myear) @ 14,45 SAY "Done"

\* Clean up NPDES numbers in working Concentration file. This involves \* replacing capital Os in the NPDES number with zeros. NPDES numbers are not \* cleaned up in the flow file, since this could result in two flow records for \* the same facility. Multiple concentration records for the same facility \* can be weighted and aggregated, but not flows. Therefore, only flow records \* with good NPDES numbers are used.

@ 15,11 SAY \*Cleaning up NPDES ids .....\* USE npdes\nconcext SCAN FOR \*O\* \$ npdes

\* Repeat this loop until there are no capital Os in the NPDES number.

DO WHILE "0" \$ npdes
 REPLACE npdes\_1 WITH stuff(npdes,at("0",npdes),1,"0")

ENDDO

ENDSCAN @ 15,45 SAY "Done"

\* Some NPDES numbers may still be nonstandard even after clean up to change \* capital Os to zeros. This section of the program lets the user examine non-\* standard NPDES numbers and either fix or delete them. This uses the \* conditional index created when the working Concentration file was created.

USE npdes\nconcext ORDER npdes\_c GO TOP && Go to the first record with a nonstandard no. IF .NOT. EOF() && If it's at the end of the file, then there are && no nonstandard NPDES nos. and nothing need be done CLEAR TYPE npdes\nvwork1.msg && Displays instructions for resolving nonstandard WAIT ""

\* Browse the nonstandard NPDES numbers; the user is not allowed to add or
\* delete records. The user has the option of either fixing the NPDES number so
\* that it is in the standard format or deleting it. Either of these actions
\* removes the record from the conditional index. When all records have been
\* removed from the index in this way, an error condition results; the ON ERROR
\* CLEAR line clears the screen when this happens, effectively ending the
\* BROWSE mode.
ON ERROR CLEAR
BROWSE NOFOLLOW NOAPPEND NODELETE FIELDS npdes
ON ERROR

ENDIF

\* Restore screen header wiped out by BROWSE command.

CLEAR

@ 0,1 SAY "WRMS DATA VERIFICATION CONVERSION UTILI TY"

CLOSE ALL RETURN && nvwork.prg

\* EOF: nvwork.prg

nvwork1.msg

WRMS DATA VERIFICATION CONVERSION UTILITY

Directions for cleaning up NPDES ids:

- \* You will be shown records with non-standard NPDES ids, i.e., those that don't start with NC. IDs containing capital O instead of zero will also be shown.
- \* If the problem is a typo (e.g., MC instead of NC) fix it.
- \* Change capital O to zero (0).
- \* Press <Ctrl-End> when finished.

Press any key to continue ...

✤ Filename: NVFAC.PRG \* Author: Anne E. Crook, Research Triangle Institute 2/15/93 Last updated: 2/23/93 Created: \* Called by:VCONVERT \* Comments: NPDES facility matching routine. NPDES facilities are matched to TRIS facilities in WRMS based on NPDES number. If no match is found, the NPDES record is deleted. Facilities are only matched \* in the working Concentration file; records in the working Flow ÷ file that correspond to unmatched facilities will simply be ignored during the emissions conversion step. \* Open working Concentration file and WRMS Facility file. Files are linked on \* NPDES number. USE npdes\nconcext IN 1 ALIAS conc USE facility IN 2 ORDER npdes ALIAS fac SET RELATION TO npdes INTO fac @ 1,0 CLEAR \* For all working Concentration file records with a corresponding WRMS Facility \* file record, copy the WRMS Facility ID from the Facility file record to the \* working Concentration file record. \* SET TALK ON echos the number of records processed to the screen so that the \* user can see the progress of the facility matching. @ 9,16 SAY "Assigning FAC\_IDs to matched facilities ..." SET TALK ON REPLACE conc->fac\_id WITH fac->fac\_id FOR fac->fac\_id <> space(10) SET TALK OFF @ 10,0 CLEAR @ 9,60 SAY "Done" \* Delete unmatched nonmanufacturing facilities from working Concentration file. @ 10,16 SAY "Deleting nonmanufacturing facilities ....." SET TALK ON DELETE FOR fac\_id = space(10) .AND.; (val(left(sic,2)) < 20 .OR. val(left(sic,2)) > 39) PACK SET TALK OFF @ 11,0 CLEAR @ 10,60 SAY "Done" CLOSE ALL USE npdes\nconcext IN 1 ORDER npdes\_u ALIAS conc USE facility IN 2 ORDER fac\_id ALIAS fac SET RELATION TO fac\_id INTO fac USE vfac IN 3 ORDER npdes ALIAS vfac \* Copy facility info for unmatched manufacturing facilities to VFAC. @ 11,16 SAY "Creating new facility records ......" newvfacs = 0SCAN && conc SELECT vfac SEEK conc->npdes IF .NOT. found() newvfacs = newvfacs + 1 APPEND BLANK REPLACE fac\_name WITH conc->fac\_name,; street\_1 WITH conc->street,; city WITH conc->city,;
county WITH conc->county,; zip WITH conc->zip,; sic\_1 WITH conc->sic,; npdes WITH conc->npdes,; year\_added WITH conc->year IF conc->fac\_id <> space(10)
 REPLACE contact WITH fac->contact, phone WITH fac->phone ENDIF @ 11,60 SAY newvfacs PICTURE \*999\* ENDIF SELECT conc ENDSCAN @ 11,60 SAY "Done" \* Update status variable to indicate facility matching is complete. l\_fac\_done = .T. @ 22,24 SAY "Facility Match Complete. Press any key to continue ..." WAIT "" CLOSE ALL RETURN && nvfac.prg \* EOF: nvfac.prg

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* Filename: NVCHEM.PRG
* Author: Anne E. Crook, Research Triangle Institute
* Created: 2/15/93 Last updated: 2/16/93
* Called by:VCONVERT
* Comments: NPDES chemical matching routine. Checks chemicals in working
          Concentration file for match in WRMS Chemical file and assigns a
          chemical id.
* Initialize variables
m_continue = "Y"
                         && does user want to continue
* The status variable 1_cauto indicates whether the user has completed auto
* chemical matching. If interactive checking of unmatched chemicals is done,
* the whole chemical matching process is complete, as indicated by l_chm_done.
* If the auto chemical match has not been done, do it.
IF .NOT. l_cauto
 DO auto
* Auto chemical match is complete. If chemical matching is not complete
 (because there are unmatched chemicals left) see if the user wants
* to quit and resume later (if they do, the ASK_QUIT procedure will return
* control to VCONVERT).
 IF .NOT. 1_chm_done
   DO ask_quit
 ENDIF
ENDIF
* If chemical matching is not done, call procedure to let user check unmatched
* chemicals.
IF .NOT. l_chm_done
 DO mis
* If user quit before chemical matching was done, return control to VCONVERT.
 IF .NOT. 1_chm_done
   RETURN
 ENDIF
* Chemical matching is complete; delete records with unmatched chemicals.
 @ 1,0 CLEAR
 @ 12,12 SAY "Deleting unmatched chemicals..."
 USE npdes\nconcext IN 1 ALIAS conc
 SET TALK ON
 DELETE FOR chem_id = space(10)
 PACK
 @ 13,0 CLEAR
 @ 12,44 SAY "Done"
 CLOSE ALL
ENDIF
RETURN && nvchem.prg
PROCEDURE ask_quit
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\* See if user wants to quit and resume later. @ 1,0 CLEAR @ 8,12 TO 15,66 @ 10,14 SAY "Press Y to continue chemical matching," @ 11,14 SAY "Press N to exit and resume chemical matching later." @ 13,32 SAY \*Continue? Y/N:\* @ 13,47 GET m\_continue PICTURE "Y" READ \* If the user does want to quit, return control to VCONVERT. IF m\_continue = "N" RETURN TO MASTER ENDIF RETURN && ask\_quit PROCEDURE auto \* Initialize variables. && Count number of NPDES concentration records matched match = 0 $no_match = 0$ && Count number of NPDES concentration records not matched && Count number of NPDES concentration records processed processed = 0\* Open working Concentration file and WRMS Chemical file. Files are unlinked. \* Chemical file order is on STORET parameter number. USE npdes\nconcext IN 1 ALIAS conc USE chemical IN 2 ALIAS chem ORDER TAG storet\_num SELECT conc \* Determine total number of NPDES Concentration records to process. total\_recs = reccount("conc") \* Display matching status report on screen. DO report \* Main loop; processes each record in working Concentration file. SCAN processed = recno("conc") && Set number of records processed SELECT chem \* Look for matching STORET parameter code in WRMS Chemical file. SEEK conc->param\_code \* If a nonblank match is found, increment match counter and copy chemical ID to \* working Concentration file. IF found() .AND. conc->param\_code <> space(2) match = match + 1REPLACE conc->chem\_id WITH chem->chem\_id ELSE \* No match found; increment no match counter and set unmatched flag in working \* Concentration file.

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no_match = no_match + 1
REPLACE conc->chem_mis WITH .T.
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ENDIF

\* Update matching status report on screen.

DO says

\* Select working Concentration file for correct functioning of SCAN loop. SELECT conc -

ENDSCAN

\* Write final chemical matching status report to file and screen.

SET DEVICE TO FILE npdes\nchmauto.txt DO report SET DEVICE TO SCREEN DO report \* Update status variables  $1_cauto = .T.$ \* IF no chemical were left unmatched, chemical matching is complete; set \* l\_chm\_done accordingly. IF no\_match = 0 $l_chm_done = .T.$ ENDIF @ 22,19 SAY "Auto Chemical Match Complete. Press any key to continue ..." WAIT "" CLOSE ALL RETURN && auto PROCEDURE report \* Displays static portion of chemical matching status report on screen. @ 1,0 CLEAR @ 4,26 SAY "NPDES DATA CONVERSION" @ 5,21 SAY "CHEMICAL MATCHING STATUS REPORT" @ 8,22 SAY "Number of records processed:" @ 8,52 SAY processed PICTURE "999999" @ 8,58 SAY "of" @ 8,61 SAY total\_recs PICTURE "999999" @ 10,24 SAY "Number of records matched:" @ 10,52 SAY match PICTURE \*99999\* @ 12,20 SAY "Number of records not matched:" @ 12,52 SAY no\_match PICTURE "999999" **RETURN && report** PROCEDURE says \* Displays dynamic portion of chemical matching status report (counter \* variables). @ 8,52 SAY processed PICTURE \*99999\* @ 10,52 SAY match PICTURE \*999999\* @ 12,52 SAY no\_match PICTURE \*99999\* **RETURN && says** 

PROCEDURE mis • User checks unmatched chemicals and either assigns a correct chemical ID \* from a printed list. \* Open working Concentration file. The index is conditional on the unmatched \* flag being true, so only unmatched chemicals are shown. USE npdes\nconcext IN 1 ORDER param\_c ALIAS conc \* Display instructions for resolving unmatched chemicals. CLEAR TYPE npdes\nvchem1.msg WAIT . \* BROWSE the unmatched chemicals; the user is not allowed to add records. \* The user will change the unmatched flag to False as each chemicla is \* resolved by assigning a chemical ID. Changing the unmatched flag to False \* removes a record from the conditional index. When all records have been \* removed from the index in this way, an error condition results; the ON ERROR \* CLEAR line clears the screen when this happens, effectively ending the \* BROWSE. The user may also exit the Browse with Esc or Ctrl-End before all \* unmatched chemicals have been dealt with. ON ERROR CLEAR BROWSE NOAPPEND NOFOLLOW FIELDS chem\_mis, chem\_id, param\_code /r, param\_name /r ON ERROR \* Restore header wiped out by BROWSE. CLEAR @ 0,1 SAY "WRMS DATA VERIFICATION CONVERSION UTILI Т Ү" \* See if any unmatched chemicals remain (user may have exited before all were \* matched). LOCATE FOR chem mis \* If no unmatched chemicals are found, then chemical matching is done. IF .not. found()  $1\_chm\_done = .T.$ \* Copy STORET numbers for manually matched chemicals to the Chemical file. @ 1,0 CLEAR @ 12,12 SAY "Copying STORET numbers to Chemical file..." SET ORDER TO USE chemical IN 2 ORDER chem\_id ALIAS chem SET RELATION TO chem\_id INTO chem SET TALK ON REPLACE chem->storet\_num WITH param\_code FOR chem->storet\_num = space(5) SET TALK OFF @ 12,54 SAY "Done" ENDIF CLOSE ALL RETURN && mis \* EOF: nvchem.prg

nvchem1.msg

WRMS DATA VERIFICATION CONVERSION UTILITY

Directions for resolution of unmatched chemical parameters:

- \* Unmatched NPDES records will be displayed, indexed on param\_code.
- \* You will only be able to edit the chem\_id and chem\_mis fields.
- \* Look up the correct chem\_id in the alphabetical listing of the WRMS chemical file and enter it in the chem\_id field. Change the chem\_mis field to F (False).
- \* If a parameter is not in the WRMS chemical file, change chem\_mis to F for all records containing that chemical, but do not enter a chem\_id.
- \* Press <Ctrl-End> when finished.

Press any key to continue ...

\* \* Filename: NVEMISS.PRG \* Author: Anne E. Crook, Research Triangle Institute 2/15/93 Last updated: 3/2/93 \* Created: \* Called by:VCONVERT \* Comments: Converts emissions data from NPDES to WRMS format. Assumes all facilities and chemicals have been correctly matched. Creates 1 Emission record from 1 or more working Concentration records. \* Open data base files: working Concentration and Flow files and units \* conversionfactor file. Concentration and Flow files are linked on NPDES \* number and pipe; Concentration and units conversion files are linked on \* parameter units code. USE npdes\nconcext IN 1 ALIAS conc USE npdes\nflowext IN 2 ORDER pipe ALIAS flow USE npdes\unitcode IN 3 ORDER unit\_code ALIAS units SET RELATION TO npdes + pipe\_num INTO flow SELECT flow SET RELATION TO conc->param\_unit INTO units SELECT conc \* Calculate annual emissions from concentration and flow using conversion \* factors in Unitcode file. The flowfactor field is 1 if flow needs to be \* included in the calculation (e.g., if avg is in concentration units like \* mg/l) and 0 if it does not (e.g., if avg is already in emission units, like \* lbs/day). The calculated emission is placed in the emiss field of the \* working Concentration file. @ 1,0 CLEAR @ 9,19 SAY "Calculating emissions ......" SET TALK ON REPLACE ALL emiss WITH conc->avg \* iif(units->flowfactor=1,flow->avg,1) \*; units->conversion REPLACE flow\_na WITH .T. FOR units->flowfactor = 1 .AND. flow->avg = 0 SET TALK OFF @ 10,0 CLEAR @ 9,56 SAY "Done" \* Close files and reopen working Concentration file. CLOSE ALL USE npdes\nconcext IN 1 ALIAS conc \* Data base in use is working Concentration file. This file may contain \* multiple records for a specific NPDES number, pipe, and parameter if data \* were originally reported in two different units. These have now been \* converted to emissions with common units, and need to be combined. \* Create a new data base file with one blank record per pipe and parameter \* to hold emissions averaged across different units. @ 11,19 SAY "Creating intermediate file ....." SET ORDER TO pipeparamu && order = npdes+pipe\_num+param\_code+param\_name UNIQUE SET TALK ON COPY TO npdes\nconcavg SET TALK OFF @ 12,0 CLEAR @ 11,56 SAY "Done" USE npdes\nconcavg IN 1 ALIAS avg BLANK ALL FIELDS param\_unit, avg, emiss, weight

```
USE npdes\nconcext IN 2 ORDER pipeparam ALIAS conc
@ 12,19 SAY "Averaging across units ....."
loopavg = 0
SCAN && avg
  loopavg = loopavg + 1
  @ 12,56 SAY loopavg PICTURE "9999"
  sum_emiss = 0
  sum_weight = 0
  missing_wt = .F.
  mflow_na = .F.
  SELECT conc
  SET KEY TO avg->npdes + avg->pipe_num + avg->param_code + avg->param_name
  COUNT TO num_conc
  GO TOP
  IF num_conc = 1
                           && only one conc record; averaging not needed
    SELECT avg
    REPLACE emiss WITH conc->emiss,;
            flow_na WITH conc->flow_na
  ELSE
                           && more than one conc records to be averaged
    IF num_conc > 1
      SCAN && conc
        IF weight = 0
                             && weight missing
          missing_wt = .T.
          EXIT
        ELSE
                             && weight not missing
          sum_emiss = sum_emiss + emiss * weight
          sum_weight = sum_weight + weight
          mflow_na = (mflow_na .OR. flow_na)
        ENDIF
      ENDSCAN && conc
      SELECT avg
      IF .not. missing_wt
        REPLACE emiss WITH sum_emiss/sum_weight,;
                flow_na WITH mflow_na
      ELSE
        REPLACE weight_na WITH .T.
      ENDIF
    ENDIF
  ENDIF
ENDSCAN && avg
@ 12,56 SAY *Done*
CLOSE ALL
* Aggregate emissions to NPDES number level (across pipes) and enter in
* Emission file.
* Initialize variables.
                     && Count number of average concentration records processed.
processed = 0
added = 0
                     && Count number of Emission records created.
* Open files: average concentration file and Emission file. Files are unlinked.
* Average concentration file is selected.
USE npdes\nconcavg IN 1 ALIAS avg
USE vemiss IN 2 ORDER source_yr ALIAS em
                                             && order = npdes+param_code+year
* Display static portions of emission conversion status report
@ 1,0 CLEAR
@ 4,26 SAY "NPDES DATA CONVERSION"
@ 5,22 SAY "EMISSION UPDATE STATUS REPORT"
@ 9,19 SAY "Number of records processed:
@ 9,48 SAY " 0"
                                               of"
```

@ 9,56 SAY reccount("avg") PICTURE "9999" @ 11,19 SAY "Number of EMISSION records created:" @ 11,56 SAY \* 0\* \* Process all average concentration file records. SCAN \* See if an Emission record already exists for the current NPDES number and \* parameter. SELECT em SEEK avg->npdes + avg->param\_code + str(avg->year,4) \* If no Emission record is found, create one and fill it in from the average \* concentration record. IF .NOT. found() APPEND BLANK REPLACE em->source WITH "N",; em->source\_id WITH avg->npdes,; em->fac\_id WITH avg->fac\_id,; em->chem\_id WITH avg->chem\_id,; em->year WITH avg->year,; em->water\_amt WITH avg->emiss,; em->flow\_na WITH avg->flow\_na,; em->weight\_na WITH avg->weight\_na added = added + 1@ 11,56 SAY added PICTURE "9999" \* If an Emission record is found, add the emission amount from the average concentration file to the quantity already in the water amount field in the \* Emission file. ELSE REPLACE em->water\_amt WITH em->water\_amt + avg->emiss,; em->flow\_na WITH (em->flow\_na .OR. avg->flow\_na),; em->weight\_na WITH (em->weight\_na .OR. avg->weight\_na) ENDIF \* Reselect the average concentration file for correct functioning of SCAN loop. SELECT avg processed = processed + 1 @ 9,48 SAY processed PICTURE \*9999\* ENDSCAN \* All average concentration records have been processed. Update status variable \* to indicate emission conversion is complete.  $1_{emiss} = .T.$ CLOSE ALL \* Delete working files. Files containing bad concentration and flow data are \* not deleted, but saved for possible future use. @ 1,0 CLEAR @ 9,12 SAY "Deleting working files ..." DELETE FILE npdes\npdesint.dbf && intermediate working file DELETE FILE npdes\nconcext.dbf && working Concentration file DELETE FILE npdes\nconcext.mdx && working Concentration file index DELETE FILE npdes\nflowext.dbf && working Flow file DELETE FILE npdes\nflowext.mdx && working Flow file index DELETE FILE npdes\nconcavg.dbf && average concentration file && average concentration file index DELETE FILE npdes\nconcavg.mdx && Saved status variables DELETE FILE npdes\nvstatus.mem

@ 9,39 SAY "Done"

\* Display message that data conversion is complete.

@ 1,0 CLEAR @ 9,25 TO 13,53 DOUBLE @ 11,27 SAY "Data conversion complete." @ 22,50 SAY "Press any key to continue ..." WAIT ""

RETURN && nvemiss

•

\* EOF: nvemiss.prg

\* Filename: PVWORK.PRG \* Author: Anne E. Crook, Research Triangle Institute \* Created: 2/15/93 Last updated: 2/23/93 \* Called by:VCONVERT \* Comments: Creates working Pretreatment files for conversion program, with structure extended to include fields used in processing. Also cleans up some fields. \* Initialize status variables.  $1_fac_done = .F.$ && Facility matching done  $1\_chm\_done = .F.$ && Chemical matching done 1 emiss = .F.&& Emissions conversion done \* Inititalize other variable. myear = space(2)&& Year to be converted \* Ask for year being converted. @ 1,0 CLEAR SET CONFIRM ON @ 9,20 SAY "Enter year of data being converted: 19" @ 9,58 GET myear PICTURE \*99\* READ SET CONFIRM OFF \* Create working data file PREEXT.DBF from saved structure PRESTRU.DBF @ 1,0 CLEAR @ 8,11 SAY "Creating working file ....." CREATE pretreat\preext FROM pretreat\prevstru \* Create index on city and facility name. This index will be used manually \* to match pretreatment facilities to TRIS facilities. INDEX ON city + fac\_name TAG city\_fac \* Create unique index on pretreatment ID (POTW NPDES number plus IU number). \* This index will be used to create new VFAC records. INDEX ON potw\_npdes + iu\_num TAG potw\_iu\_u UNIQUE \* Add records from the original data file to the working data file. SET TALK ON APPEND FROM pretreat\pretreat.dat TYPE SDF SET TALK OFF @ 9,0 CLEAR CLOSE DATABASE @ 8,45 SAY "Done" \* Enter year being converted to working file. @ 10,11 SAY "Entering year in working file ..." USE pretreat\preext REPLACE ALL year WITH val("19" + myear) @ 10,45 SAY \*Done\* CLOSE ALL RETURN && pvwork.prg \* EOF: pvwork.prg

\*

\* \* Filename: PVFAC.PRG \* \* Author: Anne E. Crook, Research Triangle Institute \* \* Created: 2/15/93 Last updated: 2/24/93 \* Called by:VCONVERT \* Comments: Pretreatment data facility matching routine. Pretreatment facilities are matched to TRIS facilities in WRMS based on the Pretreatment ID. \* Open working file and WRMS Facility file. Files are linked on pretreatment \* ID (POTW NPDES number plus facility IU number). USE pretreat\preext IN 1 ALIAS pre USE facility IN 2 ORDER pre\_id ALIAS fac SET RELATION TO potw\_npdes + iu\_num INTO fac \* Copy facility ID from Facility to working pretreatment data file for matching \* Pretreatment IDs. @ 1,0 CLEAR @ 9,16 SAY "Assigning FAC\_IDs to matched facilities ..." SET TALK ON REPLACE pre->fac\_id WITH fac->fac\_id FOR fac->fac\_id <> space(10); .AND. potw\_npdes <> space(9) SET TALK OFF @ 10,0 CLEAR @ 9,60 SAY "Done" \* Close files and reopen working Pretreatment file. CLOSE ALL USE pretreat\preext IN 1 ALIAS pre \* Delete unmatched nonmanufacturing facilities from working pretreatment file. @ 1,0 CLEAR @ 10,16 SAY "Deleting nonmanufacturing facilities ....." SET TALK ON DELETE FOR fac\_id = space(10) .AND.; (val(left(sic,2)) < 20 .OR. val(left(sic,2)) > 39) PACK SET TALK OFF @ 11,0 CLEAR @ 10,60 SAY "Done" CLOSE ALL USE pretreat\preext IN 1 ORDER potw\_iu\_u ALIAS pre USE facility IN 2 ORDER fac\_id ALIAS fac SET RELATION TO fac\_id INTO fac USE vfac IN 3 ORDER pre\_id ALIAS vfac \* Copy facility information for unmatched manufacturing facilities to VFAC. @ 11,16 SAY "Creating new facility records ....." newvfacs = 0SCAN && preext SELECT vfac SEEK pre->potw\_npdes + pre->iu\_num IF .NOT. found() newvfacs = newvfacs + 1 APPEND BLANK REPLACE fac\_name WITH pre->fac\_name,;

```
street_1 WITH pre->address,;
            city WITH pre->city,;
            zip WITH pre->zip,;
            sic_1 WITH pre->sic,;
            pre_id WITH pre->potw_npdes + pre->iu_num,;
            year_added WITH pre->year
    IF pre->fac_id <> space(10)
      REPLACE county WITH fac->county,;
              contact WITH fac->contact,;
              phone WITH fac->phone
    ENDIF
    @ 11,60 SAY newvfacs PICTURE *999*
  ENDIF
  SELECT pre
ENDSCAN
@ 11,60 SAY "Done"
CLOSE ALL
* Update status variable to indicate facility matching is complete.
l_fac_done = .T.
@ 22,24 SAY "Facility Match Complete. Press any key to continue ..."
WAIT **
RETURN && pvfac.prg
* EOF: pvfac.prg
```

\*\*\*\*\*\* \* Filename: PVCHEM.PRG \* Author: Anne E. Crook, Research Triangle Institute \* Created: 2/15/93 Last updated: \* Called by:VCONVERT \* Comments: Chemical matching is not necessary with the Pretreatment data; each of the chemicals is reported in a separate field. Chemical IDs are assigned in the emission conversion program module. \* Display message indicating that chemical matching is unnecessary. @ 1,0 CLEAR @ 9,21 TO 14,52 DOUBLE @ 11,23 SAY "It is not necessary to Match" @ 12,23 SAY "Chemicals with these data." @ 22,50 SAY "Press any key to continue ..." WAIT \*\* \* Update status variable to indicate chemical matching is done. l\_chm\_done = .T. RETURN && pvchem.prg \* EOF: pvchem.prg

\* Filename: PVEMISS.PRG \* Author: Anne E. Crook, Research Triangle Institute 2/15/93 Last updated: 3/2/93 Created: \* Called by:VCONVERT \* Comments: Converts emissions data from Pretreatment to WRMS format. Assumes all facilities have been correctly matched. Creates up to 20 Emission records from 1 working pretreatment record (one per chemical). \* Open data bases: Vemiss and working pretreatment data file. Working \* pretreatment data file is selected. USE vemiss IN 1 ALIAS em USE pretreat\preext IN 2 ALIAS pre SELECT pre \* Initialize variables. && Count Emission records added  $added_em = 0$ \* Display static portions of emission conversion status report @ 1,0 CLEAR @ 4,21 SAY "PRETREATMENT DATA CONVERSION" @ 5,21 SAY "EMISSION UPDATE STATUS REPORT" @ 9,15 SAY "Number of records processed: of" @ 9,47 SAY "0" @ 9,52 SAY reccount("pre") PICTURE "9999" @ 11,15 SAY "Number of EMISSION records added:" @ 11,52 SAY "0" \* Process each pretreatment record. SCAN && pre \* Add records to Emission and fill in data from working pretreatment file, \* using flow and concentration to calculate emissions in pounds per year. SELECT em \* Create Emission records only if the working pretreatment file has a nonzero concentration value. \* Phenol IF pre->mphe > 0APPEND BLANK REPLACE em->source WITH "P",; em->source\_id WITH pre->potw\_npdes+pre->iu\_num,; em->fac\_id WITH pre->fac\_id,; em->year WITH pre->year,; em->chem\_id WITH "CPHEN00808",; em->potw\_amt WITH pre->mflow \* pre->mphe \* 3.05,; em->flow\_na WITH (pre->mflow = 0) added\_em = added\_em + 1 @ 11,49 SAY added\_em PICTURE "9999" ENDIF \* Silver IF pre->mag > 0 APPEND BLANK REPLACE em->source WITH "P",; em->source\_id WITH pre->potw\_npdes+pre->iu\_num,;

```
em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CSILV00903",;
            em->potw_amt WITH pre->mflow * pre->mag * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
   added em = added_em + 1
   @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
* Aluminum
  IF pre->mal > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,
            em->chem_id WITH "CALUM00034",;
            em->potw_amt WITH pre->mflow • pre->mal * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
   added_em = added_em + 1
   @ 11,49 SAY added_em PICTURE "9999"
  ENDIF
* Arsenic
  IF pre->mas > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CARSE00106",;
            em->potw_amt WITH pre->mflow * pre->mas * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
   added_em = added_em + 1
   @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
* Cadmium
  IF pre->mcd > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CCADM00255",;
            em->potw_amt WITH pre->mflow * pre->mcd * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
   added_em = added_em + 1
   @ 11,49 SAY added_em PICTURE *9999*
 ENDIF
* Cobalt
 IF pre->mco > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CCOBA00335",;
            em->potw_amt WITH pre->mflow * pre->mco * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
   added_em = added_em + 1
   @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
```

```
* Chromium, total
 IF pre->mcr_tot > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CCHRO00332",;
            em->potw_amt WITH pre->mflow * pre->mcr_tot * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE *9999*
 ENDIF
* Copper
  IF pre->mcu > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CCOPP00344",;
em->potw_amt WITH pre->mflow * pre->mcu * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added em PICTURE "9999"
 ENDIF
* Mercury
  IF pre->mhg > 0
    APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CMERC00681",;
            em->potw_amt WITH pre->mflow * pre->mhg * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
* Magnesium - not a TRI chemical
   IF pre->mmg > 0
     APPEND BLANK
     REPLACE em->source WITH "P",;
             em->source_id WITH pre->potw_npdes+pre->iu_num,;
             em->fac_id WITH pre->fac_id,;
             em->year WITH pre->year,;
             em->chem_id WITH "CMAGNXXXXX",;
             em->potw_amt WITH pre->mflow * pre->mmg * 3.05,;
             em->flow_na WITH (pre->mflow = 0)
     added_em = added_em + 1
     @ 11,49 SAY added_em PICTURE *9999*
   ENDIF
* Manganese
  IF pre->mmn > 0
    APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
```

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```
em->chem_id WITH *CMANG00666*,;
            em->potw_amt WITH pre->mflow * pre->mmn • 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
   @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
* Nickel
  IF pre->mni > 0
    APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CNICK00749",;
            em->potw_amt WITH pre->mflow * pre->mni * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
   @ 11,49 SAY added_em PICTURE *9999*
  ENDIF
* Lead
  IF pre->mpb > 0
    APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CLEAD00637",;
            em->potw_amt WITH pre->mflow * pre->mpb * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE "9999"
  ENDIF
* Selenium
  IF pre->mse > 0
    APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CSELE00895",;
            em->potw_amt WITH pre->mflow * pre->mse • 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE *9999*
  ENDIF
* Zinc
  IF pre->mzn > 0
    APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CZINC01044",;
            em->potw_amt WITH pre->mflow * pre->mzn * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE "9999"
  ENDIF
* Toluene
```

```
IF pre->mtol > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CTOLU00154",;
            em->potw_amt WITH pre->mflow * pre->mtol * 3.05,;
            em \rightarrow flow na WITH (pre\rightarrow mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
* Barium
 IF pre->mba > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CBARI00123",;
            em->potw_amt WITH pre->mflow * pre->mba * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
   added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE *9999*
 ENDIF .
* Antimony
  IF pre->msb > 0
   APPEND BLANK
   REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CANTI00088",;
            em->potw_amt WITH pre->mflow * pre->msb * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE *9999*
 ENDIF
* Formaldehyde
  IF pre->mchoh > 0
   APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CFORM00575",;
            em->potw_amt WITH pre->mflow * pre->mchoh * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE "9999"
 ENDIF
* Chlorine
  IF pre->mchlor > 0
    APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CCHL000298",;
            em->potw_amt WITH pre->mflow * pre->mchlor * 3.05,;
```

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```
em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE *9999*
  ENDIF
* Cyanide
  IF pre->mcn > 0
    APPEND BLANK
    REPLACE em->source WITH "P",;
            em->source_id WITH pre->potw_npdes+pre->iu_num,;
            em->fac_id WITH pre->fac_id,;
            em->year WITH pre->year,;
            em->chem_id WITH "CCYAN00366",;
            em->potw_amt WITH pre->mflow * pre->mcn * 3.05,;
            em->flow_na WITH (pre->mflow = 0)
    added_em = added_em + 1
    @ 11,49 SAY added_em PICTURE *9999*
  ENDIF
* Update status report.
  @ 9,44 SAY recno("pre") PICTURE "9999"
  SELECT pre
ENDSCAN && pre
* Update status variable to indicate emission conversion is done.
1_{emiss} = .T.
CLOSE ALL
* Delete working files.
@ 1,0 CLEAR
@ 10,12 SAY "Deleting working files ..."
                                     && working data file
DELETE FILE pretreat\preext.dbf
DELETE FILE pretreat\preext.mdx
                                        && working data file index
DELETE FILE pretreat\pvstatus.mem
                                        && saved status variables
@ 10,39 SAY "Done"
@ 1,0 CLEAR
@ 9,25 TO 13,53 DOUBLE
@ 11,27 SAY "Data conversion complete."
@ 22,50 SAY "Press any key to continue ..."
WAIT ""
RETURN && pvemiss
```

\* EOF: pvemiss.prg

## APPENDIX B PROGRAM CODE FOR TASK6.PRG

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```
* File name: task6.prg
* Author: Anne Crook, Research Triangle Institute
* Created: 2/26/93 Updated: 4/14/93
* Make memory variables accessible to all procedures.
PUBLIC cyear, nyear, data, avail, ord_field, em_field, vem_field, rnge_field
* Initialize environment and variables
DO initial
* Get year and type of data to compare
DO getdata
* Check availability of data selected.
DO checkdata
* If data are available, generate discrepancy reports.
IF avail
 DO xdata
 DO discrep
ENDIF
* Reset environment.
CLOSE ALL
SET TITLE ON
SET SAFETY ON
SET TALK ON
RETURN && t6.prg
PROC initial
* Initialize environment and variables
 SET TITLE OFF
 SET SAFETY OFF
 SET TALK OFF
 cyear = space(2)
 nyear = 0
 data = space(1)
RETURN
PROC getdata
* Get year and type of data to compare, then set assorted variables according
* to the choice.
* Any variables set here must be in the PUBLIC statement at the beginning of
* the program!
 @ 0,0 CLEAR
 SET CONFIRM ON
 @ 8,20 SAY "Enter data to compare:"
  @ 9,20 SAY "A=AIRS, N=NPDES, P=Pretreatment"
 @ 8,43 GET data PICTURE "@M N,P,A"
 @ 11,20 SAY "Enter year to compare: 19"
 @ 11,45 GET cyear PICTURE *99*
 READ
 SET CONFIRM OFF
 nyear = 1900 + val(cyear)
 DO CASE
```

```
CASE data = "A"
     ord_field = "air_id"
     em_field = "stack_em"
     vem_field = "stack_vem"
     rnge_field = "stack_rnge"
   CASE data = "N"
     ord_field = "npdes"
     em_field = "water_em"
     vem_field = "water_vem"
     rnge_field = "water_rnge"
   CASE data = "P"
     ord_field = "pre_id"
     em_field = "potw_em"
     vem_field = "potw_vem"
     rnge_field = "potw_rnge"
  ENDCASE
RETURN
PROC checkdata
* Check availability of data selected.
 USE vemiss
 LOCATE FOR source = data .AND. year = nyear
  avail = found()
  IF .not. avail
   @ 0,0 CLEAR
   @ 7,18 TO 11,54 DOUBLE
   @ 9.20 SAY "No data available for comparison."
   @ 21,50 SAY "Press any key to continue..."
   WAIT ..
 ENDIF
 CLOSE ALL
RETURN
PROC xdata
* Extract data from Emission and Vemiss needed for discrepancy checks
* Extract TRI records from Emission for appropriate year.
  TEXT
 Extracting TRI records
 ENDTEXT
 USE emission
 COPY TO t6&data.1 FOR source = "T" .AND. year = nyear;
                 FIELDS fac_id, chem_id, year, source_id
* Extract single-media records from Vemiss for the appropriate year.
 TEXT
 Extracting single-media records
 ENDTEXT
 USE vemiss
 COPY TO t6&data.2 FOR source = data .AND. year = nyear;
                 FIELDS fac_id, chem_id, year, source_id
 CLOSE ALL
RETURN
PROC discrep
* Generate discrepancy report.
```

TEXT Combining TRI and single-media data ENDTEXT USE t6&data.1 APPEND FROM t6&data.2 FOR fac\_id <> space(10) INDEX ON fac\_id + chem\_id TAG fac\_chm\_u UNIQUE COPY TO t6&data.3 USE t6&data.3 BLANK ALL FIELDS source\_id USE t6&data.2 DELETE FOR fac\_id <> space(10) PACK INDEX ON source\_id + chem\_id TAG fac\_chm\_u UNIQUE COPY TO t6&data.4 \* Create a file with additional fields. CREATE t6&data.5 FROM t6stru APPEND FROM t6&data.3 APPEND FROM t6&data.4 \* T6&data.5 should now have one record per facility and chemical reported in either TRI or single media data base. \* Sum emissions data for each facility and chemical reported to TRI to the summary file. (Usually there is only one emission record per facility and \* chemical, but occasionally two records are reported; these are summed here.) TEXT Summing TRI emissions ENDTEXT USE t6&data.5 IN 1 ALIAS t6 USE emission IN 2 ORDER fac\_chm\_yr ALIAS em SET RELATION TO fac\_id + chem\_id + str(year,4) INTO em SET SKIP TO em REPLACE stack\_em WITH stack\_em + em->stack\_amt,; water\_em WITH water\_em + em->water\_amt,; potw\_em WITH potw\_em + em->potw\_amt,; stack\_rnge WITH (stack\_rnge .OR. em->stack\_rnge),; water\_rnge WITH (water\_rnge .OR. em->water\_rnge),; potw\_rnge WITH (potw\_rnge .OR. em->potw\_rnge); FOR em->fac\_id <> space(10) \* Sum emissions data for each facility and chemical reported to the singlemedia data base to the summary file. TEXT Summing single-media emissions ENDTEXT \* For facilities with FAC\_IDs USE vemiss IN 2 ORDER fac\_chm\_yr ALIAS vem SET RELATION TO fac\_id + chem\_id + str(year,4) INTO vem SET FILTER TO vem->source = data REPLACE stack\_vem WITH vem->stack\_amt, water\_vem WITH vem->water\_amt,; potw\_vem WITH vem->potw\_amt, flow\_na WITH vem->flow\_na,; weight\_na WITH vem->weight\_na; FOR vem->fac\_id <> space(10) \* For facilities without FAC\_IDs USE vemiss IN 2 ORDER src\_chm\_yr ALIAS vem SET RELATION TO source\_id + chem\_id + str(year, 4) INTO vem SET FILTER TO vem->source = data

REPLACE stack\_vem WITH vem->stack\_amt, water\_vem WITH vem->water\_amt,; potw\_vem WITH vem->potw\_amt, flow\_na WITH vem->flow\_na,; weight\_na WITH vem->weight\_na; FOR vem->source\_id <> space(18) CLOSE ALL \* Delete records for which data are missing for single media or reported as a \* range in TRI TEXT Deleting records with missing data or range data ENDTEXT USE t6&data.5 IN 1 ALIAS t6 DELETE FOR flow\_na .OR. weight\_na .OR. &rnge\_field PACK CLOSE ALL \* Calculate the difference in quantity reported to TRI and single-media data \* base. TEXT Calculating discrepancies ENDTEXT USE.t6&data.5 IN 1 ALIAS t6 REPLACE ALL diff\_amt WITH abs(&em\_field - &vem\_field) \* Delete records with discrepancies less than 1,000 lbs TEXT Deleting records below cutoff (1,000 lbs for NPDES and AIRS/NEDS; 10,000 lbs for Pretreatment) ENDTEXT IF data = "P" DELETE FOR diff\_amt < 10000 ELSE DELETE FOR diff\_amt < 1000 ENDIF PACK \* The total difference amount is calculated for each facility. Note that the \* difference amounts are not signed, so that equal but opposite discrepancies \* will not cancel out. TEXT Summing discrepancies to facility level ENDTEXT GO TOP  $first_rec = 1$ DO WHILE .NOT. EOF() mfac = fac\_id msrc = source\_id SUM diff\_amt TO msum WHILE fac\_id = mfac .AND. source\_id = msrc GO first\_rec REPLACE total\_diff WITH msum WHILE fac\_id = mfac .AND. source\_id = msrc first\_rec = recno() ENDDO INDEX ON total\_diff TAG total\_diff DESCENDING \* The summary file now contains only facilities with significant discrepancies.

\* Link it to the facility, and chemical files. Facilities are ordered in \* descending order of total difference in reported quantity.

TEXT Linking files ENDTEXT SET EXACT ON USE t6&data.5 IN 1 ORDER total\_diff ALIAS t6 USE facility IN 2 ORDER fac\_id ALIAS fac USE vfac IN 3 ORDER &ord\_field ALIAS vfac USE chemical IN 4 ORDER chem\_id ALIAS chem SELECT 1 SET RELATION TO t6->fac\_id INTO fac SELECT 2 SET RELATION TO t6->source\_id INTO vfac SELECT 3 SET RELATION TO t6->chem\_id INTO chem SELECT 1 GO TOP SET FIELDS TO t6->fac\_id, t6->chem\_id, t6->year, t6->source\_id,; t6->stack\_em, t6->stack\_vem, t6->water\_em, t6->water\_vem,; t6->potw\_em, t6->potw\_vem, t6->diff\_amt,; fac->fac\_name, fac->street\_1, fac->street\_2, fac->city,; fac->county, fac->zip, fac->contact, fac->phone, fac->sic\_1,; fac->air\_id, fac->npdes, fac->pre\_id,; vfac->fac\_name, vfac->street\_1, vfac->city, vfac->county,; vfac->zip, vfac->contact, vfac->phone, vfac->sic\_1,; vfac->air\_id, vfac->npdes, vfac->pre\_id,; chem->cas\_num, chem->use\_name, chem->storet\_num \* Generate report. TEXT Generating report ENDTEXT  $_{plength} = 67$ REPORT FORM t6&data TO FILE t6&data.&cyear..txt CLOSE ALL DELETE FILE t6&data.1.dbf DELETE FILE t6&data.1.mdx DELETE FILE t6&data.2.dbf DELETE FILE t6&data.2.mdx DELETE FILE t6&data.3.dbf DELETE FILE t6&data.4.dbf DELETE FILE t6&data.5.dbf DELETE FILE t6&data.5.mdx RETURN && discrep

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APPENDIX C PROGRAM CODE FOR TASK8.PRG

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```
* File name: task8.prg
* Author: Anne Crook, Research Triangle Institute
* Created: 3/16/93 Updated: 4/14/93
* List facilities reporting to single-media data bases over the TRI reporting
* threshold but not reporting to TRI.
* Make memory variables accessible to all procedures.
PUBLIC cyear, nyear, data, avail, amt_field, ord_field, em_field, vem_field,;
      rnge_field, manuf
* Initialize environment and variables
DO initial
* Get year and type of data to compare
DO getdata
* Check availability of data selected.
DO checkdata
* If data are available, generate discrepancy reports.
IF avail.
 DO discrep
ENDIF
* Reset environment.
SET TITLE ON
SET SAFETY ON
SET TALK ON
RETURN && t8.prg
PROC initial
* Initialize environment and variables
 SET TITLE OFF
 SET SAFETY OFF
 SET TALK OFF
 cyear = space(2)
 nyear = 0
 data = space(1)
RETURN
PROC getdata
* Get year and type of data to compare, then set assorted variables according
* to the choice.
* Any variables set here must be in the PUBLIC statement at the beginning of
* the program!
 @ 0,0 CLEAR
 SET CONFIRM ON
 @ 8,20 SAY "Enter data to compare:"
 @ 9,20 SAY "A=AIRS, N=NPDES, P=Pretreatment"
 @ 8,43 GET data PICTURE "@M N,P,A"
 @ 11,20 SAY "Enter year to compare: 19"
 @ 11,45 GET cyear PICTURE "99"
 READ
 SET CONFIRM OFF
```

```
nyear = 1900 + val(cyear)
 DO CASE
   CASE data = "A"
     amt_field = "stack_amt"
     ord_field = "air_id"
   CASE data = "N"
     amt_field = "water_amt"
     ord_field = "npdes"
   CASE data = "P"
     amt_field = "potw_amt"
ord_field = "pre_id"
 ENDCASE
 DO CASE
   CASE nyear = 1987
     manuf = 75000
   CASE nyear = 1988
     manuf = 50000
   OTHERWISE
     manuf = 25000
 ENDCASE
RETURN
PROC checkdata
* Check availability of data selected.
 USE vemiss
 LOCATE FOR source = data .AND. year = nyear
 avail = found()
 IF .not. avail
   @ 0,0 CLEAR
   @ 7,18 TO 11,54 DOUBLE
   @ 9,20 SAY "No data available for comparison."
   @ 21,50 SAY "Press any key to continue..."
   WAIT ""
 ENDIF
 CLOSE ALL
RETURN
PROC discrep
* Generate discrepancy report.
* Extract single-media records from Vemiss for the appropriate year.
 TEXT
 Extracting single-media records
 ENDTEXT
 USE vemiss IN 1 ALIAS vemiss
 COPY TO t8&data.1 FOR source = data .AND. year = nyear .AND.;
                     &amt_field >= 10000;
                  FIELDS source_id, year, fac_id, chem_id
* Delete records that are found in Emission (i.e., that also reported to TRI).
 TEXT
 Deleting facilities reporting to TRI
 ENDTEXT
 USE t8&data.1 IN 1 ALIAS t8
 USE emission IN 2 ORDER fac_chm_yr ALIAS em
 SET RELATION TO fac_id + chem_id + str(year, 4) INTO em
 DELETE FOR em->fac_id <> space(10)
 PACK
```

CLOSE ALL

```
* This file now indicates the chemicals reported over 10,000 lbs and not
* reported in TRI.
* Create a file with additional fields
 CREATE t8&data.2 FROM t8stru
 APPEND FROM t8&data.1
* Copy emissions from Vemiss
 USE vemiss IN 2 ORDER src_chm_yr ALIAS vem
 SET RELATION TO source_id + chem_id + str(year, 4) INTO vem
 REPLACE ALL rel_amt WITH vem->&amt_field
 CLOSE ALL
 TEXT
 Summing releases to facility level
 ENDTEXT
 USE t8&data.2 IN 1 ALIAS t8
 GO TOP
 first_rec = 1
 DO WHILE .NOT. EOF()
   msrc = source_id
   SUM rel_amt TO msum WHILE source_id = msrc
   GO first_rec
   mthresh = "Otherwise Used"
   SCAN WHILE source_id = msrc
     mthresh = iif(rel_amt > manuf, "Processed or Manufactured", mthresh)
   ENDSCAN
   GO first_rec
   REPLACE total_amt WITH msum, thresh_ex WITH mthresh WHILE source_id = msrc
   first_rec = recno()
 ENDDO
  INDEX ON thresh_ex + str(total_amt,11) TAG thresh_amt DESCENDING
 CLOSE ALL
* Link to the Vfac and Chemical files.
* Facilities are ordered in descending order of total quantity reported.
 TEXT
 Linking files
 ENDTEXT
 SET EXACT ON
 USE t8&data.2 IN 1 ORDER thresh_amt ALIAS t8
 USE vfac IN 2 ORDER &ord_field ALIAS vfac
 USE chemical IN 3 ORDER chem_id ALIAS chem
 SELECT 1
 SET RELATION TO t8->source_id INTO vfac
  SELECT 2
 SET RELATION TO t8->chem_id INTO chem
  SELECT 1
 GO TOP
  SET FIELDS TO t8->source_id, t8->year, t8->chem_id, t8->fac_id, t8->rel_amt,;
                t8->thresh_ex,;
                vfac->fac_name, vfac->street_1, vfac->city, vfac->county,;
                vfac->zip, vfac->contact, vfac->phone, vfac->sic_1,;
                vfac->air_id, vfac->npdes, vfac->pre_id, chem->cas_num,;
                chem->use_name, chem->storet_num
```

\* Generate report.

TEXT Generating report

ENDTEXT

\_plength = 53 REPORT FORM t8&data TO FILE t8&data.&cyear..txt CLOSE ALL DELETE FILE t8&data.1.dbf DELETE FILE t8&data.1.mdx DELETE FILE t8&data.2.dbf DELETE FILE t8&data.2.mdx

RETURN && discrep

.

## APPENDIX D PROGRAM CODE FOR TASK7.PRG

•
```
* File name: task7.prg
* Author: Anne Crook, Research Triangle Institute
* Created: 4/15/93 Updated: 5/28/93
* Make memory variables accessible to all procedures.
PUBLIC mfirstyear, mlastyear, numyears
DO initial
DO getyear
DO makefile
DO fillfile
IF numyears > 2
 DO pattern
 DO patt_rpt
ENDIF
DO stop
D0 stop_rpt
DO limitrecs
DO calc_discr
DO sum_discrp
DO discrp_rpt
DO reset
RETURN && t7.prg
PROC initial

    Initialize environment and variables

 SET TITLE OFF
 SET SAFETY OFF
 SET TALK OFF
RETURN
PROC reset
* Reset environment.
 SET TITLE ON
 SET SAFETY ON
 SET TALK ON
 DO delfiles
RETURN
PROC delfiles
* Delete working files
 DELETE FILE t71.dbf
 DELETE FILE t71.mdx
 DELETE FILE t72.dbf
 DELETE FILE t73.dbf
RETURN
PROC getyear
* Get years of data to compare. If more than 5 available, ask user for start
* year.
 TEXT
 Determining years of data for analysis
 ENDTEXT
```

```
USE emission ORDER vr chem
 SET FILTER TO source = "T"
 GO TOP
 mfirstyear = year
 GO BOTTOM
 mlastyear = year
 numyears = mlastyear - mfirstyear + 1
 IF numyears > 5
   DO askyear
 ENDIF
 CLOSE ALL
RETURN
PROC askyear
* Ask user for start year. Default gives last 5 years of data.
 askyear = mlastyear - 4
 @ 0,0 CLEAR
 SET CONFIRM ON
 @ 6, 6 TO 15,73 DOUBLE
 @ 7, 8 SAY "The trend analysis can only accomodate 5 years of data."
 @ 8, 8 SAY "WRMS currently contains"
   8,32 SAY numyears PICTURE *99*
 a
    8,35 SAY "years of data, from"
 a
 a
    8,55 SAY mfirstyear PICTURE *9999*
    8,60 SAY "to"
 a
 a
    8,63 SAY mlastyear PICTURE *9999*
   8,67 SAY "."
 a
 @ 10, 8 SAY "The analysis will be performed for 5 years, starting with the"
 @ 11, 8 SAY "year entered below. The default starting year shown will result"
 @ 12, 8 SAY "in the analysis of the most recent 5 years of data."
 @ 14, 8 SAY "Please enter the desired starting year:"
 @ 14,48 GET askyear PICTURE *9999* RANGE mfirstyear, mlastyear-1
 READ
 SET CONFIRM OFF
 mfirstyear = askyear
 mlastyear = min(mlastyear, askyear + 4)
 numyears = mlastyear - mfirstyear + 1
RETURN
PROC makefile
* Make the file for doing trend analysis
 TEXT
 Getting facility and chemical IDs for analysis
 ENDTEXT
 USE emission
 IF numyears <= 5
   SET FILTER TO source = "T"
 ELSE
   SET FILTER TO source = "T" .AND. ;
                year >= mfirstyear .AND. year <= mlastyear
 ENDIF
 GO TOP
 COPY TO t71 FIELDS fac_id, chem_id
 USE t71
 INDEX ON fac_id + chem_id TAG fac_chm_u UNIQUE
 COPY TO t72
 CLOSE ALL
 CREATE t73 FROM t7stru
 APPEND FROM t72
 CLOSE ALL
```

## RETURN

```
PROC fillfile
* Fill in data to trend analysis file from emission.
 TEXT
 Copying data to analysis file
 ENDTEXT
 USE t73 IN 1 ALIAS t7
 REPLACE ALL first_year WITH mfirstyear, num_years WITH numyears
 USE emission IN 2 ORDER fac_chm_yr ALIAS em
 loopnum = "1"
 loopyear = mfirstyear
 DO WHILE loopyear <= mlastyear
   @ 21,35 SAY loopyear PICTURE *9999*
   SET RELATION TO fac_id + chem_id + str(loopyear, 4) INTO em
   REPLACE total_amt&loopnum WITH em->fugit_amt + em->stack_amt +;
          em->water_amt + em->potw_amt + em->land_amt + em->off_amt;
          FOR found("em")
   loopnum = str(val(loopnum) + 1,1)
   loopyear = loopyear + 1
 ENDDO
 CLOSE ALL
RETURN .
PROC pattern
* Look for breaks in pattern of reporting
 TEXT
 Looking for reporting pattern discrepancies
 ENDTEXT
 USE t73
 SCAN
   i = "1"
   j = "2"
   DO WHILE val(i) < numyears
     bothblank = isblank(total_amt&i) .AND. isblank(total_amt&j)
     bothfull = .NOT. (isblank(total_amt&i) .OR. isblank(total_amt&j))
     IF .NOT. (bothblank .OR. bothfull)
      REPLACE num_chngs WITH num_chngs + 1
     ENDIF
     i = str(val(i)+1,1)
     j = str(val(j)+1,1)
   ENDDO
   IF num_chngs > 2 .OR. (num_chngs = 2 .AND. .NOT. isblank(total_amt1))
     REPLACE rpt_discrp WITH .T.
   ENDIF
 ENDSCAN
 CLOSE ALL
RETURN
PROC patt_rpt
* Generate pattern discrepancy report
 USE t73 IN 1 ALIAS t7
 USE facility IN 2 ORDER fac_id ALIAS fac
 USE chemical IN 3 ORDER chem_id ALIAS chem
 SET RELATION TO t7->fac_id INTO fac
 SELECT 2
```

SET RELATION TO t7->chem\_id INTO chem SELECT 1 SET FIELDS TO t7->fac\_id, t7->chem\_id, t7->first\_year, t7->num\_years,; t7->total\_amt1, t7->total\_amt2, t7->total\_amt3,; t7->total\_amt4, t7->total\_amt5, t7->rpt\_discrp,; fac->fac\_name, fac->city,; chem->use\_name  $_plength = 60$ REPORT FORM t71 TO FILE t71.txt FOR rpt\_discrp CLOSE ALL RETURN PROC stop \* Look for stopped reporting after reporting large quantities TEXT Deleting records with reporting pattern discrepancies ENDTEXT USE t73 DELETE FOR rpt\_discrp PACK TEXT Looking for stopped reporting discrepancies ENDTEXT SCAN IF num\_chngs = 2 .OR. (num\_chngs = 1 .AND. .NOT. isblank(total\_amt1)) i = "1" j = "2" stopped = .F.DO WHILE val(i) < num\_years .AND. .NOT. stopped stopped = (.NOT. isblank(total\_amt&i)) .AND. isblank(total\_amt&j)
IF i = "1" .AND. mfirstyear = 1987  $rpt_thresh = 50000$ ELSE  $rpt_thresh = 25000$ ENDIF IF stopped .AND. total\_amt&i > rpt\_thresh REPLACE stop\_discr WITH .T., last\_amt WITH total\_amt&i ENDIF i = str(val(i)+1,1)j = str(val(j)+1,1)ENDDO ENDIF ENDSCAN CLOSE ALL RETURN PROC stop\_rpt \* Generate stopped reporting discrepancy report USE t73 IN 1 ALIAS t7 INDEX ON last\_amt TAG last\_amt DESCENDING USE facility IN 2 ORDER fac\_id ALIAS fac USE chemical IN 3 ORDER chem\_id ALIAS chem SET RELATION TO t7->fac\_id INTO fac SELECT 2 SET RELATION TO t7->chem\_id INTO chem

## SELECT 1

```
SET FIELDS TO t7->fac_id, t7->chem_id, t7->first_year, t7->num_years,;
               t7->total_amt1, t7->total_amt2, t7->total_amt3,;
t7->total_amt4, t7->total_amt5, t7->rpt_discrp,;
fac->fac_name, fac->city,;
               chem->use_name
  plength = 76
  REPORT FORM t72 TO FILE t72.txt FOR stop_discr
  CLOSE ALL
RETURN
PROC limitrecs
* Delete records not suitable for quantity comparison
  TEXT
  Deleting records with stopped reporting discrepancies
  ENDTEXT
  USE t73
  DELETE FOR stop_discr
  PACK
  last = str(numyears,1)
  nextlast = str(numyears-1,1)
  TEXT
  Deleting records with only one year of data
  ENDTEXT
  SCAN
   DO CASE
     CASE (.NOT. isblank(total_amt1)) .AND. isblank(total_amt2)
       DELETE
     CASE (.NOT. isblank(total_amt&last)) .AND. isblank(total_amt&nextlast)
       DELETE
   ENDCASE
  ENDSCAN
  PACK
  CLOSE ALL
RETURN
PROC calc_discr
* Calculate discrepancies between total quantities reported over time
  TEXT
  Calculating discrepancies
  ENDTEXT
  USE t73
  SCAN
    i = "1"
    j = "2"
    DO WHILE val(i) < numyears
      IF .NOT. (isblank(total_amt&i) .OR. isblank(total_amt&j))
       discrp =
log10 (max(max(total_amt&i,total_amt&j),1)/max(min(total_amt&i,total_amt&j),1))
       absolute = abs(total_amt&i - total_amt&j)
       REPLACE max_discrp WITH max(max_discrp,round(discrp,1)),;
               max_abs WITH max(max_abs,absolute)
      ENDIF
```

```
i = str(val(i)+1,1)
     j = str(val(j)+1,1)
   ENDDO
 ENDSCAN
 CLOSE ALL
RETURN
PROC sum_discrp
* Summarize all discrepancies
 TEXT
 Summarizing discrepancies
 ENDTEXT
 USE t73
 SCAN
   IF max_discrp >= 1 .AND. max_abs >= 3000
     REPLACE signif WITH .T.
   ENDIF
 ENDSCAN
 CLOSE ALL
RETURN
PROC discrp_rpt
* Generate quantity discrepancy report
 USE t73 IN 1 ALIAS t7
 USE facility IN 2 ORDER fac_id ALIAS fac
USE chemical IN 3 ORDER chem_id ALIAS chem
 SET RELATION TO t7->fac_id INTO fac
 SELECT 2
 SET RELATION TO t7->chem_id INTO chem
 SELECT 1
 SET FIELDS TO t7->fac_id, t7->chem_id, t7->first_year, t7->num_years,;
              t7->total_amt1, t7->total_amt2, t7->total_amt3,;
t7->total_amt4, t7->total_amt5, t7->rpt_discrp,;
               t7->signif, t7->max_discrp, t7->max_abs,;
              fac->fac_name, fac->city,;
              chem->use_name
 _plength = 76
 REPORT FORM t73 TO FILE t73.txt FOR signif
 CLOSE ALL
```

```
RETURN
```

## APPENDIX E RESULTS OF COMPARISON OF TRI DATA AND SINGLE MEDIA DATA

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Page	

/93 Facilities With Signifi	lcant Discrepancies	Between TR	I and NPDES	in 1987 Page 1
Chemical	TRI	NPDES	Difference	
SODIUM HYDROXIDE	42,000	NR	42,000	
CHLORINE	1,826	NR	1,826	
AMMONIA	16,500	NR	16,500	
SODIUM SULFATE (SOLUTION)	1,500,000	NR	1,500,000	
SODIUM HYDROXIDE	711,000	NR	711,000	
SODIUM SULFATE (SOLUTION)	4,970,400	NR	4,970,400	
SODIUM SULFATE (SOLUTION)	1,155,200	NR	1,155,200	
CHLORINE AMMONIA	2,225 3,026	NR NR	2,225 3,026	
SODIUM SULFATE (SOLUTION)	4,378,000	NR	4,378,000	
SODIUM SULFATE (SOLUTION)	8,950,000	NR	8,950,000	
XYLENE 1,2,4-TRICHLOROBENZENE SODIUM SULFATE (SOLUTION)	5,540 1,100 2,300,000	NR NR NR	5,540 1,100 2,300,000	
1,2,4-TRICHLOROBENZENE SODIUM SULFATE (SOLUTION) BIPHENYL	3,500,000 2,500	NR NR NR	1,000 3,500,000 2,500	
ZINC SODIUM SULFATE (SOLUTION) COPPER	NR 6,060,000 Nr	21,385 NR 208,206	21,385 6,060,000 208,206	
SODIUM SULFATE (SOLUTION)	20,066,356	NR	20,066,356	
AMMONIUM SULFATE (SOLUTION)	30,680	NR	30,680	
SODIUM SULFATE (SOLUTION)	137,300	NR	137,300	
SODIUM SULFATE (SOLUTION)	1,000,000	NR	1,000,000	
SODIUM SULFATE (SOLUTION) CHROMIUM	759,334 NR	NR 1,566	759,334 1,566	
1,2,4-TRICHLOROBENZENE SODIUM SULFATE (SOLUTION) AMMONIUM SULFATE (SOLUTION)	6,365 341,737 11,453	NR NR NR	6,365 341,737 11,453	
SODIUM SULFATE (SOLUTION) CHLORINE	956,000 7,000	NR NR	956,000 7,000	
SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE AMMONIUM SULFATE (SOLUTION)	637,910 981,532 45,862	NR NR NR	637,910 981,532 45,862	
SODIUM SULFATE (SOLUTION)	75,000	NR	75,000	
SODIUM HYDROXIDE	3,800	NR	3,800	
1,2,4-TRICHLOROBENZENE	52,249	NR	52,249	
quantities reported in lbs.	NR = Not Repor	ted	** = Val	ues reported to both TRI and NPDES
	<pre>/93 Facilities With Signif Chemical soblum HYDROXIDE CHLORINE AMMONIA SODIUM SULFATE (SOLUTION) SODIUM SULFATE (SOLUT</pre>	/93     Facilities with significant Discrepancies       Chemical     TRI       SODIUM HYDROXIDE     42,000       CHLORINE     1,826       AMMONIA     1,500,000       SODIUM HYDROXIDE     1,826       AMMONIA     1,500,000       SODIUM HYDROXIDE     711,000       SODIUM SULFATE (SOLUTION)     1,155,200       SODIUM SULFATE (SOLUTION)     1,155,200       SODIUM SULFATE (SOLUTION)     8,950,000       SODIUM SULFATE (SOLUTION)     8,950,000       SODIUM SULFATE (SOLUTION)     3,000,000       SULPATE (SOLUTION) <t< td=""><td>/93       Facilities with Significant Discrepancies Between TR         Chemical       TRI       NPDES         SODIUM HYDROXIDE       42,000       NR         CHLORINE       1,826       NR         AMMONIA       1,500,000       NR         SODIUM HYDROXIDE       711,000       NR         SODIUM SULFATE (SOLUTION)       1,500,000       NR         SODIUM SULFATE (SOLUTION)       711,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       1,24,7130       NR         SODIUM SULFATE (SOLUTION)</td><td>/93     Facilities with Significant Discrepancies Between TR1 and NDES       Chemical     TR1     NPDES     Difference       soDINM NCPROXIDE     1,826     NR     1,826       SCHLDRINE     1,826     NR     1,826       SODINM NCPROXIDE     1,826     NR     1,6500       SODINM NCPROXIDE     1,826     NR     1,6500       SODINM NCPROXIDE     1,1000     NR     1,6500       SODINM SULFATE (SOLUTION)     1,5500000     NR     1,6500       SODINM SULFATE (SOLUTION)     1,550000     NR     1,550000       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,5520       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,155,200       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,155,200       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,155,200       SODINM SULFATE (SOLUTION)     1,155,200     NR     2,326       SODINM SULFATE (SOLUTION)     2,326,000     NR     2,326&lt;</td></t<>	/93       Facilities with Significant Discrepancies Between TR         Chemical       TRI       NPDES         SODIUM HYDROXIDE       42,000       NR         CHLORINE       1,826       NR         AMMONIA       1,500,000       NR         SODIUM HYDROXIDE       711,000       NR         SODIUM SULFATE (SOLUTION)       1,500,000       NR         SODIUM SULFATE (SOLUTION)       711,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       3,500,000       NR         SODIUM SULFATE (SOLUTION)       1,155,200       NR         SODIUM SULFATE (SOLUTION)       1,24,7130       NR         SODIUM SULFATE (SOLUTION)	/93     Facilities with Significant Discrepancies Between TR1 and NDES       Chemical     TR1     NPDES     Difference       soDINM NCPROXIDE     1,826     NR     1,826       SCHLDRINE     1,826     NR     1,826       SODINM NCPROXIDE     1,826     NR     1,6500       SODINM NCPROXIDE     1,826     NR     1,6500       SODINM NCPROXIDE     1,1000     NR     1,6500       SODINM SULFATE (SOLUTION)     1,5500000     NR     1,6500       SODINM SULFATE (SOLUTION)     1,550000     NR     1,550000       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,5520       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,155,200       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,155,200       SODINM SULFATE (SOLUTION)     1,155,200     NR     1,155,200       SODINM SULFATE (SOLUTION)     1,155,200     NR     2,326       SODINM SULFATE (SOLUTION)     2,326,000     NR     2,326<

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TRI
Between
Discrepancies
significant
With
Facilities
07/22/93

SIC	Chemical	TRI	NPDES	Difference	
2269 2269 2269 2269	TETRACHLOROETHYLENE SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE PSEUDOCUMENE BIPHENYL	42,192 259,700 68,450 17,575 13,256	N N N N N N N N N N N N N N N N N N N	42,192 259,700 68,450 17,575 13,256	•
2283	SODIUM SULFATE (SOLUTION)	180,000	NR	180,000	
2299	SODIUM SULFATE (SOLUTION)	151,148	NR	151,148	
2396	CHLORINE	42,000	NR	42,000	
2421	SODIUM HYDROXIDE	1,000	NR	1,000	
2493	FORMALDEHYDE	2,248	NR	2,248	
2600 2600 2600 2600 2600 2600	SODIUM SULFATE (SOLUTION) METHANOL CHROMIUM CHLORINE DIOXIDE CHLORINE AMMONIA ACETONE	1,608,808 187,697 3,196 4,865 1,762 13,388 2,683	N N N N N N N N N N N N N N N N N N N	1,608,808 187,697 3,196 4,865 11,762 13,388 2,683	
2610 2610 2610 2610	SODIUM SULFATE (SOLUTION) CHLOROFORM AMMONIA ACETONE	60,647,550 14,070 1,292 9,010	NR NR NR	60,647,550 14,070 1,292 9,010	
2611	AMMONIUM NITRATE (SOLUTION)	1,800	NR	1,800	
2611 2611	SODIUM SULFATE (SOLUTION) ACETONE	32,000,000 6,300	NR NR	32,000,000 6,300	
2611 2611 2611	SODIUM SULFATE (SOLUTION) AMMONIA ACETONE	16,989,000 6,050 3,900	NR NR NR	16,989,000 6,050 3,900	
2641 2641 2641 2641	SODIUM SULFATE (SOLUTION) CHLOROFORM AMMONIA ACETONE	24,000,000 28,000 490,000 8,100	NR NR NR	24,000,000 28,000 490,000 8,100	
2819	SODIUM HYDROXIDE	23,000	NR	23,000	
2819	SODIUM SULFATE (SOLUTION)	102,943	NR	102,943	
2819	SODIUM SULFATE (SOLUTION)	855,130	NR	855,130	
2819	HYDROCHLORIC ACID	12,500	NR	12,500	
2819	SODIUM SULFATE (SOLUTION)	1,900,000	NR	1,900,000	
2824	PHOSPHORIC ACID	10,000	NR	10,000	
2824 2824 2824 2824 2824	SODIUM SULFATE (SOLUTION) 2-METHOXYETHANOL METHANOL ETHYLENE GLYCOL 1,4-DIOXANE	1,187,245 2,200 5,143 19,343 2,290	NR NR NR NR	1,187,245 2,200 5,143 19,343 2,290	

\*\* = Values reported to both TRI and NPDES

NR = Not Reported

All quantities reported in lbs.

Page 2

07/22	/93 Facilities With Signific	cant Discrepancies Bet	ween TR	I and NPDES i	n 1987 Page 3
SIC	Chemical	TRI	NPDES	Difference	
2824 2824 2824	CHLORINE	1,211 8,643	NR NR	1,211 8,643	
2824 2824 2824 2824	METHANOL ETHYLENE GLYCOL 1,4-DIOXANE AMMONIA	1,936 60,990 10,606 4,180	NR NR NR	1,936 60,990 10,606 4,180	
2824 2824	ACETALDEHYDE METHANOL	9,976	NR NR	9,9/6	
2824	ETHYLENE GLYCOL	6,700	NR	9,700	
2833	ETHYLENE GLYCOL	20,000	NR	20,000	
2834	SODIUM SULFATE (SOLUTION)	369,000	NR	369,000	
2834	SODIUM SULFATE (SOLUTION)	487,746	NR	487,746	
2834	SODIUM SULFATE (SOLUTION)	124,800	NR	124,800	
2861	SODIUM SULFATE (SOLUTION)	7,272,000	NR	7,272,000	
2865 2865 2865	SODIUM SULPATE (SOLUTION) GLYCOL ETHERS ETHYLENE GLYCOL	3,800,000 3,700 1,300	NR NR NR	3,800,000 3,700 1,300	
2869 2869 2869 2869	METHANOL FORMALDEHYDE COBALT AMMONIA	1,200 1,900 1,700 7,300	NR NR NR	1,200 1,900 1,700 7,300	
2873	AMMONIUM NITRATE (SOLUTION)	42,000	NR	42,000	
2874 2874	SODIUM SULFATE (SOLUTION) n-BUTYL ALCOHOL	3,055,000 21,000	NR NR	3,055,000 21,000	
2879	FORMALDEHYDE	8,000	NR	8,000	
2899	SODIUM SULFATE (SOLUTION)	15,330	NR	15,330	
3079 3079	SODIUM SULFATE (SOLUTION) AMMONIA	1,400,000 1,800	NR NR	1,400,000 1,800	
3229	PHOSPHORIC ACID	3,746	NR	3,746	
3231 3231	SULFURIC ACID SODIUM HYDROXIDE	25,000 83,000	NR NR	25,000 83,000	
3356 3356	SODIUM SULFATE (SOLUTION) AMMONIA	170,000 4,400	NR NR	170,000 4,400	
3519	HYDROCHLORIC ACID	10,282	NR	10,282	
3552 3552	SULFURIC ACID SODIUM HYDROXIDE	45,409 33,163	NR NR	45,409 33,163	
3553 3553	GLYCOL ETHERS ETHYLENE GLYCOL	12,000 24,000	NR NR	12,000 24,000	
All q	puantities reported in lbs.	NR = Not Reported		** = Valu	les reported to both TRI and NPDES

07/22/93	Facilities with Significant	Discrepancies	Between TI	KI and NPDES	in 1987
SIC Chemical		TRI	NPDES	Difference	
3692 ZINC COMP	SUNDO	3,500	NR	3,500	

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1,400

NR NR

1,400

3861 METHANOL

4

•\* = Values reported to both TRI and NPDES

NR = Not Reported

07/22	2/93 Facilities With Significan	t Discrepancies Be	tween TF	II and NPDES	in 1988 Page 1
SIC	Chemical	TRI	NPDES	Difference	
	CHLORINE	1,500	NR	1,500	
1479	SODIUM HYDROXIDE	72,000	NR	72,000	
2200	SODIUM HYDROXIDE	527,500	NR	527,500	
2221	AMMONIA	3,479	NR	3,479	
2252	AMMONIUM SULFATE (SOLUTION)	14,062	NR	14,062	
2253	1,2,4-TRICHLOROBENZENE	2,255	NR	2,255	
2257	BIPHENYL	5,710	NR	5,710	
2257	METHANOL	4,293	NR	4,293	
2261	SODIUM HYDROXIDE	103,062	NR	103,062	
2262	1,2,4-TRICHLOROBENZENE	5,891	NR	5,891	
2262	CHLORINE	1,300	NR	1,300	
2269 2269	SODIUM HYDROXIDE AMMONIUM SULFATE (SOLUTION)	495,286 91,104	NR NR	495,286 91,104	
2272	AMMONIA	9,400	NR	9,400	
2299 2299	GLYCOL ETHERS COPPER COMPOUNDS	5,200 15,000	NR NR	5,200 15,000	
2396	CHLORINE	50,000	NR	50,000	
2421	SODIUM HYDROXIDE	3,371	NR	3,371	
2493 2493 2493	FORMALDEHYDE AMMONIA ALUMINUM OXIDE	2,000 2,900 4,600	NR NR NR	2,000 2,900 4,600	
2600 2600 2600 2600	CHROMIUM COMPOUNDS CHLORINE CATECHOL AMMONIA ACETONE	3,000 2,500 10,000 2,500	NR NR NR NR	3,000 2,100 10,000 2,500	
2610 2610 2610 2610	CHLOROFORM CATECHOL AMMONIA ACETONE	13,000 1,200 9,500	NR NR NR	13,000 1,800 1,200 9,500	
2611 2611	CATECHOL ACETONE	1,600 6,300	NR NR	1,600 6,300	
2611 2611 2611 2611	CHLOROFORM CATECHOL ACETONE	10,100 1,000 3,600	NR NR NR	10,100 1,000 3,600	
2621	AMMONIA	2,500	NR	2,500	
2641 2641	CHLOROFORM AMMONIA	28,000 8,300	NR NR	28,000 8,300	
All q	puantities reported in lbs.	NR = Not Reported		** = Val	ues reported to both TRI and NPDES

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SIC	Chemical	TRI	NPDES	Difference	
2641	ACETONE	7,500	NR	7,500	
2645	NITRIC ACID	1,750	NR	1,750	
2819	SODIUM HYDROXIDE	22,000	NR	22,000	
2819	HYDROCHLORIC ACID	12,000	NR	12,000	
2824 2824 2824 2824 2824 2824 2824 2824	2-METHOXYETHANOL METHANOL MANGANESE COMPOUNDS ETHYLENE GLYCOL COBALT COMPOUNDS CHLORINE AMMONIA	2,400 1,400 15,300 3,100 1,200	N N N N N N N N N N N N N N N N N N N	2,400 1,400 2,300 3,100 1,200 1,200	
2824 2824 2824 2824 2824	ETHYLENE GLYCOL 1,4-DIOXANE AMMONIA ACETALDEHYDE	30,560 7,494 7,414 5,220	NR NR NR	30,560 7,494 7,414 5,220	
2824 2824	METHANOL ETHYLENE GLYCOL	9,130 1,900	NR NR	9,130 1,900	
2824 2824	ETHYLENE GLYCOL 1,4-DIOXANE	6,500 17,000	NR NR	6,500 17,000	
2833	ETHYLENE GLYCOL	8,200	NR	8,200	
2865 2865	GLYCOL ETHERS ETHYLENE GLYCOL	7,700 4,400	NR NR	7,700	
2869 2869 2869 2869	TEREPHTHALIC ACID FORMALDEHYDE COBALT AMMONIA	1,800 1,500 5,470 6,800	NR NR NR	1,800 1,500 5,470 6,800	
2873	AMMONIUM NITRATE (SOLUTION)	50,735	NR	50,735	
3079	AMMONIA	1,800	NR	1,800	
3231 3231	SULFURIC ACID SODIUM HYDROXIDE	27,000 90,000	NR NR	27,000	
3245	ZINC COMPOUNDS	2,105	NR	2,105	
3351	COPPER COMPOUNDS	1,023	NR	1,023	
3356	AIMONIA	10,160	NR	10,160	
3440	ALUMINUM (FUME OR DUST)	6,395	NR	6,395	
3553 3553	GLYCOL ETHERS ETHYLENE GLYCOL	12,000 30,000	NR NR	12,000 30,000	
3573 3573	HYDROCHLORIC ACID AMMONIA	1,700 15,000	NR NR	1,700 15,000	
3675	NICKEL	NR	2,193	2,193	
3861	METHANOL	3,200	NR	3,200	
All o	quantities reported in lbs.	NR = Not Repor	ted	** = Va	lues reported to both TRI and NPDES

Facilities With Significant Discrepancies Between TRI and NPDES in 1988

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07/22/93

07/22	<pre>/93 Facilities With Significant</pre>	Discrepancies Bet	tween TRI	and NPDES 11	1 1989 Page 1
SIC	Chemical	TRI	NPDES	Difference	
2015	CHLORINE	18,750	NR	18,750	
2077	AMMONIA	29,908	NR	29,908	
2082	AMMONIA	2,100	NR	2,100	
2221	AMMONIA	13,000	NR	13,000	
2252 2252	CHLORINE AMMONIUM SULFATE (SOLUTION)	3,317 23,527	NR NR	3,317 23,527	
2253 2253	1,2,4-TRICHLOROBENZENE FORMALDEHYDE	1,500 1,200	NR NR	1,500 1,200	
2253 2253	1,2,4-TRICHLOROBENZENE FORMALDEHYDE	1,100 1,500	NR NR	1,100 1,500	
2257	METHANOL	7,662	NR	7,662	
2261	AMMONIA	5,623	NR	5,623	
2396	CHLORINE	32,000	NR	32,000	
2493 2493 2493	FORMALDEHYDE AMMONIA ALUMINUM OXIDE	1,900 2,500 4,000	NR NR NR	1,900 2,500 4,000	
2600 2600 2600 2600	CHROMIUM COMPOUNDS CATECHOL AMMONIA ACETONE	4,900 8,900 2,300	NR NR NR	4,900 8,900 3,900	
2610 2610 2610 2610	CHLOROFORM CATECHOL AMMONIA ACETONE	10,600 1,900 1,600 9,300	NR NR NR	10,600 1,900 1,600 9,300	
2611 2611	CATECHOL ACETONE	1,700 6,800	NR NR	1,700 6,800	
2611 2611 2611	CHLOROFORM CATECHOL ACETONE	16,300 1,030 3,875	NR NR NR	16,300 1,030 3,875	
2621 2621	XYLENE AMMONIA	1,335 2,088	NR NR	1,335 2,088	
2641 2641 2641 2641 2641	PHENOL CHLOROFORM CATECHOL AMMONIA ACETONE	15,000 31,000 1,500 7,600	NN NN NN NN NN	15,000 31,000 1,500 8,100 7,600	
2824 2824 2824 2824 2824 2824 2824 2824	2-METHOXYETHANOL METHANOL MANGANESE COMPOUNDS ETHYLENE GLYCOL I.4-DIOXANE COBALT COMPOUNDS CHLORINE	9,760 80,100 1,100 66,240 3,500 1,790	<b>H H H H H H H</b>	9,760 80,100 1,100 66,200 3,500 1,790	
All c	quantities reported in lbs.	VR = Not Reported		** = Valu	es reported to both TRI and NPDES

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07/22	<pre>/93 Facilities With Significant</pre>	Discrepancies	Between TF	I and NPDES	in 1989
SIC	Chemical	TRI	NPDES	Difference	
2824 2824 2824	n-BUTYL ALCOHOL AMMONIA	9,960 5,630	NR NR	5,630	
2824 2824 2824 2824	ETHYLENE GLYCOL 1,4-DIOXANE AMMONIA ACETALDEHYDE	30,468 14,958 5,843 4,348	NR NR NR	30,468 14,958 5,843 4,348	•
2824 2824 2824 2824 2824	METHANOL ETHYLENE GLYCOL 1,4-DIOXANE ACETALDEHYDE	1,800 11,000 32,000 1,500	NR NR NR	1,800 11,000 32,000 1,500	
2833	ETHYLENE GLYCOL	7,000	NR	7,000	
2865 2865	ETHYLENE GLYCOL o-ANISIDINE	4,464 4,699	NR NR	4,464 4,699	
2865 2865 2865	GLYCOL ETHERS ETHYLENE GLYCOL AMMONIA	3,900 3,700 85,000	NR NR NR	3,900 3,700 85,000	
2869 2869 2869 2869 2869	TEREPHTHALIC ACID METHANOL FORMALDEHYDE COBALT AMMONIA	1,400 1,100 2,000 9,800	N N N N N N N N N N N N N N N N N N N	1,400 1,100 2,000 9,800	
2869 2869	SULFURIC ACID AMMONIA	2,868 1,900	NR NR	2,868 1,900	
2873	AMMONIUM NITRATE (SOLUTION)	82,125	NR	82,125	
2874	AMMONIA	48,000	NR	48,000	
2875	AMMONIUM NITRATE (SOLUTION)	3,500	NR	3,500	
3229	CHLORINE	1,061	NR	1,061	
3351	COPPER COMPOUNDS	1,023	NR	1,023	
3356	AMMONIA	10,000	NR	10,000	
3553 3553	GLYCOL ETHERS ETHYLENE GLYCOL	7, <b>44</b> 7 21,300	NR NR	7,447 21,300	
3861	METHANOL	4,400	NR	4,400	
4941	CHLORINE	1,024	NR	1,024	

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\*\* = Values reported to both TRI and NPDES

1,024

1,024

4941 CHLORINE

NR = Not Reported

All quantities reported in lbs.

07/22	/93 Fa	acilities With Significant	Discrepancies	Between TF	XI and NPDES in 1990	-
SIC	Chemical		TRI	NPDES	Difference	
2016	AMMONIA		30,000	NR	30,000	
2077	AMMONIA		12,062	NR	12,062	
2082	AMMONIA		5,000	NR	5,000	•
2221 2221	CHROMIUM AMMONIA		NR 8,600	1,098 NR	1,098 8,600	
2252	CHROMIUM		NR	3,042	3,042	
2252	AMMONIUM SU	LFATE (SOLUTION)	27,700	NR	27,700	
2253	FORMALDEHYD	ы	1,188	NR	1,188	
2261	AMMONIA		7,332	NR	7,332	
2262	AMMONIA		2,618	NR	2,618	
2262	GLYCOL ETHE	RS	7,000	NR	7,000	
2262 2262	Z INC COPPER		NR NR	32,926 14,497	32,926 14,497	
2396	CHLORINE		44,000	NR	44,000	
2600 2600 2600 2600	CHROMIUM CO CATECHOL AMMONIA ACETONE	MPOUNDS	4,900 2,000 17,000 2,400	NR NR NR	4,900 2,000 17,000 2,400	
2610 2610 2610 2610	CHLOROFORM CATECHOL AMMONIA ACETONE		8,600 1,900 9,300	NN NN NN NN	8,600 1,900 9,300	
2611 2611 2611	CATECHOL AMMONIA ACETONE		1,800 26,000 7,100	NR NR NR	1,800 26,000 7,100	
2611 2611 2611	CHLOROFORM CATECHOL ACETONE		5,230 1,060 3,740	NR NR NR	5,230 1,060 3,740	
2621	AMMONIA		2,354	NR	2,354	
2641 2641 2641 2641 2641 2641	PHENOL METHANOL CHLOROFORM CATECHOL AMMONIA ACETONE		15,000 100,000 19,000 4,600 340,000	N N N N N N N N N N N N N N N N N N N	15,000 100,000 19,000 340,000 7,300	
2824 2824 2824 2824 2824 2824 2824	2-METHOXYET METHANOL MANGANESE C ETHYLENE GL 1,4-DIOXANE COBALT COMP CHLORINE	HANOL OMPOUNDS YCOL OUNDS	4,755 124,478 26,784 26,690 31,880 37,500 1,460	N N N N N N N N N N N N N N N N N N N	124,755 124,478 26,784 26,784 1,880 31,500 1,460	

÷ Page \*\* = Values reported to both TRI and NPDES

NR = Not Reported

All quantities reported in lbs.

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SIC	Chemical	TRI	NPDES	Difference
2824 2824 2824	n-BUTYL ALCOHOL AMMONIA	8,150 16,350	NR NR	8,150 16,350
2824 2824 2824 2824	ETHYLENE GLYCOL 1,4-DIOXANE AMMONIA ACETALDEHYDE	21,533 12,921 4,438 3,731	NR NR NR	21,533 12,921 4,438 3,731
282 <b>4</b> 282 <b>4</b> 2824 2824	METHANOL ETHYLENE GLYCOL 1,4-DIOXANE ACETALDEHYDE	2,300 14,000 28,000 2,300	NR NR NR	2,300 14,000 28,000 2,300
2833	ETHYLENE GLYCOL	2,100	NR	2,100
2833	AMMONIA	14,000	NR	14,000
2865	ETHYLENE GLYCOL	8,211	NR	8,211
2865 2865 2865	GLYCOL ETHERS 2,4-DINITROPHENOL AMMONIA	4,400 1,600 43,000	NR NR NR	4,400 1,600 43,000
2869 2869 2869	FORMALDEHYDE COBALT AMMONIA	1,800 2,300 12,000	NR NR NR	1,800 2,300 12,000
2873	AMMONIUM NITRATE (SOLUTION)	70,810	NR	70,810
2874	AMMONIA	80,000	NR	80,000
3079	AMMONIA	1,145	NR	1,145
3356	AMMONIA	8,300	NR	8,300
3675	ETHYLENE GLYCOL	4,057	NR	4,057
3861	METHANOL	4,300	NR	4,300
4941	CHLORINE	1,061	NR	1,061

All quantities reported in lbs.

NR = Not Reported

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07/22	/93 Facilities With Significan	it Discrepancies	Between TRI	nd Pretreatment in 1987	
SIC	Chemical	TRI	Pretreat	Difference	
20	SODIUM HYDROXIDE	608,098	NR	608,098	
20	SODIUM HYDROXIDE PHENOL	250,000 NR	NR 765,005	250,000 765,005	
2011	SODIUM HYDROXIDE	63,000	NR	63,000	
2021 2021	SODIUM HYDROXIDE PHOSPHORIC ACID	<b>4</b> 1,000 <b>4</b> 0,000	NR NR	41,000 40,000	
2023	SODIUM HYDROXIDE	10,000	NR	10,000	
2024	SODIUM HYDROXIDE	37,191	NR	37,191	
2024 2024	SODIUM HYDROXIDE PHOSPHORIC ACID	149,892 26,607	NR NR	149,892 26,607	
2026 2026	SODIUM HYDROXIDE PHOSPHORIC ACID	14,270 11,657	NR NR	14,270 11,657	
2026 2026	SODIUM HYDROXIDE PHOSPHORIC ACID	40,607 25,673	NR NR	40,607 25,673	
2026	SODIUM HYDROXIDE	20,000	NR	20,000	
2046 2046	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	7,000,000 280,000	NR NR	7,000,000 280,000	
2086 2086	SODIUM HYDROXIDE BARIUM	73,841 NR	NR 62,049	73,841 62,049	
2111	TOLUENE	NR	118,798	118,798	
2111	SODIUM HYDROXIDE	15,450	NR	15,450	
22	SODIUM HYDROXIDE	470,000	NR	470,000	
22 22	SULFURIC ACID SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	600,000 2,154,800 950,000	NR NR NR	600,000 2,154,800 950,000	
2200 2200	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	1,500,000 517,000	NR NR	1,500,000 517,000	
2200	SODIUM HYDROXIDE	711,000	NR	711,000	
2211 2211	SODIUM SULFATE (SOLUTION) GLYCOL ETHERS	1,145,300 45,800	NR NR	1,145,300 45,800	
2211	ARSENIC	NR	47,183	47,183	
2211 2211	TOLUENE ARSENIC	NR NR	33,947 46,757	33,947 46,757	
2221 2221 2221	SULFURIC ACID SODIUM HYDROXIDE AMMONIA	280,000 490,000 28,000	NR NR NR	280,000 490,000 28,000	
2221 2221	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	45,000 18,000	NR NR	45,000 18,000	
e lia	puantities reported in lbs.	NR = Not Repor	ted	<pre>** = Values reported to bot</pre>	h TRI and Pretreatment

07/22	2/93 Facilities With Significant	Discrepanc	ies Between TR)	I and Pretreatmen'	t in 1987
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2221 2221 2221 2221 2221 2221	TOLUENE SILVER SELENIUM MERCURY LEAD COPPER	NR NR NR NR NR NR	39,483 78,483 43,876 114,062 26,322 144,771	39,483 78,966 43,876 114,062 26,322 144,771	
2221 2221	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	62,500 119,500	NR NR	62,500 119,500	
2221 2221 2221 2221	TOLUENE SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE PHENOL	NR 30,800 52,994 NR	28,060 NR NR 1,220,000,000	28,060 30,800 52,994 1,220,000,000	
2221 2221 2221 2221 2221 2221 2221	TOLUENE SILVER SELENIUM MERCURY LEAD COPPER	n n n n n n n	115,138 138,165 13,817 101,321 13,817 13,817 18,422 245,625,132	115,138 138,165 13,817 13,817 101,321 13,817 13,817 245,625,132	
2221 2221 2221	SILVER PHENOL ANTIMONY	NR NR NR	1,525,701,195 50,857	1,525,701,195 50,857	
2221 2221	SODIUM HYDROXIDE ETHYLENE GLYCOL	143,000 14,785	NR NR	143,000 14,785	
2221 2221	TETRACHLOROETHYLENE SODIUM SULFATE (SOLUTION)	32,000 420,000	NR NR	32,000 420,000	
2231	SODIUM SULFATE (SOLUTION)	182,210	NR	182,210	
2231	SILVER	NR	37,988	37,988	
2251	SODIUM SULFATE (SOLUTION)	12,000	NR	12,000	
2251	PHENOL	NR	34,823	34,823	
2251	SODIUM SULFATE (SOLUTION)	105,000	NR	105,000	
2251	AMMONIUM SULFATE (SOLUTION)	60,000	NR	60,000	
2252 2252	SODIUM SULFATE (SOLUTION) AMMONIUM SULFATE (SOLUTION)	49,060 40,000	NR NR	<b>49,060</b> <b>40,000</b>	
2252	AMMONIUM SULFATE (SOLUTION)	10,800	NR	10,800	
2252	CHLORINE	171,787	NR	171,787	
2252 2252	CHLORINE AMMONIUM SULFATE (SOLUTION)	32,928 11,850	NR NR	32,928 11,850	
2252	SODIUM SULFATE (SOLUTION)	20,000	NR	20,000	
2252 2252	MERCURY Copper	NR NR	16,242 10,258	16,242 10,258	
All	quantities reported in lbs.	NR = Not R	eported	** = Values r	eported to both TRI and Pretreatment

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	TRI	Pretreat	Difference
	NR	64,050 10,169	64,050 10,169
LENE	38,853	NR	38,853
	77,242	NR	77,242
	29,412	NR	29,412
(SOLUTION)	25,100	NR	25,100
E (SOLUTION)	19,400	NR	19,400
	NR	57,645	57,645
	NR	17.080	17,080

SIC	Chemical	TRI	Pretreat	Difference	
2257	LEAD	NR	64,050	64,050	
2257	Copper	NR	10,169	10,169	
2258	XYLENE	38,853	NR	38,853	•
2258	TETRACHLOROETHYLENE	77,242	NR	77,242	
2258	BIPHENYL	29,412	NR	29,412	
2258	SODIUM SULFATE (SOLUTION)	25,100	NR	25,100	
2258	AMMONIUM SULFATE (SOLUTION)	19,400	NR	19,400	
2258	TOLUENE	NR	57,645	57,645	
2258	MERCURY	NR	17,080	17,080	
2258 2258 2258 2258 2258 2258	TOLUENE SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE SILVER AMMONIUM SULFATE (SOLUTION)	12,384 234,000 40,872 NR 67,620	NR NR NR 977,353 NR	12,384 234,000 40,872 977,353 67,620	
2258	SODIUM HYDROXIDE	29,750	NR	29,750	
2258	AMMONIUM SULFATE (SOLUTION)	290,000	NR	290,000	
226	SULFURIC ACID	271,000	NR	271,000	
2260	SILVER	NR	20,506	20,506	
2261	SODIUM SULFATE (SOLUTION)	6,050,200	NR	6,050,200	
2261	GLYCOL ETHERS	28,600	NR	28,600	
2261	ETHYLENE GLYCOL	13,300	NR	13,300	
2261	SILVER	NR	58,721	58,721	
2261	SULFURIC ACID	50,000	NR	50,000	
2261	SODIUM HYDROXIDE	167,000	NR	167,000	
2262	SODIUM HYDROXIDE	16,200	NR	16,200	
2262 2262 2262 2262	1,2,4-TRICHLOROBENZENE SODIUM SULFATE (SOLUTION) AMMONIUM SULFATE (SOLUTION) AMMONIA	23,995 68,911 26,265 29,012	NR NR NR	23,995 68,911 26,265 29,012	
2262	PHENOL	NR	18,28 <b>4</b>	18,28 <b>4</b>	
2262	MERCURY	NR	12,189	12,189	
2262	SILVER	NR	12,525	12,525	
2262	SODIUM HYDROXIDE	75,687	NR	75,687	
2262	SODIUM HYDROXIDE	400,000	NR	400,000	
2262	SODIUM SULFATE (SOLUTION)	680,000	NR	680,000	
2262	AMMONIUM SULFATE (SOLUTION)	75,000	NR	75,000	
2262	SODIUM SULFATE (SOLUTION)	84,270	NR	84,270	
2262	SODIUM HYDROXIDE	148,803	NR	148,803	
2269	TETRACHLOROETHYLENE	21,528	NR	21,528	
2269	SODIUM SULFATE (SOLUTION)	32,182	NR	32,182	
2269	SODIUM HYDROXIDE	13,777	NR	13,777	

\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

All quantities reported in lbs.

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Difference	70,786 556,175 80,898 141,572 30,337 30,337 10,112 505,613,750 505,613,750 222,470,050 10,112 222,470,050	50,000 940,000	50,349 360,000 20,979 12,587 12,587	45,600 38,475 42,000	15,000 880,000 432,000	75,000	1,575,000 63,500 351,717	116,053	14,234	21,000 31,000 54,900 14,647 28,000	16,013 24,850 26,688 64,050 21,350 10,675	17,306,707	264,900	10,235	18,304	** = Values
Pretreat	70,786 556,175 80,898 80,898 30,337 30,337 30,337 10,112 505,613,750 505,613,750 505,613,750 222,470,050	NR NR	50,349 NR 20,979 12,587	NR NR NR	NR NR NR	NR	NR NR 351,717	116,053	14,234	NR NR 54,900 14,647 NR	16,013 234,850 26,688 64,050 21,350 10,675	17,306,707	NR	NR	NR	ported
TRI	NN NN NN NN NN NN NN NN NN NN NN NN NN	50,000 940,000	NR 360, 000 NR NR NR	45,600 38,475 42,000	15,000 880,000 432,000	75,000	1,575,000 63,500 NR	NR	NR	21,000 31,000 NR NR 28,000	NN NN NN NN NN NN NN NN NN	NR	264,900	10,235	18,304	NR = Not Rep
Chemical	TOLUENE SILVER SILVER MERCURY LEAD COPPER COPPER CORALT CORALT CORALT CADMIUM BARIUM ALUMINUM (FUME OR DUST)	TETRACHLOROETHYLENE SODIUM SULFATE (SOLUTION)	TOLUENE SODIUM HYDROXIDE SELENIUM MERCURY COPPER	1,1,1-TRICHLOROETHANE SODIUM HYDROXIDE MELAMINE	SULFURIC ACID SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	SODIUM SULFATE (SOLUTION)	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE PHENOL	COBALT	PHENOL	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE SILVER COPPER AMMONIUM SULFATE (SOLUTION)	TOLUENE SILVER SELENIUM MERCURY LEAD CADMIUM	PHENOL	SODIUM SULFATE (SOLUTION)	SODIUM HYDROXIDE	XYLENE	puantities reported in lbs.
SIC	66666666666666666666666666666666666666	2269 2269	2269 2269 2269 2269 2269	2269 2269 2269	2269 2269 2269	2269	2269 2269 2269	2269	2269	2269 2269 2269 2269 2269	2269 2269 2269 2269 2269 2269 2269	2269	2269	2269	2272	All q

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07/22	/93 Facilities With	Significant	Discrepancies	Between TRI	and Pretreatment	in 19
SIC	Chemical		TRI	Pretreat	Difference	
2272 2272 2272 2272	1,2,4-TRICHLOROBENZENE SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE ETHYLENE GLYCOL		29,767 196,605 105,810 11,673	N N N N N N N N N N N N N N N N N N N	29,767 196,605 105,810 11,673	•
2281 2281 2281	TOLUENE PHENOL MERCURY		NR 2, 7, NR 2, 7, NR	11,895 45,000,000 2, 14,640	11,895 745,000,000 14,640	
2281 2281	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE		669,300 18,690	NR NR	669,300 18,690	
2281	SODIUM SULFATE (SOLUTION)		170,000	NR	170,000	
2281 2281 2281 2281 2281	TOLUENE SELENIUM MERCURY LEAD COPPER		K K K K K	13,366 13,366 35,643 28,960 13,366	13,366 13,366 35,643 28,960 13,366	
2281	SODIUM HYDROXIDE		70,000	NR	70,000	
2282 2282 2282	SULFURIC ACID SODIUM HYDROXIDE ETHYLENE GLYCOL		12,100 34,000 19,200	NR NR NR	12,100 34,000 19,200	
2283	SODIUM SULFATE (SOLUTION)		180,000	NR	180,000	
2283	SODIUM SULFATE (SOLUTION)		11,400	NR	11,400	
2284 2284	METHANOL FORMALDEHYDE		209,400 62,800	NR NR	209,400 62,800	
2293 2293 2293	SULFURIC ACID SODIUM HYDROXIDE CHLORINE		62,479 1,858,724 330,000	NN NN NN	62,479 1,858,724 330,000	
2295 2295	SULFURIC ACID SODIUM HYDROXIDE		59,800 55,500	NR NR	59,800 55,500	
2296	SULFURIC ACID		13,256	NR	13,256	
2297 2297 2297	TOLUENE LEAD COPPER		NR NR NR	58,403 14,601 58,403	58,403 14,601 58,403	
2298	SODIUM SULFATE (SOLUTION)		542,975	NR	542,975	
2299 2299	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE		1,779,000 232,000	NR NR	1,779,000 232,000	
2300	SODIUM HYDROXIDE		201,719	NR	201,719	
2300	SODIUM HYDROXIDE		189,000	NR	189,000	
2321 2321	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE		620,000 55,000	NR NR	620,000 55,000	
2321	PHENOL		NR	85,044	85,044	

\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

All quantities reported in lbs.

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Difference	120,032 578,336 87,296 87,296 52,736 65,472 10,912 10,912	176,000	235,000	760,500 23,200	97,750 28,809	12,161	18,203	16,475	20,000	12,894	55,484	12,919 * 62,000	43,736	17,000	23,000	206,000 10,800 29,200	23,900	280,000 54,000 13,000 21,000	81,700	21,960	67,500 13,807	12,554	** = Values
Pretreat	120, 032 578, 336 87, 296 87, 296 32, 736 52, 736 52, 472 10, 912 10, 912	NR	NR	NR NR	NR NR	12,161	NR	16,475	NR	12,894	NR	81 NR	NR	NR	NR	NR NR NR	NR	NR NR NR NR	NR	21,960	NR NR	NR	orted
TRI	n n n n n n n n n n n n n n n n n n n	176,000	235,000	760,500 23,200	97,750 28,809	NR	18,203	NR	20,000	NR	55,484	13,000 62,000	43,736	17,000	23,000	206,000 10,800 29,200	23,900	280,000 54,000 13,000 21,000	81,700	NR	67,500 13,807	12,554	NR = Not Rep
Chemical	TOLUENE SILVER SILVER SELENIUM MERCURY MERCURY COPPER COPPER COPPER COBALT CADMIUM (FUME OR DUST)	SODIUM SULFATE (SOLUTION)	SODIUM SULFATE (SOLUTION)	SODIUM SULFATE (SOLUTION) BIPHENYL	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	CHLORINE	n-BUTYL ALCOHOL	MERCURY	TOLUENE	PHENOL	ISOPROPYL ALCOHOL	TOLUENE METHANOL	NITRIC ACID	AMMONIA	SODIUM HYDROXIDE	SODIUM SULFATE (SOLUTION) METHANOL ETHYLENE GLYCOL	ETHYLENE GLYCOL	SODIUM SULFATE (SOLUTION) METHYL TERT-BUTYL ETHER ETHYLENE GLYCOL ACETONE	ETHYLENE OXIDE	MERCURY	METHANOL DIMETHYL PHTHALATE	SODIUM HYDROXIDE	puantities reported in lbs.
SIC	232222222222222222222222222222222222222	2322	2322	2322 2322	2341 2341	2439	2511	2511	2641	2653	2741	2754 2754	2793	2819	2819	2821 2821 2821	2821	2833 2833 2833 2833 2833	2834	2834	2834 2834	2843	All c

07/22	://93 Facilities With Significa	nt Discrepanci	es Between TRI	and Pretreatment in 1987
SIC	Chemical	TRI	Pretreat	Difference
2843	SODIUM SULFATE (SOLUTION)	419,000	NR	419,000
2843 2843	SULFURIC ACID SODIUM HYDROXIDE	31,000 12,900	NR NR	31,000 12,900
2865	AMMONIA	17,000	NR	17,000
2869	SODIUM SULFATE (SOLUTION)	2,055,470	NR	2,055,470
2869	SODIUM SULFATE (SOLUTION)	137,500	NR	137,500
2899	SODIUM SULFATE (SOLUTION)	15,330	NR	15,330
3052 3052	TOLUENE SODIUM SULFATE (SOLUTION)	11,600 230,000	90,774 NR	79,174 ** 230,000
3069	MERCURY	NR	20,714	20,714
3231	SODIUM HYDROXIDE	10,590	NR	10,590
329	ASBESTOS	140,000	NR	140,000
3321 3321	TOLUENE	NR NR	74,291 134,633	74,291 134,633
3339	CYANIDE COMPOUNDS	NR	497,394	497,394
3351	METHANOL	10,000	NR	10,000
3353	PHENOL	NR	47,742	47,742
3354 3354 3354	PHENOL COPPER ARSENIC	NR NR NR	319,170,240 13,654 234,725	319,170,240 13,654 234,725
3411	SILVER	NR	47,565	47,565
3412	SODIUM HYDROXIDE	58,500	NR	58,500
3423 3423 3423	SULFURIC ACID SODIUM HYDROXIDE HYDROCHLORIC ACID	163,000 242,000 225,000	NR NR NR	163,000 242,000 225,000
3425 3425 3425	TOLUENE MERCURY COPPER	NR NR NR	10,120 10,795 62,072	10,120 10,795 62,072
3429	SODIUM SULFATE (SOLUTION)	110,000	NR	110,000
3429	SODIUM SULFATE (SOLUTION)	205,900	NR	205,900
3444	ARSENIC	NR	31,959	31,959
3448 3448 3448	TOLUENE SELENIUM LEAD	0 NR NR	13,517 26,282 16,520	13,517 ** 26,282 16,520
3471 3471	SELENIUM COBALT	NR NR	132,844 1,195,592	132,844 1,195,592
Allo	quantities reported in lbs.	NR = Not Rep	ported	<pre>** = Values reported to both TRI and Pretreatment</pre>

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/22,	793 Facilities With Significant	Discrepancies	Between TRI	and Pretreatment in 1987
υÏ	Chemical	TRI	Pretreat	Difference
71	SODIUM SULFATE (SOLUTION)	57,000	NR	57,000
71	SULFURIC ACID	65,000	NR	65,000
11	SODIUM HYDROXIDE NITRIC ACID	<b>4</b> 7,000 12,000	NR NR	47,000
79	HYDROCHLORIC ACID	27,000	NR	27,000
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SIC	Chemical	TRI	Pretreat	Difference	
3471	SODIUM SULFATE (SOLUTION)	57,000	NR	57,000	
3471 3471 3471	SULFURIC ACID SODIUM HYDROXIDE NITRIC ACID	65,000 47,000 12,000	NR NR NR	65,000 47,000 12,000	
3479	HYDROCHLORIC ACID	27,000	NR	27,000	
3479	SODIUM HYDROXIDE	112,000	NR	112,000	
3499	SULFURIC ACID	33,680	NR	33,680	
3519 3519 3519 3519 3519	TOLUENE MERCURY COPPER COBALT ARSENIC ALUMINUM (FUME OR DUST)	NN NN NN NN NN O	58,772 16,456 423,157 23,509 2,565,977 2,262,715	58,772 16,456 423,157 23,509 2,565,977 2,262,715 **	
3536	SILVER	NR	89,213	89,213	
3599	ALUMINUM OXIDE	12,000	NR	12,000	
3613 3613 3613 3613	TOLUENE MERCURY COPPER	NR NR	52,342 11,632 13,958	52,342 11,632 13,958	
3621 3621	TOLUENE MERCURY	NR NR	45,857 45,857	45,857 45,857	
3621 3621 3621 3621 3621	TOLUENE SILVER SELENIUM MERCURY	NN NN NN NN	42,700 109,800 30,500 36,600	42,700 109,800 30,500 36,600	
3622	GLYCOL ETHERS	21,000	NR	21,000	
3622 3622 3622	SULFURIC ACID SODIUM HYDROXIDE HYDROCHLORIC ACID	59,517 26,054 25,374	NR NR NR	59,517 26,054 25,374	
3672	MERCURY	NR	10,614	10,614	
3674 3674 3674	SULFURIC ACID SODIUM HYDROXIDE PHOSPHORIC ACID	76,083 33,600 11,228	NR NR NR	76,083 33,600 11,228	
3674 3674 3674	SULFURIC ACID SODIUM HYDROXIDE HYDROCHLORIC ACID	11,718 74,355 85,010	NR NR NR	11,718 74,355 85,010	
3675	ETHYLENE GLYCOL	33,500	NR	33,500	
3676	LEAD	NR	40,275	40,275	
3678	SODIUM SULFATE (SOLUTION)	65,000	NR	65,000	
3691	SODIUM SULFATE (SOLUTION)	163,000	NR	163,000	
All c	quantities reported in lbs.	NR = Not Repo	orted	** = Values reported to bot	r,

TRI and Pretreatment

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SIC	Chemical	TRI	Pretreat	Difference	
3695	TOLUENE	NR	22,326	22,326	
3700	SODIUM SULFATE (SOLUTION)	183,251	NR	183,251	•
3714 3714 3714 3714 3714	TOLUENE SILVER SELENIUM MERCURY	K K K K K	383,690 1,899,266 115,107 57,554 172,661	383,690 1,899,266 115,107 57,554 172,661	
3714 3714 3714 3714	COFFER COBALT CADMIUM ALUMINUM (FUME OR DUST)	NN NN NN NN	57,554 19,185 19,185 19,185	57,554 19,185 19,185	
3714	SODIUM HYDROXIDE	152,530	NR	152,530	
3714 3714	SULFURIC ACID SODIUM HYDROXIDE	94,500 150,500	NR NR	94,500 150,500	
3714 3714	SULFURIC ACID SODIUM HYDROXIDE	88,200 80,200	NR NR	88,200 80,200	
3964	XYLENE	30,000	NR	30,000	
3999	SODIUM HYDROXIDE	22,000	NR	22,000	
8608 8098	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	500,000 100,000	NR NR	500,000 100,000	

All quantities reported in lbs.

NR = Not Reported

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07/22	/93 Facilities With Significant	Discrepancies	Between TRI	and Pretreatment in 1988
SIC	Chemical	TRI	Pretreat	Difference
20 20	SULFURIC ACID SODIUM HYDROXIDE	80,700 448,851	NR NR	80,700 448,851
20	SODIUM HYDROXIDE	93,000	NR	93,000
2016	PHENOL	NR	16,144	16,144
2024 2024	SODIUM HYDROXIDE AMMONIA	89,254 82,560	NR NR	89, 254 82, 560
202 <b>4</b> 2024	SODIUM HYDROXIDE PHOSPHORIC ACID	148,365 48,364	NR NR	148,365 48,364
2026 2026	SODIUM HYDROXIDE PHOSPHORIC ACID	18,846 18,674	NR NR	18, 846 18, 674
2026	PHOSPHORIC ACID	11,000	NR	11,000
2026	TOLUENE	NR	64,604	64,604
2032	PHENOL	NR	606,762	606,762
2046	SODIUM HYDROXIDE	1,200,000	NR	1,200,000
2077	AMMONIA	90,000	NR	<b>90</b> ,000
2077	AMMONIA	67,800	NR	67,800
2086	SODIUM HYDROXIDE	45,674	NR	45,674
2086	SODIUM HYDROXIDE	37,363	NR	37,363
2111	TOLUENE	0	40,375	40,375 **
2111	SODIUM HYDROXIDE	13,000	NR	13,000
22	SODIUM HYDROXIDE	920,000	NR	920,000
22 22	SULFURIC ACID SODIUM HYDROXIDE	675,3 <b>4</b> 2 757, <b>4</b> 70	NR NR	675,342 757,470
2200	SODIUM HYDROXIDE	527,500	NR	527,500
2200	SODIUM HYDROXIDE	1,105,558	NR	1,105,558
2211	SODIUM HYDROXIDE	11,250	NR	11,250
2221 2221 2221	SULFURIC ACID SODIUM HYDROXIDE AMMONIA	350,000 711,000 31,500	NR NR NR	350,000 711,000 31,500
2221	SODIUM HYDROXIDE	101,150	NR	101,150
2221	SODIUM HYDROXIDE	27,500	NR	27,500
2221	SODIUM HYDROXIDE	125,000	NR	125,000
2221	SODIUM HYDROXIDE	32,650	NR	32,650
2221	PHENOL	NR	42,122,025	42,122,025
All c	quantities reported in lbs.	NR = Not Repor	ted	<pre>** = Values reported to both TRI and Pretreatment</pre>

07/22	793 Facilities With Significant	Discrepancies	Between TRI a	and Pretreatment in 1988
SIC	Chemical	TRI	Pretreat	Difference
2221	PHENOL	NR	37,028	37,028 .
2221	SODIUM HYDROXIDE	115,065	NR	115,065
2221	TETRACHLOROETHYLENE	102,600	NR	102,600
2231 2231	SILVER ARSENIC	NR NR	11,712 223,769	11,712 223,769
2251 2251	SODIUM HYDROXIDE AMMONIUM SULFATE (SOLUTION)	10,260 10,800	NR NR	10,260 10,800
2251	CHLORINE	231,000	NR	231,000
2251 2251	TOLUENE MERCURY	NR NR	51,806 10,361	51,806 10,361
2251	AMMONIUM SULFATE (SOLUTION)	60,000	NR	60,000
2252 2252 2252	SODIUM HYDROXIDE CHLORINE AMMONIUM SULFATE (SOLUTION)	12,000 56,560 29,500	NR NR NR	12,000 56,560 29,500
2252	CHLORINE	252,670	NR	252,670
2252	CHLORINE	46,300	NR	46,300
2252 2252	MERCURY ARSENIC	NR NR	10,190 10,190	10,190 10,190
2252	BIPHENVL	26,995	NR	26,995
2252 2252	SODIUM HYDROXIDE AMMONIUM SULFATE (SOLUTION)	12,587 17,000	NR NR	12,587 17,000
2252	CHLORINE	33,000	NR	33,000
52555 525555 52555	TOLUENE SILVER SELENIUM PHENOL MERCURY PHENOL	NN NN NN NN NN NN	31,034 10,345 12,414 3,055,841 22,758 10,201	31,034 10,345 12,414 3,055,841 10,201
2252	AMMONIUM SULFATE (SOLUTION)	14,062	NR	14,062
2252	CHLORINE	25,063	NR	25,063
2253	PHENOL	NR	59,870	59,870
2253 2253	SULFURIC ACID SODIUM HYDROXIDE	120,000 240,000	NR NR	120,000 240,000
2253 2253 2253 2253	XYLENE SODIUM HYDROXIDE CHLORINE BIPHENYL	97,744 11,700 73,000 15,883	NR NR NR NR	97,744 11,700 73,000 15,883
2257	SODIUM HYDROXIDE	58,825	NR	58,825
All (	quantities reported in lbs.	NR = Not Repo	rted	<pre>** = Values reported to both TRI and Pretreatment</pre>

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07/22	/93 Facilities Wi	ith Significant Div	screpancies Be	etween TRI	and Pretreatment in 1988	
SIC	Chemical		TRI	Pretreat	Difference	
2257	SILVER		NR	13,491	13,491	
2257	AMMONIUM SULFATE (SOLU	LION)	10,000	NR	10,000	
2257	SODIUM HYDROXIDE		47,500	NR	47,500	
2257 2257 2257 2257	SODIUM HYDROXIDE PSEUDOCUMENE FORMALDEHYDE BIPHENYL AMMONIA		144,800 33,220 144,611 90,775 71,545	NN NN NN NN	144,800 33,220 144,611 90,775 71,545	
2258 2258 2258 2258 2258 2258 2258	XYLENE TETRACHLOROETHYLENE SODIUM HYDROXIDE ETHYLENE GLYCOL BIPHENYL BMMONIUM SULFATE (SOLU'	(NOI	27,265 61,864 49,006 42,071 18,273 24,048	NN NN NN NN NN NN	27,265 61,864 49,006 42,071 18,273 24,048	
2258	AMMONIUM SULFATE (SOLU'	LION)	19,950	NR	19,950	
2258 2258 2258 2258	TOLUENE SODIUM HYDROXIDE SILVER AMMONIUM SULFATE (SOLU'	(ION)	18,006 29,885 NR 63,285	NR NR 89, 222 NR	18,006 29,885 89,222 63,285	
2258 2258	SODIUM HYDROXIDE AMMONIUM SULFATE (SOLU	TION)	26,600 152,000	NR NR	26,600 152,000	
226	SULFURIC ACID		271,000	NR	271,000	
2260	SILVER		NR	25,639	25,639	
2261	SODIUM HYDROXIDE		213,338	NR	213,338	
2261	SODIUM HYDROXIDE		59,700	NR	59,700	
2262	SODIUM HYDROXIDE		13,660	NR	13,660	
2262 2262 2262 2262	XYLENE 1,2,4-TRICHLOROBENZENE AMMONIUM SULFATE (SOLU AMMONIA	TION)	88,496 10,201 22,994 28,295	NR NR NR	88,496 10,201 22,994 28,295	
2262 2262	PHENOL MERCURY		NR NR	14,903 13,334	14,903 13,334	
2262	SODIUM HYDROXIDE		91,397	NR	91,397	
2262	SODIUM HYDROXIDE		280,000	NR	280,000	
2262	AMMONIUM SULFATE (SOLU	TION)	57,600	NR	57,600	
2269 2269	TETRACHLOROETHYLENE SODIUM HYDROXIDE		13,856 23,663	NR NR	13,856 23,663	

\*\* = Values reported to both TRI and Pretreatment

58,000 12,000

> NR NR

58,000 12,000 NR = Not Reported

All quantities reported in lbs.

2269 TETRACHLOROETHYLENE 2269 BIPHENYL

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07/22	/93 Facilities With Significan	t Discrepancie	s Between TRI	and Pretreatme	nt in 1988
SIC	Chemical	TRI	Pretreat	Difference	
2269	SODIUM HYDROXIDE	34,500	NR	34,500	
2269	AMMONIUM SULFATE (SOLUTION)	11,000	NR	11,000	•
2269 2269 2269 2269	1,2,4-TRICHLOROBENZENE TETRACHLOROETHYLENE SODIUM HYDROXIDE PHENOL	14,848 10,150 335,000 NR	NR NR NR 247,297	14,848 10,150 335,000 247,297	
2269	SODIUM HYDROXIDE	36,500	NR	36,500	
2269 2269	SODIUM HYDROXIDE PHENOL	290,000 NR	NR 14,870	290,000 14,870	
2269	GLYCOL ETHERS	15,200	NR	15,200	
2269	PHENOL	NR	14,889	14,889	
2269	SODIUM HYDROXIDE	11,650	NR	11,650	
2269	SODIUM HYDROXIDE	26,600	NR	26,600	
2269 2269 2269 2269	1,2,4-TRICHLOROBENZENE TETRACHLOROETHYLENE SODIUM HYDROXIDE PSEUDOCUMENE BIPHENYL	59,704 42,074 82,200 23,496 13,256	N N N N N N N N N N N N N N N N N N N	59,704 42,074 82,200 23,496 13,256	
2272 2272 2272 2272	1,2,4-TRICHLOROBENZENE SODIUM HYDROXIDE PHOSPHORIC ACID ETHYLENE GLYCOL	12,952 114,908 15,232 16,747	NR NR NR	12,952 114,908 15,232 16,747	
2281	SODIUM HYDROXIDE	18,957	NR	18,957	
2281	PHENOL	NR	6,717,625	6,717,625	
2281	SODIUM HYDROXIDE	150,000	NR	150,000	
2282 2282	SULFURIC ACID SODIUM HYDROXIDE	20,564 157,971	NR NR	20,564 157,971	
2282	1,2,4-TRICHLOROBENZENE	59,922	NR	59,922	
2284 2284	METHANOL FORMALDEHYDE	226,000 117,900	NR NR	226,000 117,900	
2293 2293 2293	SULFURIC ACID SODIUM HYDROXIDE CHLORINE	79,842 1,857,734 332,000	NR NR NR	79,842 1,857,734 332,000	
2296	SULFURIC ACID	13,100	NR	13,100	
2299	SODIUM HYDROXIDE	348,070	NR	348,070	
2300	SODIUM HYDROXIDE	305,680	NR	305,680	
2321	SODIUM HYDROXIDE	82,250	NR	82,250	
2321	PHENOL	NR	78,927	78,927	
All ç	nuantities reported in lbs.	NR = Not Repo	rted	** = Values	reported to both TRI and Pretreatment

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07/22	:/93 Facilities With Significar	t Discrepancies	Between TRI	and Pretreatment in 1988	
SIC	Chemical	TRI	Pretreat	Difference	
2322	SODIUM HYDROXIDE	183,144	NR	183,144	
2322 2322	SODIUM HYDROXIDE BIPHENYL	170,023 28,667	NR NR	170,023 28,667	
2322	SODIUM HYDROXIDE	86,688	NR	86,688	
2341	SODIUM HYDROXIDE	26,606	NR	26,606	
2439	CHLORINE	NR	26,873	26,873	
2641	TOLUENE	14,200	NR	14,200	
2653	PHENOL	NR	11,011	11,011	
2754 2754	TOLUENE METHANOL	13,886 14,674	56 NR	13,830 ** 14,674	
2793	NITRIC ACID	27,300	NR	27,300	
2819	AMMONIA	31,455	NR	31,455	
2819	SODIUM HYDROXIDE	22,000	NR	22,000	
2821 2821 2821	METHANOL FORMALDEHYDE ETHYLENE GLYCOL	50,128 48,465 152,388	NR NR NR	50,128 48,465 152,388	
2824	ACETONE	22,609	NR	22,609	
2833 2833 2833 2833 2833 2833	TOLUENE METHYLENE CHLORIDE CHLOROFORM ACETONE	68,000 12,000 65,000 170,000	309 NR NR NR	67,691 ** 12,000 65,000 170,000	
2834 2834	ETHYLENE ØXIDE ETHYLENE GLYCOL	25,000 100,000	NR NR	25,000 100,000	
2834 2834 2834 2834 2834	SODIUM HYDROXIDE METHANOL DIMETHYL PHTHALATE n-BUTYL ALCOHOL	22,602 19,772 10,229 12,608	<u> </u>	22,602 19,772 10,229 12,608	
2834	GLYCOL ETHERS	27,000	NR	27,000	
2843	PHENOL	NR	61,268	61,268	
2844 2844	SULFURIC ACID SODIUM HYDROXIDE	25,517 71,525	NR NR	25,517 71,525	
2844	GLYCOL ETHERS	56,000	NR	56,000	
2865	AMMONIA	16,000	NR	16,000	

\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

All quantities reported in lbs.

SODIUM HYDROXIDE n-BUTYL ALCOHOL

2869 2869

2869 ETHYLENE GLYCOL

54,495 19,042 14,887

NR NR

54,495 19,042 14,887

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SIC	Chemical	TRI	Pretreat	Difference	
3052 3052	ZINC COMPOUNDS	26,839 8,900	NR 28,212	26,839 19,312	:
3231	SODIUM HYDROXIDE	12,238	NR	12,238	
3321	PHENOL	NR	2,948,806	2,948,806	
3354	SODIUM HYDROXIDE	148,000	NR	148,000	
3354 3354	PHENOL ARSENIC	NR NR	1,323,022,282 71,039	1,323,022,282 71,039	
3423 3423 3423	SULFURIC ACID SODIUM HYDROXIDE HYDROCHLORIC ACID	155,000 230,000 225,000	NR NR NR	155,000 230,000 225,000	
3425	PHENOL	NR	16,145,938	16,145,938	
3471 3471	SULFURIC ACID SODIUM HYDROXIDE	14,632 14,632	NR NR	14,632 14,632	
3471	SODIUM HYDROXIDE	18,000	NR	18,000	
3471 3471 3471	SULFURIC ACID SODIUM HYDROXIDE NITRIC ACID	36,283 31,452 10,334	NR NR NR	36,283 31,452 10,334	
3499 3499	SULFURIC ACID SODIUM HYDROXIDE	60,749 33,191	NR NR	60,7 <b>4</b> 9 33,191	
3519 3519	SULFURIC ACID DIETHANOLAMINE	12,000 12,000	NR NR	12,000	
3519	ARSENIC	NR	135,974	135,974	
3531	DIETHANOLAMINE	30,607	NR	30,607	
3622	GLYCOL ETHERS	14,732	NR	14,732	
3622 3622 3622 3622	SULFURIC ACID SODIUM HYDROXIDE HYDROCHLORIC ACID	65,179 37,847 41,901	NR NR NR	65,179 37,847 41,901	
3661	GLYCOL ETHERS	15,785	NR	15,785	
3674 3674	PHOSPHORIC ACID AMMONIA	10,104 11,154	NR NR	10,104 11,154	
3674 3674 3674	SULFURIC ACID SODIUM HYDROXIDE HYDROCHLORIC ACID	12,349 31,086 67,663	NR NR NR	12,349 31,086 67,663	
3675	ETHYLENE GLYCOL	42,783	NR	42,783	
3676	PHENOL	NR	2,812,890	2,812,890	
3713	PHOSPHORIC ACID	16,000	NR	16,000	
3714	SODIUM HYDROXIDE	390,267	NR	390,267	

\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

All quantities reported in lbs.

SIC	Chemical	TRI	Pretreat	Difference
3714 3714 3714	SULFURIC ACID SODIUM HYDROXIDE	26,000	NR NR	126,000 228,000
3714 3714	SULFURIC ACID SODIUM HYDROXIDE	38,200 30,200	NR NR	88,200 80,200
3963	ALUMINUM OXIDE	16,000	NR	16,000
3999	SODIUM HYDROXIDE	10,000	NR	10,000
3999	SODIUM HYDROXIDE	15,750	NR	15,750

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07/22	/93 Facilities With Significant	Discrepancie	s Between TRI	and Pretreatment in 1989
SIC	Chemical	TRI	Pretreat	Difference
20 20	SULFURIC ACID CHLORINE	166,358 15,295	NR NR	166, 358 15, 295
20	AMMONIA	257,500	NR	257,500
2016 2016	TOLUENE CYANIDE COMPOUNDS	NR NR	660,428 8,521,649	660,428 8,521,649
2021	PHOSPHORIC ACID	11,200	NR	11,200
2024	PHOSPHORIC ACID	47,560	NR	47,560
2026	PHOSPHORIC ACID	12,000	NR	12,000
2026	PHENOL	NR	7,840,475	7,840,475
2070	PHOSPHORIC ACID	13,700	NR	13,700
2086	PHOSPHORIC ACID	10,660	NR	10,660
22 22	XYLENE TETRACHLOROETHYLENE 1,2-DICHLOROBENZENE	62,775 20,925 52,313	NR NR NR	62, 775 20, 925 52, 313
2200	ARSENIC	NR	80,057	80,057
2200	TOLUENE	NR	25,206	25,206
2200	AMMONIUM SULFATE (SOLUTION)	17,600	NR	17,600
2205	CHROMIUM	NR	11,311	11,311
2211	ARSENIC	NR	39,166	39,166
2211	PHENOL	NR	20,053,549	20,053,549
2221 2221 2221 2221 2221	TETRACHLOROETHYLENE SULFURIC ACID METHANOL ETHYLENE GLYCOL AMMONIA	12,600 16,500 81,164 53,200 26,575	N N N N N N N N N N N N N N N N N N N	12,600 16,500 81,164 53,200 26,575
2221	TETRACHLOROETHYLENE	44,000	NR	44,000
2228	SULFURIC ACID	14,000	NR	14,000
2241 2241 2241 2241	TOLUENE SELENIUM MERCURY COPPER	NR NR NR	30,053 10,018 43,410 46,749	30,053 10,018 43,410 46,749
2251	AMMONIUM SULFATE (SOLUTION)	57,800	NR	57,800
2252 2252	CHLORINE AMMONIUM SULFATE (SOLUTION)	78,000 45,000	NR NR	78,000 45,000
2252	CHLORINE	11,250	NR	11,250
2252 2252	CHLORINE AMMONIUM SULFATE (SOLUTION)	323,184 12,000	NR NR	323,184 12,000
All c	nuantities reported in lbs.	NR = Not Repo	rted	<pre>** = Values reported to both TRI and Pretreatment</pre>

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and Pretreatment	Difference	55,192	11,628	10,200	39.500

SIC	Chemical	TRI	Pretreat	Difference	
2252	CHLORINE	55,192	NR	55,192	
2252	AMMONIUM SULFATE (SOLUTION)	11,628	NR	11,628	
2252	XYLENE	10,200	NR	10,200	
2252	AMMONIUM SULFATE (SOLUTION)	39,500	NR	39,500	
2252	CHLORINE	67,732	NR	67,732	
2252	AMMONIUM SULFATE (SOLUTION)	23,527	NR	23,527	
2252	CHLORINE	21,160	NR	21,160	
2252	CHLORINE	20,000	NR	20,000	
2253	PHENOL	NR	63,993	63,993	
2253	SILVER	NR	14,738	14,738	
2253	ETHYLENE GLYCOL	10,000	NR	10,000	
2257 2257 2257 2257	GLYCOL ETHERS FORMALDEHYDE 1,2-DICHLOROBENZENE AMMONIA	100,000 177,820 150,327	N N N N N N N N N N N N N N N N N N N	100,000 77,820 100,000 15,327	
22588 22588 22588 22588 22588 25288 25288 25258 25258 25258 25258 25258 25258 25258 25258 25258 25258 25258 25258 25258 25258 25258 2558 2558 2558 2558 2558 2558 2558 2558 2558 2558 2558 25557 25557	XYLENE TETRACHLOROETHYLENE SILVER ETHYLENE GLYCOL BIPHENYL AMMONIUM SULFATE (SOLUTION)	19,325 43,737 43,437 43,488 13,488 14,389	NR NR 38, 406 NR NR NR	19,325 43,737 38,406 43,488 13,430 14,389	
2258	AMMONIUM SULFATE (SOLUTION)	25,850	NR	25,850	
2258 2258	SULFURIC ACID HYDROCHLORIC ACID	48,183 12,865	NR NR	<b>4</b> 8,183 12,865	
2258 2258 2258	TETRACHLOROETHYLENE PSEUDOCUMENE AMMONIUM SULFATE (SOLUTION)	21,030 27,631 57,824	NR NR NR	21,030 27,631 57,824	
2258	AMMONIUM SULFATE (SOLUTION)	184,800	NR	184,800	
226	SULFURIC ACID	271,000	NR	271,000	
2261 2261	GLYCOL ETHERS ETHYLENE GLYCOL	38,660 55,294	NR NR	38,660 55,294	
2262 2262	XYLENE AMMONIA	131,088 30,351	NR NR	131,088 30,351	
2262	PHENOL	NR	21,941	21,941	
2262	1,1,1-TRICHLOROETHANE	18,000	NR	18,000	
2262	AMMONIUM SULFATE (SOLUTION)	48,000	NR	48,000	
All	quantities reported in lbs.	NR = Not Repo	orted	** = Values repo	rted to both TRI and Pretreatment

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Between TRI and Pretreatment in 1989	Pretreat Difference	62,357 62,357	NR 14,243	NR 12,000	NR 33,000 405,488 405,488	NR 125,000 140,898 140,898	18, 321 18, 321	14,644 14,644	NR 11,854 NR 36,728 NR 19,949 NR 17,432	NR 12,000 113,266,576 313,266,576 NR 17,600	NR 56,047	NR 248,000 NR 60,000 165,870 165,870	10,195 10,195	NR 55,800	NR 46,640	13,465 13,465
lficant Discrepancies	TRI	NR	14,243	12,000	33,000 NR	125,000 NR	NR	NR	11,854 36,728 19,949 17,432	12,000 NR 3 17,600	56,047	248,000 60,000 NR	NR	55,800	46,640	NR
/93 Facilities With Sign	Chemical	TOLUENE	<b>TETRACHLOROETHYLENE</b>	BIPHENYL	1,2,4-TRICHLOROBENZENE PHENOL	FLUOMETURON ARSENIC	PHENOL	PHENOL	1,2,4-TRICHLOROBENZENE TETRACHLOROETHYLENE PSEUDOCUMENE BIPHENYL	SULFURIC ACID PHENOL ETHYLENE GLYCOL	1,2,4-TRICHLOROBENZENE	METHANOL FORMALDEHYDE CHLORINE	PHENOL	CHLORINE	PHOSPHORIC ACID	TOLUENE
07/22/	SIC	2262	2269	2269	2269 2269	2269 2269	2269	2269	2269 2269 2269 2269	2282 2282 2282	2282	2284 2284 2284	2293	2300	2321	2321

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\*\* = Values reported to both TRI and Pretreatment

16,500 58,750 43,500

NR NR

16,500 58,750 43,500

169,494 114,607 171,779 43,527

169,494 114,607 171,779 43,527

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TOLUENE MERCURY LEAD COPPER

2514 2514 2514 2514 219,335 959,204

219,335 959,204 23,100 20,100

NR NR

23,100 20,100

NITRIC ACID

2793 2819

SILVER BARIUM

2641 2641 AMMONIA

NR = Not Reported

All quantities reported in lbs.

METHANOL FORMALDEHYDE ETHYLENE GLYCOL

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and Pretreatment	Difference	10,502 17,723
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SIC	Chemical	TRI	Pretreat	Difference	
2821 2821	TOLUENE ALUMINUM (FUME OR DUST)	NR NR	10,502 17,723	10,502 17,723	
2821	FORMALDEHYDE	10,575	NR	10,575	
2821	ETHYLENE GLYCOL	13,762	NR	13,762	
2821	AMMONIA	25,000	NR	25,000	
2833 2833 2833 2833 2833	TOLUENE METHYLENE CHLORIDE CHLOROFORM ACETONE	41,000 13,000 52,000 140,000	10, 559 NR NR NR	30,441 ** 13,000 52,000 140,000	
2834 2834	ETHYLENE OXIDE ETHYLENE GLYCOL	18,000 76,000	NR NR	18,000 76,000	
2834	GLYCOL ETHERS	54,000	NR	54,000	
2841	TOLUENE	NR	3,878,378	3,878,378	
2844 2844	TOLUENE MERCURY	NR NR	132,097 22,355	132,097 22,355	
2844	GLYCOL ETHERS	125,000	NR	125,000	
2865	AMMONIA	15,000	NR	15,000	
2869	ETHYLENE GLYCOL	42,310	NR	42,310	
2869 2869 2869 2869	TOLUENE METHANOL FORMALDEHYDE DIMETHYL PHTHALATE	17,163 15,919 12,340 90,712	22 NR NR NR	17,141 ** 15,919 12,340 90,712	•
3052 3052	ZINC COMPOUNDS TOLUENE	22,423 2,703	NR 18,536	22,423 15,833 **	
3069	METHANOL	16,600	NR	16,600	
3281	PHENOL	NR	16,592	16,592	
329	PHENOL	38,200	NR	38,200	
3321	PHENOL	NR	5,606,623	5,606,623	
3354	PHENOL	NR	700,444	700,444	
3356	MANGANESE AND COMPOUNDS	NR	21,781	21,781	
3425	PHENOL	NR	26,980,692	26,980,692	
3444	ARSENIC	NR	10,331	10,331	
3471 3471 3471 3471 3471	TOLUENE MERCURY LEAD SELENIUM MERCURY	NR NR NR NR	2,208,580 1,099,699 409,861 438,807 487,337	2,208,580 1,099,699 409,861 487,337 487,337	
All o	quantities reported in lbs.	NR = Not Rep	orted	<pre>• = Values reported to both TRI and Pretreat.</pre>	ent

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SIC	Chemical	TRI	Pretreat	Difference
3471 3471 3471	SULFURIC ACID NITRIC ACID		NR NR	18,513 14,058
3499	SULFURIC ACID	88,257	NR	88,257
3553	ETHYLENE GLYCOL	10,700	NR	10,700
3613	PHENOL	NR	4,605,043	4,605,043
3622 3622	SILVER GLYCOL ETHERS	NR 21,500	129,797 NR	129,797 21,500
3646	ETHYLENE GLYCOL	83,000	NR	83,000
3672	PHENOL	NR	140,504	140,504
3675	ETHYLENE GLYCOL	19,226	NR	19,226
3676	PHENOL	NR	348,081	348,081
3692	PHENOL	NR	12,485,779	12,485,779
3713	PHOSPHORIC ACID	19,000	NR	<b>19,00</b> 0
3714	TOLUENE	NR	10,093	10,093
3714	PHENOL	NR	45,334,578	45,334,578
3732	TOLUENE	0	10,763	10,763 **
3964	XYLENE	27,000	NR	27,000
3999	ARSENIC	NR	16,894	16,894

\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

07/22	<pre>/93 Facilities With Significant</pre>	Discrepancies	Between TRI	and Pretreatment in 1990	
SIC	Chemical	TRI	Pretreat	Difference	
 	CHLORINE AMMONIA	52,463 16,145	NR NR	52,463 16,145	
20	SULFURIC ACID	30,000	NR	30,000	
20	AMMONIA	228,500	NR	228,500	
2015	AMMONIA	29,000	NR	29,000	
2015	AMMONIA	31,000	NR	31,000	
2016	CYANIDE COMPOUNDS	NR	12,067	12,067	
2016	AMMONIA	38,000	NR	38,000	
2021	PHOSPHORIC ACID	21,000	NR	21,000	
2024	PHENOL	NR	1,492,663	1,492,663	
2024	PHOSPHORIC ACID	26,464	NR	26,464	
2026	PHOSPHORIC ACID	11,600	NR	11,600	
2070	PHOSPHORIC ACID	20,000	NR	20,000	
2077	AMMONIA	60,824	NR	60,824	
2082	AMMONIA	18,000	NR	18,000	
22	XYLENE 1, 2-DICHLOROBENZENE	24,300 20,250	NR NR	2 <b>4</b> ,300 20,250	
2211	ARSENIC	NR	55,992	55,992	
2221 2221	METHANOL AMMONIA	66,973 24,433	NR NR	66,973 24,433	
2221	TETRACHLOROETHYLENE	28,000	NR	28,000	
2228	SULFURIC ACID	10,500	NR	10,500	
2231	CADMIUM	NR	10,162	10,162	
2251	AMMONIUM SULFATE (SOLUTION)	59,350	NR	59,350	
2251	PHENOL	NR	664,365	664,365	
2251	PHENOL	NR 2	30,844,779	230,844,779	
2251	AMMONIA	20,038	NR	20,038	
2251	AMMONIUM SULFATE (SOLUTION)	60,000	NR	60,000	
2252	PHENOL	NR	10,707	10,707	
2252 2252	XYLENE TETRACHLOROETHYLENE	13,600 15,500	NR NR	13,600 15,500	
2252	AMMONIUM SULFATE (SOLUTION)	34,080	NR	34,080	
All q	Nantities reported in lbs.	R = Not Repor	ted	<pre>** = Values reported to</pre>	

\*\* = Values reported to both TRI and Pretreatment

Page 1

07/22	/93 Facilities With Significa	ant Discrepancies	s Between TRI a	and Pretreatmen	t in 1990.
SIC	Chemical	TRI	Pretreat	Difference	
2252	AMMONIUM SULFATE (SOLUTION)	39,000	NR	39,000	
2252	PHENOL	NR	11,332	11,332	
2252	AMMONIUM SULFATE (SOLUTION)	27,700	NR	27,700	
2252	CHLORINE	23,250	NR	23,250	
2253	PHENOL	NR	53,577	53,577	
2253	ETHYLENE GLYCOL	10,000	NR	10,000	
2253	MERCURY	NR	20,359	20,359	
2257	BIPHENYL	14,000	NR	14,000	
2257	FORMALDEHYDE	102,000	NR	102,000	
2258 2258	ETHYLENE GLYCOL XYLENE	13,000 28,000	NR NR	13,000 28,000	
2258 2258	ZINC CADMIUM	NR NR	76,191 35,556	76,191 35,556	
2258	TETRACHLOROETHYLENE	31,161	NR	31,161	
2258	AMMONIUM SULFATE (SOLUTION)	60,400	NR	60,400	
2258	XYLENE	22,500	NR	22,500	
2261	AMMONIUM SULFATE (SOLUTION)	32,000	NR	32,000	
2261	AMMONIA	147,000	NR	147,000	
2261 2261	GLYCOL ETHERS ETHYLENE GLYCOL	33,026 28,907	NR NR	33,026 28,907	
2262 2262	XYLENE AMONIA	131,088 31,764	NR NR	131,088 31,764	
2262	TETRACHLOROETHYLENE	10,000	NR	10,000	
2262	PHENOL	NR	31,378	31,378	
2262	MERCURY	NR	15,500	15,500	
2262	DECABROMODIPHENYL OXIDE	14,800	NR	14,800	
2262	1, 1, 1-TRICHLOROETHANE	20,000	NR	20,000	
2262	AMMONIUM SULFATE (SOLUTION)	43,165	NR	43,165	
2262	AMMONIA	14,000	NR	14,000	
2269	TETRACHLOROETHYLENE	11,713	NR	11,713	
2269	BIPHENYL	12,800	NR	12,800	
2269	PHENOL	NR	495,427	495,427	
All c	quantities reported in lbs.	NR = Not Repo	rted	** = Values 1	reported to both TRI and Pretreatment

E-33

and Pretreatment	Difference
Between TRI	Drotroat
Discrepancies	T chris
Facilities With Significant	
/93	lerimodr
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SIC	Chemical	TRI	Pretreat	Difference
2269	FLUOMETURON	81,000	NR	81,000
2269	PHENOL	NR	18,385	18,385
2269	PHENOL	NR	8,400,459	8,400,459
2269	PHENOL	NR	17,658	17,658
2269 2269 2269	1,2,4-TRICHLOROBENZENE PSEUDOCUMENE BIPHENYL	20,817 16,946 26,164	NR NR NR	20,817 16,946 26,164
2282 2282	PHENOL ETHYLENE GLYCOL	NR 18,000	264,679,317 NR	264,679,317 18,000
2282	1,2,4-TRICHLOROBENZENE	51,000	NR	51,000
2284 2284 2284	METHANOL FORMALDEHYDE CHLOR INE	208,000 46,000 NR	NR NR 156,388	208,000 46,000 156,388
2300	CHLORINE	42,000	NR	42,000
2321 2321	PHENOL MERCURY	NR NR	62,995 25,475	62,995 25,475
2322	PHENOL	NR	3,515,921	3,515,921
2439	CHLORINE	NR	10,289	10,289
2793	NITRIC ACID	26,880	NR	26,880
2821 2821 2821	METHANOL Formaldehyde Ethylene glycol	13,040 90,520 48,180	NR NR NR	13,040 90,520 48,180
2821	AMMONIA	25,000	NR	25,000
2833 2833 2833 2833 2833 2833 2833 2833	TOLUENE METHYLENE CHLORIDE CHLOROFORM AMMONIA ACETONE	53,000 23,000 75,000 40,000 141,000	69, 077 NR NR NR NR	16,077 •• 23,000 75,000 40,000 141,000
2834	METHANOL	38,297	NR	38,297
2834	ARSENIC	NR	14,736	14,736
2834	GLYCOL ETHERS	50,000	NR	50,000
2843	PHENOL	NR	65,801	65,801
2844	GLYCOL ETHERS	74,000	NR	74,000
2865	AMMONIA	12,000	NR	12,000
2869	ETHYLENE GLYCOL	15,570	NR	15,570
2869 2869	TOLUENE METHANOL	10,297 13,560	21 NR	10,276 ** 13,560

\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

All quantities reported in lbs.

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in 1990

SIC	Chemical	TRI	Pretreat	Difference	
2869 2869 2869	FORMALDEHYDE DIMETHYL PHTHALATE n-BUTYL ALCOHOL		NR NR NR		
3069	METHANOL	22,173	NR	22,173	
3188	HYDROCHLORIC ACID	22,000	NR	22,000	
329 329	PHENOL BARIUM COMPOUNDS 1,	46,000 900,000	NR NR	46,000 1,900,000	
3321	PHENOL	NR	3,506,294	3,506,294	
3411	PHENOL	NR	154,537	154,537	
3425	PHENOL	NR	28,328,606	28,328,606	
3471	TOLUENE	N I	190,005	190,005	
3471 3471	SELENTUM	¥ ¥ !	19,388	19,388	
34/1 3471 3471	MERCURY LEAD COPPER	NN NN NN	20,681 20,681 23,266	19, 803 20, 681 23, 266	
3496	PHENOL	NR	98,391	98,391	
3499	SULFURIC ACID	91,650	NR	91,650	
3519 3519	SULFURIC ACID DIETHANOLAMINE	57,243 14,456	NR NR	57,243 14,456	
3553	GLYCOL ETHERS	37,239	NR	37,239	
3612 3612 3612 3612 3612	TOLUENE SELENIUM MERCURY LEAD COPPER	0 NR NR NR	54,936 13,734 68,670 20,601 13,734	54,936 *• 13,734 *• 68,670 20,601 13,734	
3622	GLYCOL ETHERS	26,687	NR	26,687	
3676	PHENOL	NR	83,356	83,356	
3692	PHENOL	NR	19,666,082	19,666,082	
3713	PHOSPHORIC ACID	12,000	NR	12,000	

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\*\* = Values reported to both TRI and Pretreatment

NR = Not Reported

All quantities reported in lbs.

66,831,600 15,981

66,831,600 15,981

NR NR

ARSENIC

3999

3714 PHENOL

07/23/93 Page 1	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Constituent/Was	te Code	Lbs Tr 1988	ansferred O 1989	ff-site 1990	
Facility 000002					
F-Waste Cons	tituents Reported to TRI:				
1,1,1-TRI	CHLOROETHANE	0	0	0	
	Constituent Total	0	0	0	
F-Wastes Rep	orted to NCAR:		0	3 561	
F001 F002		14,500	10,120	14,273	
F005		500			
	Waste Total	15,000	10,120	17,834	
Facility 000002					
F-Waste Cons FREON 113	tituents Reported to TRI:	2,034	5,710	13,800	
METHYLENE	CHLORIDE	2,033	5,710	0	
	Constituent Total	4,067	11,420	13,800	
F-Wastes Rep	orted to NCAR:	4 067	22 826	10 210	
F002		26,120	32,505	51,300	
	Waste Total	30,187	55,341	70,610	
Facility 000007					
F-Waste Cons	tituents Reported to TRI:				
BENZENE n-BUTYL A	LCOHOL	0 0	0 0	0 0	
METHANOL TOLUENE		0 0	9,360 0	10,415 0	
	- Constituent Total	0	9,360	10,415	
F-Wastes Rep	orted to NCAR:				
F001	-	72,320	0	0	
	Waste Total	72,320	0	0	
Facility 000009					
F-Waste Cons FRFON 113	tituents Reported to TRI:	48.670	88.085	198.212	
METHYLENE	CHLORIDE	11,000	0	0	
	Constituent Total	59,670	88,085	198,212	
F-Wastes Rep	orted to NCAR:	0	0	E 500	
D001F003F D003F002		0	0	172,682	
F002	-	108,242	229,271	348,692	
	Waste Total	108,242	229,271	526,954	
Facility 000010					
F-Waste Cons METHYLENF	tituents Reported to TRI: CHLORIDE	0	0	0	
1,1,1-TRI	CHLOROETHANE	231	0	0	
	Constituent Total	231	0	0	

07/23/93 Page 2	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			1990
Constituent/Wast	e Code	Lbs Tra 1988	ansferred Of 1989	f-site 1990
F-Wastes Repo D001F002F0 D001F003F0 D005F002 D007F001 F001 F002	orted to NCAR: 003 005	0 0 0 73,170 11,541	0 0 0 1,237 1,541	335 459 250 139 627 3,376
	Waste Total	84,711	2,778	5,186
Facility 000011				
F-Waste Const METHANOL	tituents Reported to TRI:	0	0	0
	Constituent Total	0	0	0
F-Wastes Repo D001F003F0 D019F002 D035F005 •F001 F002 F003	orted to NCAR: 005U117	0 0 0 293,690 0	0 0 800 58,104 300	100 417 1,850 0 0
	Waste Total	293,690	59,204	2,367
Facility 000012				
F-Waste Const 1,1,1-TRIC	tituents Reported to TRI: CHLOROETHANE	0	0	0
	- Constituent Total	0	0	0
F-Wastes Repo F001 F002	orted to NCAR:	99 <b>4</b> 0	0 35	0 0
	Waste Total	994	35	0
Facility 000017				
F-Waste Const ACETONE BENZENE METHANOL TOLUENE XYLENE	tituents Reported to TRI:	0 0 0 0	0 26 0 130 800	0 0 0 0
	Constituent Total	0	956	0
F-Wastes Repo F003F005 F005	orted to NCAR:	0 1, <b>4</b> 75	2,888 0	0 0
	Waste Total	1,475	2,888	0
Facility 000024				
F-Waste Const 1,1,1-TRI	tituents Reported to TRI: CHLOROETHANE	0	0	0
	- Constituent Total	0	0	0

07/23/93 Page 3	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Constituent/Wast	e Code	Lbs Tra 1988	ansferred Off 1989	-site 1990	
F-Wastes Repo F001	rted to NCAR:	8,065	0	0	
	Waste Total	8,065	0	0	
Facility 000028					
F-Waste Const	ituents Reported to TRI:	:			
ACETONE	-	0	0	0	
n-BUTYL AL	COHOL	0	0	0	
METHYL ETH	YL KETONE (MEK)	0	0	0	
METHANOL		0	0	0	
TOLUENE	BOTTL RETONE	0	0	0	
XYLENE		ŏ	Ö	õ	
	Constituent Total				
	constituent iotai	Ū	Ũ	v	
F-Wastes Repo	orted to NCAR:	٥	178 900	0	
D001F003 D035F003F0	05	0	178,900	199 838	
F005	05	293,920	219,585	155,058	
	Waste Total	293.920	398,485	199.838	
		2307520	,	200,000	
Facility 000032					
F-Waste Const	ituents Reported to TRI:				
ACETONE		2,678	4,553	3,050	
n-BUTYL AL	COHOL	560	1,468	0	
CYCLOHEXAN	E	3,896	0	6,664	
METHYL ETH	YL KETONE (MEK)	3,653	8,813	7,682	
METHYL ISO	BUTYL KETONE	730	1,908	1,743	
TOLUENE		3,409	5,973	7,407	
XYLENE	-	1,461		0	
	Constituent Total	16,387	22,715	26,546	
F-Wastes Repo	orted to NCAR:				
F003		40,590	69,955	43,576	
	Waste Total	40,590	69,955	43,576	
Facility 000033					
F-Waste Const	ituents Reported to TRI:	:	•	0	
METHANOL XVI ENE		2 052	1 969	1 979	
AILENE		2,052	1,000	1,0/0	
	Constituent Total	2,052	1,868	1,878	
F-Wastes Repo	orted to NCAR:				
F003		2,052	1,868	1,878	
	Waste Total	2,052	1,868	1,878	
Facility 000034					
INCITICA 000034					
F-Waste Const	ituents Reported to TRI:	:			
METHANOL	_	0	0	0	
1,1,1-TRIC	HLOROETHANE	0	0	0	
	Constituent Total	0	0	0	

07/23/93 Page	Comparison of F-Wastes F-Waste Constituents Re	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Constitu	lent/Waste Code	Lbs Tra 1988	ansferred Off 1989	-site 1990		
F-Wa: F F	stes Reported to NCAR: 001 005	1,407 173	0 0	0 0		
	- Waste Total	1,580	0	0		
Facility	Y 000038					
F-Wa	ste Constituents Reported to TRI:					
F	REON 113	19,200	0	0		
M 1	ETHYL ETHYL KETONE (MEK) ,1,1-TRICHLOROETHANE -	21,600 10,800	13,275 1,200	11,931 0		
	Constituent Total	51,600	14,475	11,931		
F-Wa	stes Reported to NCAR:					
F	001	3,150	1,210	1,300		
F	002	19,200	14,571	1,850		
F F	003 003F005	21,600	26,060	99,770		
	- Waste Total	43,950	41,841	102,920		
Facilit	Y 000040					
F-Wa 1	ste Constituents Reported to TRI: ,1,1-TRICHLOROETHANE	0	5,337	7,339		
	- Constituent Total	0	5,337	7,339		
F-Wa	stes Reported to NCAR:					
D	001D002F003F005	0	0	1,520		
D	007F005	0	800	0		
F	001	729	5,337	7,339		
F F	003 003F005	1,750	1,167	3,723		
	- Waste Total	2,479	9,204	12,582		
Facilit	y 000042					
F-Wa	ste Constituents Reported to TRI:					
М	ETHYL ETHYL KETONE (MEK)	0	0	0		
Т	RICHLOROETHYLENE	692	541	610		
	Constituent Total	692	541	610		
F-Wa	stes Reported to NCAR:					
D	001F001	0	9,016	0		
D	001F005		6,462	6,928		
न न	001 005	5,878	0	10,893		
	Waste Total	17,415	15,478	17,821		
Facilit	y 0000 <b>4</b> 6					
F-Wa	ste Constituents Reported to TRI:					
A	CETONE	0	0	0		
n	-BUTYL ALCOHOL	0	0	0		
E	THYLBENZENE Fruvi Fruvi Fruvie (MER)	U 0	0	0		
M M	ETHANOL	0	õ	õ		
M	ETHYL ISOBUTYL KETONE	Ō	0	0		
Т	OLUENE	0	0	0		

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Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990 \_

Constituent Wester Cale	Lbs Tra	ansferred Off	E-site
Constituent/Waste Code	1988	1989	
XYLENE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:			
F003	0	117 020	85,173
F005	101,640	0	0
- Waste Total	101,640	117,920	85,173
Facility 000047			
E-Wasto Constituents Perorted to TRI.			
n-BUTYL ALCOHOL	13,027	750	250
METHYL ETHYL KETONE (MEK)	12,261	750	250
METHANOL TOLUENE	26,859	750	250 750
XYLENE	5,364	750	250
Constituent Total	74,333	3,750	1,750
F-Wastes Reported to NCAR:			
F003	0	270,102	192,154
Waste Total	0	270,102	192,154
Facility 000048			
F-Waste Constituents Reported to TRI:			
FREON 113	11,586	13,544	6,532
1, 1, 1-TRICHLOROETHANE	4,920		4,400
Constituent Total	16,514	13,544	10,932
F-Wastes Reported to NCAR:	•		
D001F002F005 F001	0	4,400	0 000
F001F003F005	ŏ	1,800	4,400
F002	19,800	0	0
F002F003 F002F003F004	0	11,900	0
F002F005	Ő	1,200	7,100
F003	0	4,000	0
F005 -	0	800	0
Waste Total	19,800	39,180	12,400
Facility 000052			
F-Waste Constituents Reported to TRI:			-
METHYL ETHYL KETONE (MEK) TOLUENE	0	0	0
Constituent Total	0	0	0
E Master Deported to MORD.			
F-wastes reported to NCAR: F003	25,560	0	0
F003F005	0	40,200	90,800
- Waste Total	25,560	40,200	90,800

## Facility 000053

F-Waste Constituents Reported to TRI:

07/23 Page	/93 6	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Const	ituent/Was	- te Code	Lbs Tra 1988	ansferred Off 1989	f-site 1990	
	1.1.1-TRI	CHLOROETHANE			0	
		Constituent Total	U	U	0	
F-	Wastes Repo F001	orted to NCAR:	0	59,739	0	
		Waste Total	0	59,739	0	
Facil	ity 000056					
F-	Waste Cons	tituents Reported to TRI:				
	METHYLENE	CHLORIDE	0	0	0	
	TOLUENE		0	750	250	
	1,1,1-1RI					
		Constituent Total	0	750	250	
F-	Wastes Rep	orted to NCAR:				
	F001		4,390	3,044	2,656	
	F003 F003F005		0	587	341 1 840	
	• F005		17,117	12,469	9,684	
		 Waste Total	21,507	17,455	14,521	
Facil	ity 000059					
F -	D002F003	cituents Reported to IRI:	0	0	758	
	F002		2,574	0	0	
	F003		0	328	0	
		Waste Total	2,574	328	758	
Facil	ity 000060					
F-	Waste Const	tituents Reported to TRI:				
	ACETONE		0	4,300	10	
	METHANOL	LCOHOL	1 142	2,250	500	
	TOLUENE		15,197	2,250	500	
	XYLENE		498	2,250	500	
		Constituent Total	16,837	13,300	2,010	
F-	Wastes Repo	orted to NCAR:				
	F003		32,560	31,640	35,840	
		Waste Total	32,560	31,640	35,840	
Facil	ity 000061					
F-	Waste Const	tituents Reported to TRI:				
	TETRACHLO	ROETHYLENE	1,564	0	0	
	1.1.1-TRT	CHLOROETHANE	6,912	2,/12	6,484 0	
	1,1,1 11(1)					
		Constituent Total	8,476	2,712	6,484	
F-	Wastes Repo	orted to NCAR:	•	· ·	20.000	
	D001F002		U O	U N	32,800 400	
	F002		146,400	258,908	229,000	
	F002F005		0	0	32,000	

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Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

Constituent/Waste Code	Lbs Tr 1988	ansferred O 1989	ff-site 1990
F003 F005	2,400 3,200	6,731 392	28,400
Waste Total	152,000	266,031	322,600
Facility 000075			
F-Waste Constituents Reported to TRI ACETONE n-BUTYL ALCOHOL METHANOL TOLUENE	: 0 0 0	0 0 0	0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:	<u>^</u>	20.047	• 05 405
F003F005 F005	28,490	39,847	25,426
Waste Total	28,490	39,847	25,426
Facility 000078			
F-Waste Constituents Reported to TRI ACETONE n-BUTYL ALCOHOL METHANOL TOLUENE 1,1,1-TRICHLOROETHANE XYLENE	: 2,760 0 12,200 94,905 5,450	2,328 0 0 55,200 0	0 0 0 0 0 0
Constituent Total	115,315	57,528	0
F-Wastes Reported to NCAR: F001 F002 F002F003 F002F004	93,520 1,400 0 0	78,712 0 10,000 0	60,500 0 1,200
Waste Total	94,920	88,712	61,700
Facility 000080 F-Waste Constituents Reported to TRI XYLENE	: 0	11,000	12,000
Constituent Total	0	11,000	12,000
F-Wastes Reported to NCAR: D001F003 F003 F003F005 F005	0 14,000 0 20,020	5,995 6,400 7,339 0	0 458 3,720 0
Waste Total	34,020	19,734	4,178
Facility 000081			
F-Waste Constituents Reported to TRI FREON 113 1,1,1-TRICHLOROETHANE	: 0 0	46 0	0 0
Constituent Total	0	46	0

07/23/93 Page 8	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Constituent/Wast	e Code	Lbs T 1988	ransferred 0 1989	ff-site 1990	
F-Wastes Repo	rted to NCAR:	7 579	10 212	8 205	
F001 F003		7,578	10,212	385	
F003F005		õ	Ő	1,142	
	- Waste Total	7,578	10,212	9,732	
Facility 000083					
F-Waste Const	ituents Reported to TRI:				
F001		7,285	2,566	0	
	Waste Total	7,285	2,566	0	
Facility 000086					
F-Waste Const	ituents Reported to TRI:	:			
FREON 113		0	0	0	
	Constituent Total	0	0	0	
F-Wastes Repo	rted to NCAR:				
• F002		263,968	160,768	150,644	
	Waste Total	263,968	160,768	150,644	
Facility 000087					
F-Waste Const	ituents Reported to TRI:				
METHANOL		0	0	0	
TOHOLINE	-		°		
	Constituent Total	U	0	0	
F-Wastes Repo	rted to NCAR:				
F003F005		0	12,800	17,600	
F005	_	11,600	0	U	
	Waste Total	11,600	12,800	17,600	
Facility 000090					
E-Waste Const	ituents Reported to TRI				
ACETONE	Trance Reported to INT	0	0	0	
n-BUTYL AL	COHOL	0	0	0	
METHANOL		0	0	0	
TOLUENE		0	0	0	
XYLENE	_		0		
	Constituent Total	0	0	0	
F-Wastes Repo	rted to NCAR:				
F003F005		0	15,200	14,400	
F005	-	20,000		6,797	
	Waste Total	20,000	15,200	21,197	
Facility 000091					
F-Waste Const	ituents Reported to TRI:	:			
ACETONE		0	0	0	
n-BUTYL AL	COHOL	0	0	0	
METHANOL TOLUENE		0	0	0	
XYLENE		ő	ő	ŏ	

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07/23/93 Page 9	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Constituent/Was	ste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
	- Constituent Total			0	
F-Wastes Rep F003F005 F005	ported to NCAR:	0 43,200	38,800 0	42,000 0	
	Waste Total	43,200	38,800	42,000	
Facility 000092	2				
F-Waste Cons ACETONE n-BUTYL A METHYL E METHANOL TOLUENE XYLENE	stituents Reported to TRI: ALCOHOL THYL KETONE (MEK)	0 0 0 0 0		0 0 0 0 0	
F-Wastes Rep F003F005 F005	ported to NCAR:	0 40,800 	70,400	40,000 0 	
Facility 000093	3			,	
F-Waste Cons ACETONE n-BUTYL 2 METHYL E METHANOL TOLUENE XYLENE	stituents Reported to TRI: ALCOHOL THYL KETONE (MEK)	0 0 0 0 0	0 0 0 0 0	0 0 0 0 0	
	Constituent Total	0	0	0	
F-Wastes Rep F003F005 F005	ported to NCAR: - Waste Total	64,000 64,000	66,000 0 66,000	86,000 0  86,000	
Facility 000094	4				
F-Waste Cons ACETONE n-BUTYL X METHYL E METHANOL TOLUENE XYLENE	stituents Reported to TRI: ALCOHOL THYL KETONE (MEK)	0 0 0 0 0	0 0 0 0 0	0 0 0 0 0 0	
	Constituent Total	0	0	0	
F-Wastes Rej F003F005 F005	ported to NCAR:	0 44,800	62,000	0 0	
	Waste Total	44,800	62,000	0	

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07/23 Page	/93 10	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Const	ituent/Wast	e Code	Lbs Tra 1988	ansferred Of: 1989	f-site 1990	
	ity 000098					
F-	Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL TOLUENE XYLENE	ituents Reported to TRI: COHOL YL KETONE (MEK)	0 0 0 0 0	0 0 0 0 0	0 0 0 0 0	
		- Constituent Total	0	0	0	
F-	Wastes Repo F003F005 F005	rted to NCAR:	0 56,000	67,200 0	86,800 0	
		Waste Total	56,000	67,200	86,800	
Facil F-	Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL METHYL ISO TOLUENE XYLENE	ituents Reported to TRI: COHOL YL KETONE (MEK) BUTYL KETONE	130,000 250 0 0 110,000 120,000	11,750 750 14,750 7,750 14,750 53,750 35,750	19,255 5 23,155 9,355 2,900 60,255 20,955	
		Constituent Total	360,250	139,250	135,880	
F-	Wastes Repo D001D009F0 D001F003F0 F005	rted to NCAR: 03F005 05	0 0 625,193	0 352,220 0	8,800 304,989 0	
		Waste Total	625,193	352,220	313,789	
Facil F-	ity 000100 Waste Const FREON 113 METHYLENE	ituents Reported to TRI: CHLORIDE Constituent Total	7,500 86,400 93,900	10,265 32,100 42,365	5,075 22,800 27,875	
F-	Wastes Repo F001 F002	rted to NCAR:	0 115,250	2,000 38,400	2,400 19,800	
		Waste Total	115,250	40,400	22,200	
Facil	ity 000105					
F-	Waste Const D001F001 F001 F002 F003	ituents Reported to TRI:	0 2,100 900 0	0 1,000 0 542	7,480 0 880	
		Waste Total	3,000	1,542	8,360	

Facility 000106

F-Waste Constituents Reported to TRI:

Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Tra	ansferred Off	-site
Constituent/Waste Code	1988	1989	1990
ACETONE METHYL ETHYL KETONE (MEK) METHANOL TOLUENE	0 0 0	0 0 0	0 0 0
XYLENE			0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F003F005 F005	12,400 0 21,200	26,000 0	0 0 18,400
- Waste Total	33,600	26,000	18,400
Facility 000109			
F-Waste Constituents Reported to TRI:			
METHYLENE CHLORIDE 1,1,1-TRICHLOROETHANE	0 0	0 0	0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F001 F002	1,368 184,964	0 198,175	0 69,467
- Waste Total	186,332	198,175	69,467
Facility 000115			
F-Waste Constituents Reported to TRI:	1		
1,1,1-TRICHLOROETHANE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:			
F001 F003	5,850	4,219	20,620
F005	0	1,236	100
Waste Total	6,046	5,455	23,098
Facility 000117			
F-Waste Constituents Reported to TRI:	:		
METHYL ETHYL KETONE (MEK)	0	0	0
TOLUENE XYLENE	0	0	0
Constituent Total	0		0
E-Wastes Reported to NCAR.			
F001	21,800	0	400
F002 F002F002F005	6,600	2,500	500
F003	õ	0	400
F003F005	0	16,100	8,800
F005	15,900	800	1,200
Waste Total	44,300	20,200	11,300

Facility 000120

F-Waste Constituents Reported to TRI:

07/23/93 Page 12		Comparison of F-Wastes R F-Waste Constituents Rep	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/W	aste Code	Lbs Tra 1988	ansferred Off 1989	-site 1990	
		ETUVI RETONE (MER)		750		
	TOLUENE		0 0	750 1,500	250 0	
		Constituent Total	0	3,000	250	
F-	Wastes F	eported to NCAR:				
	F002F00		0	77,360	65,814	
	F003F00 F005	15	116,640	3 <b>2,4</b> 00 0	õ	
		 Waste Total	116,640	109,760	65,814	
Facil	ity 0001	.23				
F-	Waste Co	onstituents Reported to TRI:				
	ACETONE	NI COHOI	0	0	0	
	METHYL	ETHYL KETONE (MEK)	Ň	Õ	ŏ	
	METHANC		0	0	0	
	TOLUENE	2	0	0	0	
	•	 Constituent Total	0		0	
F-	Wastes H	Reported to NCAR:				
-	F005		38,800	57,200	56,100	
		Waste Total	38,800	57,200	56,100	
Facil	ity 0001	124				
F-	-Waste Co	onstituents Reported to TRI:			4 645	
	1,1,1-7	TRICHLOROETHANE	2,958	3,383	1,605	
		Constituent Total	2,958	3,383	1,605	
F-	Wastes H	Reported to NCAR:	2 050	2 202	1 605	
	F001		2,958	3,383	1,605	
		Waste Total	2,958	3,383	1,605	
Facil	lity 0001	131				
F-	-Waste Co	onstituents Reported to TRI:	0	750	500	
	FREON	(5)	0	/50	0	
	METHYL	ETHYL KETONE (MEK)	13,632	6,245	8,350	
	TETRACI XYLENE	ILOROETHYLENE	14,725 0	2,562	11,000 0	
		 Constituent Total	28,357	9,557	19,850	
- च	-Wastes H	Reported to NCAR:				
•	F001		18,406	50,033	13,595	
	F002		0 २ 14२	11,365	7,761 3,477	
	F003		1,040	2,750	8,414	
		 Waste Total	22,589	65,940	33,247	
Facil	Lity 0003	134				
F-	-Waste Co	onstituents Reported to TRI:				
-	ACETON	-	1,979	750	0	

07/23 Page	13 13	Comparison of F-Wastes F-Waste Constituents Re	Reported to	NCAR and TRI: 1988 to	1990
Const	ituent/Wast	e Code	Lbs T 1988	ransferred O 1989	ff-site 1990
	n-BUTYL AL ETHYLBENZE METHYL ETH METHANOL METHYL ISO TOLUENE XYLENE	COHOL NE YL KETONE (MEK) BUTYL KETONE	0 0 15,804 10,346 0 3,631 1,246	750 750 750 0 0 0 750	0 0 0 0 0 0 0
		Constituent Total	33,006	3,750	0
F-	Wastes Repo F003 F005	rted to NCAR:	1,280,234 751,832	1,717,351 0	0 1,027,695
		Waste Total	2,032,066	1,717,351	1,027,695
Facil	lity 000137				
F-	Waste Const ACETONE n-BUTYL AL CYCLOHEXAN ETHYLBENZE METHYL ETH METHYLENE METHYL ISO TETRACHLOR TOLUENE 1,1,1-TRIC TRICHLOROE XYLENE	ituents Reported to TRI COHOL E NE YL KETONE (MEK) CHLORIDE BUTYL KETONE OETHYLENE HLOROETHANE THYLENE	: 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0
		Constituent Total	0	0	0
F	-Wastes Repo F003 F003F005	rted to NCAR:	187,110 0	0 339,041	0 295,827
		Waste Total	187,110	339,041	295,827
Facil	lity 000139				
F·	-Waste Const METHANOL TOLUENE	ituents Reported to TRI	: 0 2	0 16,373	0 370
		Constituent Total	2	16,373	370
F	-Wastes Repo F005	rted to NCAR:	16,100	24,640	110,120
		Waste Total	16,100	24,640	110,120
Faci	lity 000141				
F	-Waste Const METHYL ETH TOLUENE	ituents Reported to TRI YL KETONE (MEK)	: 0 0	0 0	0 0
		Constituent Total	0	0	0
F	-Wastes Repo F005	rted to NCAR:	0	5,540	0

07/23 Page	/93 14	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Ti 1988	ransferred Of 1989	f-site 1990
		Waste Total	0	5.540	
Facil	ity 000142	huste iotui	Ũ	0,010	, i i i i i i i i i i i i i i i i i i i
	-				
F-1	Waste Const 1,1,1-TRIC	HLOROETHANE	0	0	0
		Constituent Total	0	0	0
F-1	Wastes Repo	rted to NCAR:			
	D001D022D0	23F003	0	0	4,000
	D001F001		400	0	1,500
	F001		1,050	õ	2,350
	F003		800	0	0
	F003F004 F005		0 6,600	3,500 0	0 0
		- Waste Total	8,850	3,500	11,650
Facil	ity 000143				
F-1	Waste Const	ituents Reported to TRI:			
-	1,1,1-TRIC	HLOROETHANE	0	0	0
		Constituent Total	0	0	0
F-'	Wastes Repo	orted to NCAR:			
	F001	_	208,250	0	0
		Waste Total	208,250	0	0
Facil	ity 000144				
F-	Waste Const	ituents Reported to TRI:			
	n-BUTYL AL	COHOL	7,792	0	0
	METHANOL		43,781	0	0
	TOLUENE		0	0	0
		-			
		Constituent Total	51,573	0	0
F-	Wastes Repo	orted to NCAR:	70 000	^	70 100
	F003 F005		70,320 0	1.392	79,100 84,750
	1005	-	·		
		Waste Total	70,320	1,392	163,850
Facil	ity 000148				
F-	Waste Const	ituents Reported to TRI:		-	
	D001D039F0	03F005	0	0	5,021
	D001F003		7 440	8,340 0	417
	F003	-			
		Waste Total	7,440	8,340	5,438
Facil	ity 000149				
F-	Waste Const	ituents Reported to TRI:	^	^	^
	ACETONE	COHOI	0	U N	0
	METHYL ETH	IVL KETONE (MEK)	1,579	ő	ŏ
	METHANOL		0	0	0
	METHYL ISC	BUTYL KETONE	0	0	0

## Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs T:	ransferred (	Off-site
Constituent/Waste Code	1988	1989	1990
TETRACHLOROETHYLENE	0	0	0
1,1,1-TRICHLOROETHANE	0	872	0
XYLENE	178	394	0
Constituent Total	1,757	1,266	0
F-Wastes Reported to NCAR:			
D001D002D007F002 D001F001F003F005	0	0 24,690	2,507
D002D007D008F002	0	40,876	18,585
F001 F002	5,600 31,376	0	0
- Waste Total	36,976	65,566	21,092
Facility 000151			
E-Waste Constituents Reported to TRI:			
FREON 113	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:			
F002	6,232	1,802	1,588
F003	8	4,136	4,634
Waste Total	6,240	5,938	6,222
Facility 000156			
F-Waste Constituents Reported to TRI: METHYLENE CHLORIDE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:			
F001	0	42,940	32,696
Waste Total	0	42,940	32,696
Facility 000157			
F-Waste Constituents Reported to TRI:	:		
1,1,1-TRICHLOROETHANE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:			
F001 F003F005	0	982 683	0
Waste Total		1 665	
waste Iotai	Ŭ	1,005	Ŭ
Facility 000163			
F-Waste Constituents Reported to TRI: F002	: 0	3,576	1,787
- Waste Total	0	3,576	1,787
Facility 000164		-	
E Waste Constituents Dependent to MDT.			
FREON 113	46	750	5

Page :	16 F	-Waste Constituents Rep	orted to T	RI: 1988 to 1	1990
Constit	tuent/Waste	- Code	Lbs Tra 1988	ansferred Of: 1989	f-site 1990
	1,1,1-TRICHL	OROETHANE	71	750	5
		Constituent Total	117	1,500	10
F-Wa	astes Report	ed to NCAR:			
]	D001F002U075	U121	0	0 196	161
1	D001F003F005		0	466	219
1	F001		77,034	103,784	63,315
1	F002		113,160	104,704	89,691
1	F002U226		0	173	0
1	F003 F005		1,799	0	0
		 Waste Total	193,387	209,323	153,386
acili	ty 000165				
F-Wa	aste Constit	uents Reported to TRI:	•	<u> </u>	
	ACETONE		0	0	0
1	METHYLENE CH.	LORIDE	0	0	0
•	1,1,1-TRICHL	OROETHANE	ő	0 0	Ő
		Constituent Total	0	0	0
F-Wa	astes Report	ed to NCAR:			
1	F001 F003		244,464 395,733	147,613 306,882	0 352,849
		 Waste Total	640,197	454,495	352,849
acili	ty 000170				
F-W	aste Constit	uents Reported to TRI:			
]	F001	-	0	550	125
1	F003		1,720	0	0
		Waste Total	1,720	550	125
acili	ty 000171				
F-W	aste Constit	uents Reported to TRI:	850	0	
]	F001		750		
		Waste Total	750	0	0
acili	ty 000174				
F-Wa	aste Constit	uents Reported to TRI:			
1	D001D005F002	F005	0	1,000	0
		Waste Total	0	1,000	0
acili	ty 000176				
F-Wa	aste Constit	uents Reported to TRI:	•	•	-
1	ACETONE METHYLENE CH	LORIDE	0	0 0	0 0
		 Constituent Total	0	0	0
F-W	astes Report	ed to NCAR:			
1	F002		107,500	68,961	43,355
1	F003		26,600	21,063	23,001

Comparison of F-Wastes Reported to NCAR and

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07/23, Page	/93 17	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	- ce Code	Lbs Tr. 1988	ansferred Of 1989	f-site 1990
	 F005		482	<b>44</b> 0	1,213
	1005	 Waste Total	134,582	90,464	67,569
Facil	ity 000177				
F-1	Waste Const	tituents Reported to TRI:			
	ACETONE	COHOL	0	80 80	0 78
	METHANOL		0	140	0
	TOLUENE XYLENE		0	40	0
		 Constituent Total	0	1,140	643
F-	Wastes Repo	orted to NCAR:			
	D001F003F0 F002	005K086	0 0	0 0	857,997 1,320
		Waste Total	0	0	859,317
Facil	ity 000180				
F-	Waste Const	tituents Reported to TRI:			
	ACETONE	HYL KETONE (MEK)	7,327 26,308	0	0
	METHYL IS	DBUTYL KETONE	0	0	0
	TOLUENE		7,327	0	0
		 Constituent Total	48,289	0	0
F-	Wastes Rep	orted to NCAR:	•	•	111 600
	D001D034F	003F005 005	0	0	29,400
	F001		0	7,030	0
	F003 F003F005		25,528	146,700	ŏ
		Waste Total	25,528	153,730	141,000
Facil	ity 000184				
F-	Waste Cons	tituents Reported to TRI:			
	n-BUTYL A	LCOHOL ENE	8,200 0	5,500 750	4,551 750
	METHYL ET	HYL KETONE (MEK)	0	0	0
	METHYL IS TOLUENE	OBUTYL RETONE	10,600	7,000	3,850
	XYLENE		0	750	3,760
		Constituent Total	18,800	14,000	12,916
F-	Wastes Rep	orted to NCAR: 005	0	0	6,600
	F003		0	0	14,107
	F003F005 F005		996,197	884,181	001,945
		 Waste Total	996,197	884,181	882,652
Facil	lity 000185				
F-	-Waste Cons METHYL ET	tituents Reported to TRI: HYL KETONE (MEK)	827	900	890

07/23 Page	/93 18	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Const	ituent/Wast	e Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
	METHANOL TOLUENE XYLENE		2,388 1,194 695	1,500 1,200 700	1,660 1,170 620	
		- Constituent Total	5,104	4,300	4,340	
F-	Wastes Repo	rted to NCAR:	_			
	F003 F003F005		0	6,930 0	0 6,504	
	F005		13,600	0	0	
		- Waste Total	13,600	6,930	6,504	
Facil	ity 000187					
F-	Waste Const ACETONE	ituents Reported to TRI:	0	0	0	
	n-BUTYL AL	COHOL	0	0	0	
	METHYL ETH METHANOL	YL KETONE (MEK)	0	0	0	
	METHYL ISO	BUTYL KETONE	0	0	0	
	* XYLENE		0	0	0	
		- Constituent Total	0	0	0	
F-	Wastes Repo	orted to NCAR:				
	D001D035F0 F003F005	03F005	0	0 64,050	38,850	
	F005		50,050	0	Ō	
		- Waste Total	50,050	64,050	38,850	
Facil	ity 000195.					
F-	Waste Const D001F003	ituents Reported to TRI:	0	25	0	
		- Waste Total	0	25	0	
Facil	ity 000196.					
F-	Waste Const F003	ituents Reported to TRI:	20,753	459	0	
		- Waste Total	20,753	459	0	
Facil	ity 000199.					
F-	Waste Const	ituents Reported to TRI:	0	0	0	
	METHANOL		0	0	0	
	TETRACHLOR	OETHYLENE	0	750	5	
		Constituent Total	0	750	5	
F-	Wastes Repo	orted to NCAR:	<u>^</u>	550	<u> </u>	
	D002F003 F001		ں 66,680	550 64,066	63,972	
	F002		410	0	5,855	
	F003 F003F005		22,978	0 11,676	40 11,167	
	F005		790	0	0	
	F005F006	_	U	5,753	0	

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07/2 Page	23/93 2 19	Comparison of F-Wastes F-Waste Constituents Re	Reported to eported to I	NCAR and RI: 1988 to	1990
Cons	stituent/Wast	e Code	Lbs Tr 1988	ansferred Of 1989	f-site 1990
		Waste Total	90 858	82 045	81 034
Faci	lity 000200	waste local	50,050	02,045	01,034
F	-Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL TOLUENE XYLENE	ituents Reported to TRI COHOL YL KETONE (MEK)	: 425 1,187 781 746 2,092 1,662	1,275 2,864 1,762 1,749 2,253 0	5 250 250 250 250 250
		Constituent Total	6,893	9,903	1,255
F	F-Wastes Repc D001F003F0 F003 F003F005	orted to NCAR: 05	0 12,000 0	17,700 0 15,000	137,400 0 137,400
		waste Total	12,000	32,700	137,400
Facı	· · ·				
F	F-Waste Const METHYLENE METHANOL	ituents Reported to TRI CHLORIDE	: 4,423 0	1,000 0	0 0
		Constituent Total	4,423	1,000	0
F	F-Wastes Repo F002 F003 F005	orted to NCAR:	5,555 200 0	1,000 480 60	2,515 500 750
		Waste Total	5,755	1,540	3,765
Faci	ility 000205				
I	F-Waste Const METHANOL	ituents Reported to TRI	: 4,482	4,611	0
		Constituent Total	4,482	4,611	0
I	F-Wastes Repo F003 F003F005	orted to NCAR:	160,160 0	137,2 <b>4</b> 0 3,670	35,530 0
		Waste Total	160,160	140,910	35,530
Fac:	ility 000207				
1	F-Waste Const n-BUTYL AI METHYL ETH TOLUENE XYLENE	tituents Reported to TRI LCOHOL HYL KETONE (MEK)	: 3,510 10,200 32,700 3,510	3,713 10,313 33,000 3,548	0 0 0 0
		Constituent Total	49,920	50,574	0
1	F-Wastes Repo F003 F003F005 F005	orted to NCAR:	364,995 0 44,550	935 513,975 0	0 551,775 0
		Waste Total	409,545	514,910	551,775

07/23 Page	/93 20	Comparison of F-Wastes I F-Waste Constituents Rep	Reported to ported to TH	NCAR and RI: 1988 to 1	.990
Const	ituent/Was	te Code	Lbs Tra 1988	ansferred Off 1989	-site 1990
Facil	1CY 000208				
F-	Waste Cons n-BUTYL A ETHYLBENZ METHYL ET METHANOL METHYL IS TOLUENE	COBUTYL KETONE	0 0 0 0	107 141 1,766 0 140 2,889	228 294 3,420 930 601 4,500
	1,1,1-TRI XYLENE	CHLOROETHANE	0	2,748	4,500
		- Constituent Total	0	7,974	15,093
F-	Wastes Per	ported to NCAR.			
r	D001F003 F003 F003F005 F005		0 812,371 0 2,295	<b>41,043</b> 720,336 0 0	0 0 664,569 0
		- Waste Total	814,666	761,379	664,569
• Facil	ity 000210	)			
F-	Waste Cons TOLUENE	stituents Reported to TRI:	0	0	0
		- Constituent Total	0	0	0
F-	Wastes Rep D002D007F F001F002F F002 F003 F005U0190	oorted to NCAR: 7005 7003F005 J228 - Waste Total	0 0 0 0 0	140 4,950 0 500 5,590	0 20 250 0 270
Facil	ity 000211	· · · · · · · · · · · · · · · · · · ·		·	
F-	Waste Cons ACETONE n-BUTYL / METHYL E METHANOL TOLUENE XYLENE	stituents Reported to TRI: ALCOHOL THYL KETONE (MEK)	2,700 0 28,000 0 12,300 0	4,200 0 2,500 11,000 750	3,846 1,051 250 2,299 10,951 0
		- Constituent Total	43,000	19,200	18,397
F-	Wastes Rep F003F005 F005	ported to NCAR:	0 70,854	45,790 0	53,502 0
		Waste Total	70,854	45,790	53,502
Facil	lity 000212	2			
F-	Waste Con: METHYLEN 1,1,1-TR	stituents Reported to TRI: E CHLORIDE ICHLOROETHANE	0 0	18,200 0	28,400 0
		Constituent Total	0	18,200	28,400

See .....

07/23/93 Page 21	Comparison of F-Wastes   F-Waste Constituents Rep	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituen	t/Waste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
F-Waste F002 F003	s Reported to NCAR:	10,907 420	19,705 0	32,754 0	
	Waste Total	11,327	19,705	32,754	
Facility 0	00214				
F-Waste	Constituents Reported to TRI:				
ACET	ONE	0	0	0	
n-BU	TYL ALCOHOL	0	0	0	
METH METH	ANOL	0	0	0	
TOLU	ENE	õ	õ	Ő	
XYLE	NE	0	0	0	
	Constituent Total	0	0	0	
F-Waste	s Reported to NCAR:				
F003		78,860	0	0	
F003 F005	F005	18,686	40,400	37,200	
•	- Waste Total	97,546	40,400	37,200	
Facility 0	00217				
F-Waste	Constituents Reported to TRI:				
FREC	N 113	0	0	0	
METH	ANOL	0	0	0	
1,1,	1-TRICHLOROETHANE -	0	0	0	
	Constituent Total	0	0	0	
F-Waste F002	s Reported to NCAR:	57,758	13,000	5,481	
	- Waste Total	57,758	13,000	5,481	
Facility 0	00218				
F-Waste	Constituents Reported to TRI:	٥	٥	0	
P REC METH		õ	õ	õ	
1,1,	1-TRICHLOROETHANE	Ō	Ō	Ō	
	- Constituent Total	0		0	
	- Dependent to NCAR.				
F-waste	F002	0	9.368	11.778	
F002		20,681	17,860	9,984	
	- Waste Total	20,681	27,228	21,762	
Facility 0	000225				
F-Waste	Constituents Reported to TRI.				
1,1,	1-TRICHLOROETHANE	0	2,752	3,327	
	Constituent Total	0	2,752	3,327	
F-Waste	es Reported to NCAR:				
F001		64,527	49,275	0	
F003		7,200	2,400	0	
F003	J	0,000	v		

07/23/93Comparison of F-Wastes Reported to NCAR andPage 22F-Waste Constituents Reported to TRI: 1988 to 1			
Constituent/Waste Code	Lbs Transferred Off-site - 1988 1989 19		
Waste Total	77,727	51,675	440
Facility 000229			
F-Waste Constituents Reported to TRI:			
ACETONE	0	0	0
METHYL ETHYL KETONE (MEK)	0	0	0
TOLUENE	Ő	Ō	0
XYLENE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:			
D001D008F003F005	0	0	75,600
F003F005	48.400	42,000	0
		42,000	75 600
Waste Total	48,400	42,000	/5,600
acility 000233			
F-Waste Constituents Reported to TRI:	0	0	0
ACETONE METHYL, FTHYL, KETONE (MEK)	ő	Ö	C
METHANOL	Ő	0	Ċ
TOLUENE	0	0	C
XYLENE	0	0	C 
Constituent Total	0	0	C
F-Wastes Reported to NCAR:			
D035F003F005	0	0	136,320
	140,197	132 210	(
F003F005			
Waste Total	140,197	132,210	136,320
acility 000235			
F-Waste Constituents Reported to TRI:	0	4.0	,
	0	40	
F005058	ő	90	(
- Waste Total	0	140	(
Facility 000236			
F-Waste Constituents Reported to TRI:			
n-BUTYL ALCOHOL	0	0	(
CYCLOHEXANE	0	0	(
METHYL ETHYL KETONE (MEK) TOLUENE	0 0	0 0	(
- Constituent Total	0	0	
E-Wastes Reported to NCAD.			
F002	2,440	0	(
F005	159,497	114,538	86,609
- Waste Total	161,937	114.538	86,609

07/23/93 Page 23	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Waste	e Code	Lbs Tr 1988	ansferred O 1989	ff-site 1990
Facility 000244				
F-Waste Const:	ituents Reported to TRI	:	0	٥
METHYL ETHY	YL KETONE (MEK)	ŏ	ŏ	0
METHANOL		0	0	0
XYLENE		0	0	0
	Constituent Total	0	0	0
F-Wastes Repo	rted to NCAR:			
D001D008F0	03F005	0	0 29,200	38,094
F005		30,471	0	0
	Waste Total	30,471	29,200	38,094
Facility 000247				
F-Waste Const	ituents Reported to TRI	: 55 700	70 217	1 461
ACETONE		55,722	/9,34/	1,401
	Constituent Total	55,722	79,347	1,461
F-Wastes Repo	rted to NCAR:	0	0	115
D002F003		Ő	800	0
F001		5 900	7 810	0
F002		199,684	204,153	88,400
F003F005		0	440	0
F004 F005		4,440	120	0
	Waste Total	210,104	211,430	88,515
Facility 000250				
F-Waste Const	ituents Reported to TRI	:	10 205	2 160
1,1,1-TRIC	HLOROETHANE	16,215	10,285	2,160
	Constituent Total	16,215	10,285	2,160
F-Wastes Repo F001	rted to NCAR:	6,600	0	560
F001F002		0	8,415	3,645
F003 F003F005		5,325	25,712	0 14,958
	Waste Total	11,925	34,127	19,163
Facility 000251				
F-Waste Const	ituents Reported to TRI	:		
n-BUTYL AL	COHOL	29,094	0	0
NITROBENZE	NE	24,530	7,146	6,850
TOLUENE		19,020	5,490	2,750
	Constituent Total	105,854	13,386	9,850
F-Wastes Repo	rted to NCAR:	^	^	100
F002		0	240	490

07/23 Page	/93 24	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituont/Wast	a Code	Lbs Tr 1988	ansferred Ofi 1989	E-site 1990
	F002F005		0	0	240
	F003		58,187	860	180
	F003F005 F005		64,456	39,134	122,282 <sup>.</sup>
		Waste Total	122,643	244,554	123,312
Facil	ity 000253				
F-	Waste Const	ituents Reported to TRI	:		
	ACETONE		0	0	0
	METHYL ETHY	YL KETONE (MEK)	830		
		Constituent Total	830	0	0
F-'	Wastes Repo	rted to NCAR:			
	D001F003		4 200	737	7,315
	F002 F003		14,200	737	0
	F003F005		0	1,320	3,828
		Waste Total	18,517	2,794	11,143
Facil	ity 000255				
_		the set a Demonstral to MDI	r.		
F	METHYL ETH	YL KETONE (MEK)	75,750	64,000	500
		Constituent Total	75,750	64,000	500
	Wastas Bono	rtod to NCAR.			
r -	FOO1	ILEG CO NCAR:	161,570	1,964	1,727
	F002		4,400	800	14,500
	F003F005		0	0 20 601	5,013 4 320
	FUUS		165 070		
		Waste Total	105,970	23,305	25,560
Facil	ity 000257				
F-	Waste Const TETRACHLOR	ituents Reported to TRI OETHYLENE	[: 0	0	3,713
		Constituent Total	0	0	3.713
-	Marker Demo	wheel the NGNR.	-		- •
F	Wastes Repo	05	0	0	33,000
	F001		Ō	0	4,900
	F003		402	0	0
		Waste Total	402	0	37,900
Facil	ity 000258				
F-	Waste Const	ituents Reported to TRI	Ι:		
	ACETONE		7,102	8,465	1 007
	n-BUTYL AL	COHOL	14,203	U A	1,007 1 074
	METHIL ETH METHINOI	IL RETONE (MER)	0	8,465	250
	METHYL ISO	BUTYL KETONE	Ō	1,679	0
	TOLUENE		10,340	11,395	2,367
	XYLENE		7,030	10,403	250
		Constituent Total	38,675	40,407	5,800

Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Transferred Off-site		-site
Constituent/Waste Code	1988 	1989	1990
F-Wastes Reported to NCAR: D001D007F003F005 F003F005 F005	0 0 138,862	0 84,650 0	110,563 0 0
	138,862	84,650	110,563
Facility 000260			
F-Waste Constituents Reported to TRI:			
ACETONE	6,283	7,716	0
n-BUTYL ALCOHOL METHYI, ETHYI, KETONE (MEK)	14,557	0	0
METHANOL	0	7,716	250
TOLUENE	10,598	10,588	1,263
XYLENE		<i>2,372</i>	
Constituent Total	31,438	28,392	2,513
F-Wastes Reported to NCAR:	0	0	26 914
D001F005 F003F005	0	77,167	36,814
F005	118,783	0	0
Waste Total	118,783	77,167	36,814
Facility 000261			
F-Waste Constituents Reported to TRI:			
ACETONE	5,713	0	0
METHYL ETHYL KETONE (MEK) METHANOI.	0	0	0
METHYL ISOBUTYL KETONE	3,955	0	0
TOLUENE	8,332	0	0
- XYLENE	5,005		
Constituent Total	23,665	0	0
F-Wastes Reported to NCAR:	66 050	0	٥
F005 -			
Waste Total	66,950	0	0
Facility 000262			
F-Waste Constituents Reported to TRI:	2 012	2 101	0
ACETONE D-BUTYL ALCOHOL	3,813	3,101	750
METHYL ETHYL KETONE (MEK)	0	0	1,855
METHANOL	2 2 2 0	3,101	250
METHYL ISOBUTYL KETONE TOLUENE	5,209	6,408	2,328
XYLENE	3,782	4,645	250
- Constituent Total	16,133	18,567	5,433
F-Wastes Reported to NCAR:			
D001D007F003F005	0	0	35,476
F003F005 F005	36,875	31,010	0
-		21 010	2E 476
Waste Total	36,8/5	31,010	37,4/0

07/23/93 Page 26	<ul> <li>93 Comparison of F-Wastes Reported to NCAR and</li> <li>26 F-Waste Constituents Reported to TRI: 1988 to 1990</li> </ul>				
Constituent/Wast	e Code	Lbs Tra 1988	nsferred Off 1989	f-site 1990	
Facility 000263					
F-Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL METHYL ISO TOLUENE XYLENE	ituents Reported to TRI: COHOL YL KETONE (MEK) BUTYL KETONE	0 0 0 0 0 0	0 0 0 0 0 0	0 0 0 0 0 0	
	Constituent Total	0	0	0	
F-Wastes Repo F005	rted to NCAR:	18,800	21,800	125,200	
	Waste Total	18,800	21,800	125,200	
Facility 000264					
F-Waste Const CRESOL(S)	ituents Reported to TRI:	0	0	0	
	Constituent Total	0	0	0	
F-Wastes Repo F002 F003 F003F005 F005	rted to NCAR:	0 51,800 0 0	0 0 9,091 0	7 0 0 2,919	
	Waste Total	51,800	9,091	2,926	
Facility 000270					
F-Waste Const ACETONE TETRACHLOR	ituents Reported to TRI: OETHYLENE	0 0	0 0	0 0	
	Constituent Total	0	0	0	
F-Wastes Repo F003	rted to NCAR:	0	500	0	
	Waste Total	0	500	0	
Facility 000273					
F-Waste Const 1,1,1-TRIC	ituents Reported to TRI: HLOROETHANE	0	0	0	
	Constituent Total	0	0	0	
F-Wastes Repo F001 F003	rted to NCAR:	85,518 0	81,760 3,620	73,319 3,575	
	- Waste Total	85,518	85,380	76,894	
Facility 000275					
F-Waste Const ACETONE CHLOROBENZ FTHYLBENZE	ituents Reported to TRI: ENE NE	0 0 0	0 0 750	0 0 0	

## Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

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Construct Wester Orde	Lbs 1	ransferred (	Off-site
Constituent/Waste Code		1989	1990
METHANOL XYLENE	59,900 3,315	24,000 1,000	0 0
Constituent Total	63,215	25,750	0
F-Wastes Reported to NCAR: D001F003 D001F003F005 F002 F002F003F005 F003 F005	0 4,486 0 528,297 1,800 	3,000 0 163,387 0 166,387	28,380 450 0 3,902,160 84,660 0 
Facility 000277	551,505	100,507	1,010,000
Facility 0002//			
F-Waste Constituents Reported to TRI: ACETONE	520	7,560	0
Constituent Total	520	7,560	0
F-Wastès Reported to NCAR: F002 F003	1,350 75,960	0 69,269	0 15,937
Waste Total	77,310	69,269	15,937
Facility 000278			
F-Waste Constituents Reported to TRI: FREON 113 METHYLENE CHLORIDE	0 0	0 0	0 0
- Constituent Total	0	0	0
F-Wastes Reported to NCAR: D001F002 F002	0	<b>4</b> ,215 12,886	487 0
Waste Total	0	17,101	487
Facility 000279			
F-Waste Constituents Reported to TRI:			
FREON 113 1,1,1-TRICHLOROETHANE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR:	15 050	26.400	44.000
F001 F002	11,495	26,400	44,226
F005	0	824	
Waste Total	27,445	27,224	44,226
Facility 000280			
F-Waste Constituents Reported to TRI: n-BUTYL ALCOHOL CYCLOHEXANE METHYLENE CHLORIDE	990 0 0	1,590 0 0	2,190 0 5,100
- Constituent Total	990	1,590	7,290

07/23 Page	/93 28	Comparison of F-Wastes F-Waste Constituents H	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wa	aste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
F-1	Wastes R D001F00 D001F00 F001 F001F00 F001F00 F002 F002F00 F003	eported to NCAR: 1 1F003 2F003 3 3F005	0 0 0 18,025 0 452	0 18,052 5,580 240 0 4,191 0 240	3,440 13,762 2,027 0 5,404 4,645 16,473 430	
		Waste Total	18,477	28,303	46,181	
Facil	ity 0002	81				
F-	Waste Co METHANO 1,1,1-T	nstituents Reported to TR L RICHLOROETHANE	I: 0 	750	755 0	
		Constituent Total	0	750	755	
F-	Wastes R •F001	eported to NCAR:	0	0	910	
		Waste Total	0	0	910	
Facil	ity 0002	87				
F-	Waste Co METHYLE 1,1,1-T	nstituents Reported to TR NE CHLORIDE RICHLOROETHANE	I: 2,000 0	7,820 0	<b>4</b> ,860 0	
		Constituent Total	2,000	7,820	4,860	
F-	Wastes R F002	eported to NCAR:	0	50,675	27,690	
		Waste Total	0	50,675	27,690	
Facil	ity 0002	89				
F-	Waste Co ACETONE CYCLOHE METHANO TETRACH	nstituents Reported to TR XANE L LOROETHYLENE	I: 14,892 53,337 41,933 66,350	14,800 0 0 0	0 0 0 0	
		Constituent Total	176,512	14,800	0	
F-	Wastes R D001F00 D001F00 D001F00 D002F00 D022D02 D035F00 F002 F002F00 F003	eported to NCAR: 2 3F005 3F005U159 3U002U213 2 7D040F002 3F005 3F005	0 0 0 0 101,600 0 136,380	0 0 0 0 0 70,480 0 102,733	$ \begin{array}{r} 160\\ 44,896\\ 675\\ 50\\ 59,200\\ 240\\ 4,020\\ 10,419\\ 38,745\\ 0 \end{array} $	
		Waste Total	237,980	173,213	158,405	

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07/23/93 Page 29	F-Waste Constituents Rep	orted to TR	I: 1988 to 1	990
Constituent/Waste	- Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
Facility 000298				
F-Waste Consti D001F003F00	tuents Reported to TRI: 5	0	8,500	3,713
	 Waste Total	0	8,500	3,713
Facility 000300				
F-Waste Consti XYLENE	tuents Reported to TRI:	0	0	0
	Constituent Total	0	0	0
F-Wastes Repor D001F003 D001F003F00 F003	ted to NCAR:	0 0 34,000	26,000 0 0	5,450 30,800 0
	 Waste Total	34,000	26,000	36,250
Facility 000303				
F-Waste Consti 1,1,1-TRICE	ituents Reported to TRI: HLOROETHANE	0	0	0
	Constituent Total	0	0	0
F-Wastes Repor D001F001 D008F003F00 F001 F003 F005	rted to NCAR: 05  Waste Total	0 2,420 0 0 2,420	0 2,736 317 379 3,432	708 709 2,260 0 0 
Facility 000311				·
F-Waste Const: METHANOL	ituents Reported to TRI:	0	0	0
	Constituent Total	0	0	0
F-Wastes Repo: F001 F001F005 F002 F003 F004 F005	rted to NCAR:	0 0 0 0 0	0 937 670 380 252 740 2,979	1,028 0 439 0 1,380
Escility 000214	habbe redar	·	_,	_,
F-Wasto Const	ituants Penorted to TRI.			
CYCLOHEXAN METHYL ETH TOLUENE	E YL KETONE (MEK)	0 0 0	0 0 3,78 <b>4</b>	250 250 0
	 Constituent Total	0	3,784	500
F-Wastes Repo D001F005	rted to NCAR:	0	6,186	0

07/23 Page	/93 30	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	ce Code	Lbs Tra 1988	ansferred Off 1989	-site 1990
	F002 F002F003 F003F005		2,424 0 0	3,017 0 0	250 170 93,296
		Waste Total	2,424	9,203	93,716
Facil	ity 000315				
F-	Waste Const ACETONE n-BUTYL AI	cituents Reported to TRI: LCOHOL	0 0	0 0	0 0
	METHANOL METHYL ISC TOLUENE XYLENE	DBUTYL KETONE	0 0 6,180 0	0 0 0 0	0 0 220 0
		- Constituent Total	6,180	0	220
F-	Wastes Repo D001F003 D001F003F( D001F003F( D001F003U( D001F005 F003 F005	orted to NCAR: 005 005U077 077 - Waste Total	0 0 0 108,425 853,845  962,270	480,080 211,220 3,920 0 146,740 2,380 1,850 	36,260 64,620 0 79,840 2,275 18,180 1,050 202,225
Facil	ity 000317				
F-	Waste Const ACETONE n-BUTYL AI METHYL ETH METHANOL TOLUENE XYLENE	Lituents Reported to TRI: LCOHOL HYL KETONE (MEK)	0 0 5,932 13,842 0	0 0 0 0 0 0	0 0 0 0 0
		- Constituent Total	19,774	0	0
F-	Wastes Repo D001D035F( F003F005 F005	orted to NCAR: 003F005	0 0 39,550	0 31,710 0	24,150 0 0
		Waste Total	39,550	31,710	24,150
Facil	ity 000327				
F-	Waste Const 1,1,1-TRIC TRICHLOROP	tituents Reported to TRI: CHLOROETHANE ETHYLENE	80 2,000	250 2,500	80 4,100
		- Constituent Total	2,080	2,750	4,180
F-	Wastes Repo D001F001 F001 F005	orted to NCAR:	0 174,532 0	8,960 173,565 750	7,940 94,990 445
		Waste Total	174,532	183,275	103,375

07/2 Page	3/93Comparison of F-Wastes Reported to NCAR and31F-Waste Constituents Reported to TRI: 1988 to 1990				1990
Cons	tituent/Wast	e Code	Lbs Tr 1988	ansferred Of 1989	f-site 1990
 Faci	lity 000328				
F	-Waste Const n-BUTYL AL METHYL ISC	ituents Reported to TRI: COHOL BUTYL KETONE	0	0	0
	1,1,1-TRIC	HLOROETHANE -	U		
		Constituent Total	0	0	0
F	Y-Wastes Repo D001D002F0 D002F003F0 F001 F002 F003 F003F005 F005	orted to NCAR: 03F005 05	0 0 0 1,000 0 0	0 900 30,380 1,550 0 1,400	400 500 0 400 2,000 1,600
		Waste Total	1,000	34,230	4,900
Faci	llity 000329 F-Waste Const METHANOL TOLUENE XYLENE	ituents Reported to TRI:	82 10,560 240	81 32,100 300	79 29,100 0
		Constituent Total	10,882	32,481	29,179
F	F-Wastes Repo F005	orted to NCAR:	98,500	159,000	162,875
		Waste Total	98,500	159,000	162,875
Faci	ility 000330				
I	F-Waste Const 1,1,1-TRIC	ituents Reported to TRI: CHLOROETHANE	0	0	0
		Constituent Total	0	0	0
I	F-Wastes Repo D001D027D( D001F002F( D001F003F( D002F003F( F003 F003F005 F005	orted to NCAR: 029F003 003F005 005	0 0 0 17,451 0 4,290	0 350 0 49,490 1,330 0	4,320 700 14,600 500 0 0
		Waste Total	21,741	51,170	20,120
Fac:	ility 000334				
]	F-Waste Const CRESOL(S) METHANOL TETRACHLOI XYLENE	cituents Reported to TRI	: 19,765 0 0	0 26,117 0 44,064	380 61,000 5,205
		Constituent Total	19,765	70,181	66,585
]	F-Wastes Repo F003 F005	orted to NCAR:	0 0	55,080 81,660	0 36,000

07/23/93 Page 32	Compar F-Wast	ison of 1 e Constit	F-Wastes tuents Re	tes Reported to NCAR and s Reported to TRI: 1988 to 1990			
Constitu	ent/Waste Code			Lbs Tr 1988	ansferred Of 1989	f-site 1990	
		Waste	- Total	0	136,740	36,000	
Facility	000335						
F-Was	te Constituents	Reported	d to TRI:				
AC	ETONE			0	0	0	
ME	THIL ETHIL KETO THANOL	NE (MER)		0	0	0	
TO	LUENE			õ	õ	Ő	
	Con	stituent	Total	0	0	0	
F-Was	tes Reported to	NCAR:		101 055	0	0	
FO	03 2005			121,055	116 760	0	
F0	05			ő	110,700	167,914	
		Waste	- Total	121,055	116,760	167,914	
Facility	000336						
			_				
F-Ŵas TO	te Constituents LUENE	Reported	d to TRI:	0	0	0	
	Con	stituent	Total	0	0	0	
F-Was	tes Reported to	NCAR:					
F0	01			3,000	3,600	0	
F0	05		_	0	0	184,980	
		Waste	Total	3,000	3,600	184,980	
Facility	000337						
F-Was	te Constituents	Reported	d to TRI:				
FR	EON 113			0	0	0	
XY	LENE			0	0	0	
	Con	stituent	- Total	0	0	0	
E-Was	tos Poported to	NCAP.					
D0	01F003F005	NCAN:		0	27,710	27,350	
FO	01			64,879	10,982	1,200	
F0	02			3,314	3,599	5,500	
FO	03		-	6,108			
		Waste	Total	74,301	42,291	34,050	
Facility	000342						
F-Was	te Constituents	Reported	i to TRI:				
CY	CLOHEXANE		-	0	0	0	
	Con	stituent	Total	0	0	0	
F-Was	tes Reported to	NCAR:					
FO	01			0	3,340	0	
FO	02		-	U 	1,920	0	
		Waste	Total	0	5,260	0	

- with

## Facility 000346

F-Waste Constituents Reported to TRI:

Lis Transferred Off-site	07/23/93 Page 33	Comparison of F-Wa F-Waste Constituen	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
F005         500         1,040         760           Waste Total         500         1,040         760           Facility 000347         F-Waste Constituents Reported to TRI: ACETONE         0         0         0         0           P-Waste Constituents Reported to TRI: ACETONE         2,700         3,000         750           MDIFIL: METHIL: ACETONE         2,700         3,000         7,50           MDIFIL: METHIL: METHIL: ACETONE         0         12,800         5,255           13,000         10,600         7,350           Constituent Total         17,700         34,400         17,805           P-Wastes Reported to NCAR: D001D035F003         0         0         7,755           Fool         S5,890         86,021         80,550           Facility 000350         Constituent Total         950         3,311         0           F-Wastes Reported to NCAR: F001         800         0         0         0           F-Waste Constituents Reported to TRI: FRECN 113         800         0         0         0           Fool         Constituent Total         950         3,311         0         0           F-Waste Constituents Reported to TRI: FRECN 113         0         0         0 <t< th=""><th>Constitu</th><th>ent/Waste Code</th><th> Lbs 1 1988</th><th>Fransferred C 1989</th><th>ff-site 1990</th></t<>	Constitu	ent/Waste Code	Lbs 1 1988	Fransferred C 1989	ff-site 1990	
Waste Total         500         1,040         760           Facility 000347         F-Waste Constituents Reported to TRI: ACETOME         0         0         0         0           F-Waste Constituents Reported to TRI: ACETOME         2,700         3,000         750           METHYL ALCHORL         2,700         3,000         750           METHYL ETHYL KETONE (MEK)         2,000         8,000         4,450           TOLUEME         13,000         10,660         7,350           Constituent Total         17,700         34,400         17,805           DO1D035F003         0         0         77,750           D01F003         55,890         86,021         0           F-Wastes Reported to NCAR:         950         3,311         0           F-Waste Constituents Reported to TRI:         1,1,1-TRICHLOROETHANE         950         3,311         0           F-Wastes Reported to NCAR:         950         3,311         0         0         0           F-Wastes Reported to NCAR:         950         3,311         0         0         0           F-Waste Constituents Reported to TRI:         950         3,311         0         0         0           F-Wastes Constituents Reported to TRI: <t< th=""><th> F(</th><th></th><th>500</th><th>1.040</th><th>760</th></t<>	 F(		500	1.040	760	
Facility 000347         F-Waste Constituents Reported to TRI: ACETOME (MEK)       2,700       3,000       750         METHYL ALCHOL       2,700       8,000       4,450         TOLUENE       0       0       0       7,255         1,1,1-TRICHLOROETHANE       0       0       7,755         D001D035F003       0       0       7,750         D001D035F003       0       0       2,800         F-Wastes Reported to NCAR:       0       0       2,800         D001D035F003       0       0       2,800         Foldstruct       55,890       86,021       0         -Waste Total       55,890       86,021       80,550         Facility 000350		Waste Tot	al 500	1,040	760	
F-Waste Constituents Reported to TRI:       0       0       0       0         ACETONE       2,000       3,000       750         METHYL ETNYL KETONE (MEK)       2,000       8,000       4,450         TOLUENE       0       12,800       5,255         1,1.1-TRICHLOROETHANE       0       0       0       7,300         Constituent Total       17,700       34,400       17,605         F-Wastes Reported to NCAR:       0       0       2,800       7,750         D001D035F003       0       0       0       2,800         F003       55,890       86,021       0       0         Waste Total       950       3,311       0       0         Constituent Total       950       3,311       0       0         F-Wastes Reported to NCAR:       950       3,311       0       0         Fo03       800       0       0       0       0         Fo11       1,600       0<	Facility	000347				
ACETONE 0 0 0 0 0 n-BUTYL ALCOHOL 2,700 3,000 750 METHYL ETHYL KETONE (MEK) 2,000 8,000 4,450 TOLUENE 0 12,800 5,255 1,1,1-TRICHLOROETHANE 0 12,800 17,805 Constituent Total 17,700 34,400 17,805 F-Wastes Reported to NCAR: D001D035F003 0 0 0 7,755 D001F003 55,890 86,021 80,550 F-Waste Constituents Reported to TRI: 1,1,1-TRICHLOROETHANE 955 3,311 0 F-Wastes Reported to NCAR: F001 503 800 0 0 F-Wastes Reported to NCAR: F001 0 0 0 0 For a structure total 950 3,311 0 F-Wastes Reported to NCAR: F001 0 0 F-Waste Constituents Reported to TRI: 1,1,1-TRICHLOROETHANE 950 3,311 0 F-Wastes Reported to NCAR: F001 0 0 Facility 000353 F-Waste Constituents Reported to TRI: FEEDN 113 0 F-Waste Constituents Reported to TRI: FEEDN 113 0 F-Waste Constituents Reported to TRI: FEEDN 113 0 F-Waste Reported to NCAR: F005 0 0 0 TOLUENE 0 F-Waste Reported to NCAR: F005 0 0 TOLUENE 0 F-Waste Reported to NCAR: F005 0 0 Waste Total 0 F-Waste Reported to NCAR: F005 0 TOLUENE 0 F-Waste Constituent Total 0 F-Waste Reported to NCAR: F005 0 TOLUENE 0 TOLUENE 0 TOLUENE 0 F-Waste Reported to NCAR: F005 0 TOLUENE 0 F-Waste Constituent Reported to TRI: FREON 113 0 METHYL ETHYL KETONE (MEK) 385,728 0 TOLUENE 0 TOLUENE 0 TOLUENE 0 TOLUENE 0 TOLUENE 0 TOLUENE 0 F-Waste Reported to NCAR: F005 0 METHYL KETONE (MEK) 117,640 3,657 0 METHYL ETHYL KETONE (MEK) 117,640 3,657 0 METHYL CHYL KETONE (MEK) 117,780 37,733 130,3	F-Was	te Constituents Reported to	TRI:			
H-BUTL ALCONCL TOLUENE         2,000 0         3,000 12,800 0         4,450 5,255 0           TOLUENE         0         12,800 0         5,255 0           I, 1, 1-TRICHLOROETHANE         0         0         7,350 0           Constituent Total         17,700         34,400         17,805           F-Wastes Reported to NCAR: D0010355003         0         0         77,750 0         2,800 0         0         2,800 0         0         0         2,800 0         0	AC	ETONE	0	0	0	
TOLLENE         0         12,800         5,255           1,1-TRICHLOROETHANE         0         0         0           XYLENE         13,000         10,600         7,350           Constituent Total         17,700         34,400         17,805           F-Wastes Reported to NCAR:         0         0         77,750           D001D0357003         0         0         2,800           F003         55,890         86,021         0           Waste Total         55,890         86,021         0           Facility 000350	n- MH	THYL ETHYL KETONE (MEK)	2,700	8,000	4,450	
1.1.1-TRICHLOROFITANE       13,000       10,600       7,350         Constituent Total       17,700       34,400       17,805         F-Wastes Reported to NCAR:       0       0       7,750         D001D035F003       0       0       2,800         F003       55,890       86,021       0         Waste Total       55,890       86,021       80,550         Facility 000350       F-Waste Constituents Reported to TRI:       1,1,1-TRICHLOROETHANE       950       3,311       0         F-Wastes Reported to NCAR:       0       0       0       0       0         F001       800       0       0       0       0         F001       800       0       0       0       0         Fordate Constituents Reported to TRI:       FRENN 113       0       0       0         FWaste Reported to NCAR:       1,600       0       0       0         Waste Total       1,600       0       0       0         FWastes Reported to NCAR:       11,600       0       0       0         FWaste Stituents Reported to TRI:       FREON 113       0       0       0       0         F-Wastes Reported to NCAR:       17,640	TC		0	12,800	5,255	
Constituent Total 17,700 34,400 17,805 P-Wastes Reported to NCAR: D001D035F003 0 0 0 2,800 F003 55,890 86,021 0 Waste Total 55,890 86,021 0 Waste Total 55,890 86,021 0 F-Waste Constituents Reported to TRI: 1,1,1-TRICHLOROETHANE 550 3,311 0 F-Wastes Reported to NCAR: F001 800 0 0 0 F003 800 0 0 F-Wastes Reported to NCAR: FREON 113 0 ToLUENE 191,032 404,060 0 ToLUENE 191,032 404,060 0 F-Wastes Reported to NCAR: F005 0 F-Waste Streported to NCAR: F005 0 ToLUENE 191,032 404,060 0 F-Wastes Reported to NCAR: F005 0 ToLUENE 191,032 404,060 0 F-Wastes Reported to NCAR: F005 0 ToLUENE 191,032 404,060 0 ToLUENE 191,032 404,060 0 ToLUENE 191,032 404,060 0 F-Wastes Reported to NCAR: F005 0 ToLUENE 191,032 404,060 0 ToLUENE 191,032 191,785 104,307 0 Waste Total 162,920 41,628 130,353 104,307 0 Waste Total 297,357 191,785 104,307 0 Waste Total 297,357 191,785 104,307 0 Waste Total 297,357 191,785 104,307 0 Total 191,785 104	, T X	LENE	13,000	10,600	7,350	
F-Wastes Reported to NCAR: D001D03SF003 F003       0       0       77,750 0         Waste Total       55,890       86,021       0         Waste Total       55,890       86,021       0         Waste Total       55,890       86,021       0         Facility 000350       F-Waste Constituents Reported to TRI: 1,1,1-TRICHLORCETHANE       950       3,311       0         Constituent Total       950       3,311       0       0       0         F-Wastes Reported to NCAR: F001       800       0       0       0         Waste Total       1,600       0       0       0         F-Wastes Constituents Reported to TRI: FREON 113       0       0       0       0         Maste Total       1,600       0       0       0       0         F-Wastes Constituents Reported to TRI: FREON 113       0       191,032       404,060       0       0         Constituent Total       576,760       -464,188       0       0       0       0         F-Wastes Reported to NCAR: F005       0       37,325       36,058       0       0       0         Generative Maste       117,640       3,657       0       0       0       0         Mast		Constituent Tot	al 17,700	34,400	17,805	
D001D035F003         0         0         77,750           D001F003         0         2,800         0         2,800           F003	F-Was	stes Reported to NCAR:				
D001F003 F003 F003 Waste Total 55,890 B6,021 B0,550 Facility 000350 F-Waste Constituents Reported to TRI: 1,1,1-TRICHLOROETHANE Constituent Total 950 F-Wastes Reported to NCAR: F001 F003 B00 F-Wastes Reported to NCAR: F001 Waste Total 1,600 F-Waste Constituents Reported to TRI: FRON 113 Constituent Total 576,760 F-Wastes Reported to NCAR: F005 F-Waste Constituents Reported to TRI: FRON 113 Constituent Total 576,760 F-Wastes Reported to NCAR: F005 F-Waste Constituents Reported to TRI: FRON 113 Constituent Total 576,760 METHYL ETHYL KETONE (MEK) Maste Total 0 TOLUENE F-Wastes Reported to NCAR: F005 F-Waste Constituents Reported to TRI: FRON 113 Constituent Total 576,760 METHYL ETHYL KETONE (MEK) Maste Total 0 TOLUENE F-Wastes Constituents Reported to TRI: FRON 113 Constituent Total 576,760 METHYL ETHYL KETONE (MEK) Maste Total 0 TOLUENE Constituent Total 162,920 H1,628 130,353 F-Wastes Reported to NCAR: F001 METHYL ETHYL KETONE (MEK) METHYNOL Constituent Total 162,920 METHYNOL Constituent Total 162,920 METHYNOL Constituent Total 162,920 METHYNOL Constituent Total 162,920 Maste Total 297,357 J91,785 J04,307 F-69	D	01D035F003	0	0	77,750	
Waste Total         55,890         86,021         80,550           Facility 000350         F-Waste Constituents Reported to TRI:         1,1,1-TRICHLOROETHANE         950         3,311         0           Constituent Total         950         3,311         0         0           F-Wastes Reported to NCAR:         500         0         0         0           F001         800         0         0         0           Waste Total         1,600         0         0         0           FREON 113         0         0         0         0         0           MeTHYL ETHYL KETONE (MEK)         385,728         60,128         0         0         0           Constituent Total         576,760         464,188         0         <	D( F(	001F003 003	55,890	86,021	2,800	
Facility 000350         F-Waste Constituents Reported to TRI:         1,1,1-TRICHLOROETHANE       950         Constituent Total       950         950       3,311         0         F-Wastes Reported to NCAR:         F001       800         F003       800         Waste Total       1,600         0       0         Waste Total       1,600         FREON 113       0         0       0         METHYL ETHYL KETONE (MEK)       385,728         60,128       0         0       0         0       0         0       0         0       0         0       0         METHYL ETHYL KETONE (MEK)       385,728         60,128       0         0       0         0       37,325         36,058         F-Wastes Reported to NCAR:         F005       0         0       37,325         36,058         Facility 000357         F-Waste Constituents Reported to TRI:         FREON 113       7,500         0       238         0		Waste Tot	al 55,890	86,021	80,550	
F-Waste Constituents Reported to TRI:       950       3,311       0         Constituent Total       950       3,311       0         F-Wastes Reported to NCAR:       950       3,311       0         F001       800       0       0         Waste Total       1,600       0       0         Facility 000353       0       0       0         F-Waste Constituents Reported to TRI:       0       0       0         FREON 113       0       0       0       0         METHYL ETHYL KETONE (MEK)       385,728       60,128       0         TOLUENE       191,032       404,060       0       0         Constituent Total       576,760       464,188       0         F-Wastes Reported to NCAR:       0       37,325       36,058         Facility 000357	Facility	• • • • • • • • • • • • • • • • • • • •				
P-Waste Constituents Reported to TK1:       950       3,311       0         Constituent Total       950       3,311       0         F-Wastes Reported to NCAR:       800       0       0         F001       800       0       0         Waste Total       1,600       0       0         Facility 000353            Freen 113       0       0       0       0         METHYL ETHYL KETONE (MEK)       385,728       60,128       0         ToLUENE       191,032       404,060       0       0         Constituent Total       576,760       464,188       0       0         F-Wastes Reported to NCAR:       0       37,325       36,058         Facility 000357						
Constituent Total         950         3,311         0           F-Wastes Reported to NCAR: F001         800         0         0         0           Waste Total         1,600         0         0         0           Waste Total         1,600         0         0         0           Facility 000353         0         0         0         0         0           FREON 113         0         0         0         0         0         0           METHYL ETHYL KETONE (MEK)         385,728         60,128         0         0         0           TOLUENE         191,032         404,060         0         0         0         0           Constituent Total         576,760         464,188         0	F-Was 1,	te Constituents Reported to 1,1-TRICHLOROETHANE	950 950	3,311	0	
F-Wastes Reported to NCAR:       800       0       0         F001       800       0       0         Waste Total       1,600       0       0         Pacility 000353       0       0       0       0         FECN 113       0       0       0       0         METHYL ETHYL KETONE (MEK)       385,728       60,128       0         TOLUENE       191,032       404,060       0         Constituent Total       576,760       464,188       0         F-Wastes Reported to NCAR:       0       37,325       36,058         Facility 000357       0       0       0       0         FREON 113       7,500       0       0       0         Maste Total       0       37,325       36,058         Facility 000357       136,058       0       0       0         F-Waste Constituents Reported to TRI:       7,500       0       0       0         METHANOL       238       0       0       0       238       0         TOLUENE       37,780       37,733       130,353       0       0       0       130,353         F-Wastes Reported to NCAR:       8,525       0		Constituent Tot	al 950	3,311	0	
F001 F003       800 Waste Total       0 1,600       0 0       0 0         Facility 000353       F-Waste Constituents Reported to TRI: FREON 113 TOLUENE       0 0       0 0       0 0       0 0       0 0         F-Waste Constituents Reported to TRI: FREON 113 TOLUENE       191,032       404,060 0       0 0       0 0       0 0       0 0         F-Wastes Reported to NCAR: F005       0 0       37,325       36,058 0	F-Was	stes Reported to NCAR:		<u>^</u>	0	
Waste Total         1,600         0         0           Facility 000353         F-Waste Constituents Reported to TRI: FREON 113         0         0         0         0           METHYL ETHYL KETONE (MEK)         385,728         60,128         0         0         0           TOLUENE         191,032         404,060         0         0         0         0           Constituent Total         576,760         464,188         0         0         37,325         36,058           F-Wastes Reported to NCAR: F005         0         37,325         36,058         36,058           Waste Total         0         37,325         36,058         36,058           Facility 000357	F( F(	)03	800	0	0	
Facility 000353         F-Waste Constituents Reported to TRI: FREON 113       0       0       0         METHYL ETHYL KETONE (MEK)       385,728       60,128       0         TOLUENE       191,032       404,060       0         Constituent Total       576,760       464,188       0         F-Wastes Reported to NCAR: F005       0       37,325       36,058         Waste Total       0       37,325       36,058         Facility 000357       0       0       0         F-Waste Constituents Reported to TRI: FREON 113       7,500       0       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHYNU ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F001       8,525       0       0       0         F005       288,832       191,785       104,307		Waste Tot	al 1,600	0	0	
F-Waste Constituents Reported to TRI: FREON 113       0       0       0         METHYL ETHYL KETONE (MEK)       385,728       60,128       0         TOLUENE       191,032       404,060       0         Constituent Total       576,760       464,188       0         F-Wastes Reported to NCAR: F005       0       37,325       36,058         Waste Total       0       37,325       36,058         Facility 000357	Facility	000353				
FREDN 113       0       0       0       0         METHYL ETHYL KETONE (MEK)       385,728       60,128       0         TOLUENE       191,032       404,060       0         Constituent Total       576,760       464,188       0         F-Wastes Reported to NCAR:       0       37,325       36,058         Fo05       0       37,325       36,058         Waste Total       0       37,325       36,058         Facility 000357            FREON 113       7,500       0       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F001       8,525       0       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	F-Was	ste Constituents Reported to	TRI:			
MEININ FINIL KETONE (MEK)       191,032       404,060       0         TOLUENE	FI	REON 113	0 385 728	0 60 128	. 0	
Constituent Total 576,760 464,188 0 F-Wastes Reported to NCAR: F005 0 37,325 36,058 Waste Total 0 37,325 36,058 Facility 000357 F-Waste Constituents Reported to TRI: FREON 113 7,500 0 0 METHYL ETHYL KETONE (MEK) 117,640 3,657 0 METHANOL 0 238 0 TOLUENE 37,780 37,733 130,353 Constituent Total 162,920 41,628 130,353 F-Wastes Reported to NCAR: F001 60 288,832 191,785 104,307 Waste Total 297,357 191,785 104,307 F-68	T	DLUENE	191,032	404,060	ő	
F-Wastes Reported to NCAR: F005       0       37,325       36,058         Waste Total       0       37,325       36,058         Facility 000357       0       37,500       36,058         F-Waste Constituents Reported to TRI: FREON 113       7,500       0       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F001       8,525       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307		Constituent Tot	al 576,760	464,188	0	
F005       0       37,325       36,058         Waste Total       0       37,325       36,058         Facility 000357       0       37,325       36,058         Facility 000357       7,500       0       0       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F001       8,525       0       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	F-Was	stes Reported to NCAR:				
Waste Total       0       37,325       36,058         Facility 000357         F-Waste Constituents Reported to TRI:         FREON 113       7,500       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	F	)05		37,325	36,058	
Facility 000357 F-Waste Constituents Reported to TRI: FREON 113 7,500 0 0 METHYL ETHYL KETONE (MEK) 117,640 3,657 0 METHANOL 0 238 0 TOLUENE 37,780 37,733 130,353 Constituent Total 162,920 41,628 130,353 F-Wastes Reported to NCAR: F001 8,525 0 0 F005 288,832 191,785 104,307 Waste Total 297,357 191,785 104,307 E-68		Waste Tot	al 0	37,325	36,058	
F-Waste Constituents Reported to TRI:         FREON 113       7,500       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:         F001       8,525       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	Facility	y 000357				
FREON 113       7,500       0       0         METHYL ETHYL KETONE (MEK)       117,640       3,657       0         METHANOL       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:         F001       8,525       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	F-Was	ste Constituents Reported to	TRI:	0	0	
METHANOL TOLUENE       0       238       0         TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F001       8,525       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	F1 M	EON 113 ETHYL ETHYL KETONE (MEK)	7,500 117,640	3,657	0	
TOLUENE       37,780       37,733       130,353         Constituent Total       162,920       41,628       130,353         F-Wastes Reported to NCAR:       8,525       0       0         F001       8,525       0       0         F005       288,832       191,785       104,307         Waste Total       297,357       191,785       104,307	М	ETHANOL	0	238	0	
Constituent Total 162,920 41,628 130,353 F-Wastes Reported to NCAR: F001 F005 Waste Total 297,357 191,785 104,307 F-68	T	DLUENE	37,780	37,733	130,353	
F-Wastes Reported to NCAR: F001 F005 Waste Total 297,357 191,785 104,307 F-68		Constituent Tot	al 162,920	41,628	130,353	
F001 8,525 0 0 F005 288,832 191,785 104,307 Waste Total 297,357 191,785 104,307 F-68	F-Wa:	stes Reported to NCAR:		•	•	
Waste Total 297,357 191,785 104,307	F F	001 005	8,525 288,832	0 191,785	0 104,307	
waste 10tal 257,357 131,765 104,307 F_60			 al 207 257	 101 705	104 307	
		waste Tot	т_29	191,100	104,007	

07/23/93 Page 34	<ul> <li>Comparison of F-Wastes Reported to NCAR and</li> <li>F-Waste Constituents Reported to TRI: 1988 to 1990</li> </ul>			
Constituent	/Waste Code	Lbs Tra 1988	nsferred Off 1989	f-site 1990
Facility 00	0358			
F-Waste	Constituents Reported to TRI	:	1 000	
n-BUT METHY	YL ALCOHOL I ETHYI KETONE (MEK)	4,672	1,298	0
METHA	NOL	4,672	6,993	ŏ
TOLUE	NE	4,672	3,180	0
XYLEN	E	0	1,554	0
	Constituent Total	18,688	14,113	0
F-Wastes	Reported to NCAR:			
F005		52,400	46,900	43,268
	Waste Total	52,400	46,900	43,268
Facility 00	0360			
F-Waste	Constituents Reported to TRI	[:		
TOLUE	NE	6,364	358	0
XYLEN	E	5,162	288	
	Constituent Total	11,526	646	0
F-Wastes	Reported to NCAR:	40 450	•	•
F003 F003F	005	12,150	900	0
	Waste Total	12,150	900	0
	0261	,	2	-
Facility 00	0361			
F-Waste	Constituents Reported to TRI	[:		5 9 6 9
METHY METHA	L ETHYL KETONE (MEK)	0	3,229	5,363
TOLUE	INCE INE	2,888	3,229	5,363
XYLEN	E	0	0	0
	Constituent Total	2,888	6,458	10,726
F-Wastes	Reported to NCAR:			
D001F	003F005	0	10,764	17,878
	Waste Total	0	10,764	17,878
Eacility 00	0366			
ractifey ou				
F-Waste	Constituents Reported to TRI	[:	0	0
METHY	NE L ETHYL KETONE (MEK)	925	338	ŏ
METHA	NOL	0	787	0
TOLUE	NE .	1,094	2,002	0
	Constituent Total	3,250	3,127	0
F-Wastes D001F	Reported to NCAR:	0	0	25,948
20011				
	waste Total	U	U	20,948
Facility 00	0372			
F-Waste	Constituents Reported to TRI	I:	•	•
METHA	NOL	323	U	0

07/23/93 Page 35 Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

Constituent/Waste Code	Lbs Tra 1988	ansferred Of: 1989	f-site 1990
TOLUENE	15.604	2.970	3.400
Constituent Total	15,927	2,970	3,400
	13,527	2,510	5,400
D001D008F005 D001F005 F003 F005	0 0 43,724 107,550	0 0 138,322	67,490 88,580 0 0
Waste Total	151,274	138,322	156,070
Facility 000373			
F-Waste Constituents Reported to TRI FREON 113	: 0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: D001F002 D008F001 D008F002 F002 F003	0 0 971 486	0 0 1,650 1,834 917	4,462 1,043 0 570 0
Waste Total	1,457	4,401	6,075
Facility 000376			
F-Waste Constituents Reported to TRI FREON 113 1,1,1-TRICHLOROETHANE	: 0 0	0 0	0 1
Constituent Total	0	0	1
F-Wastes Reported to NCAR: D001F005 F001 F002 Waste Total	0 11,160 10,235	2,038 5,550 1,129	0 8,324 2,237
Facility 000277	21,393	0,111	10,501
F-Waste Constituents Reported to TRI ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	: 0 0 0 0 0 0	0 0 0 0 0 0	0 0 0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005	0 26,000	20,285 0	33,719 0
Waste Total	26,000	20,285	33,719
Facility 000379			
F-Waste Constituents Reported to TRI n-BUTYL ALCOHOL	: 0	0	0

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Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Tra	ansferred Off	-site
Constituent/Waste Code			
METHANOL TOLUENE	0 0	0 0	0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005	0 22,440	15,315 0	17,276 0
Waste Total	22,440	15,315	17,276
Facility 000380			
F-Waste Constituents Reported to TRI: ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	0 0 0 0 0 0	0 0 0 0 0 0	0 0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005 - Waste Total	0 28,600 28,600	27,200 0 	32,781 0  32,781
Facility 000381	-		
F-Waste Constituents Reported to TRI: ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	0 0 0 0 0 0	0 0 0 0 0	0 0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005	32,568	26,760	10,808
Waste Total	32,000	20,700	10,808
Facility 000382			
F-Waste Constituents Reported to TRI: ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	0 0 0 0 0 0	0 0 0 0 0 0	0 0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005	0 29,760	28, <b>4</b> 27 0	40,397 0
Waste Total	29,760	28,427	40,397

07/23/93Comparison of F-WasterPage 37F-Waste Constituents 1	s Reported to Reported to TR	NCAR and NI: 1988 to 1	.990
Constituent/Waste Code	Lbs Tra 1988	insferred Off 1989	-site 1990
Facility 000383			
F-Waste Constituents Reported to TR ACETONE n-BUTYL ALCOHOL	I: 0 0	0	0
METHIL ETHIL RETONE (MER) METHANOL TOLUENE XYLENE	0 0 0	0 0 0	0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005	0 62,864	57,805 0	117,502 0
Waste Total	62,864	57,805	117,502
Facility 000384			
F-Waste Constituents Reported to TR ACETONE METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	I: 0 0 0 0 0	0 0 0 0	0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003	0	0	26,698
Waste Total	0	0	26,698
Facility 000385			
F-Waste Constituents Reported to TR ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	I: 0 0 0 0 0 0	0 0 0 0 0	0 0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003 F005	0 54,808	50,185 0	66,115 0
Waste Total	54,808	50,185	66,115
Facility 000386			
F-Waste Constituents Reported to TR ACETONE n-BUTYL ALCOHOL METHANOL TOLUENE XYLENE	I: 0 0 0 0 0	0 0 0 0	0 0 0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003	0	6,685	0

07/23/93 Page 38	)7/23/93Comparison of F-Wastes Reported to NCAR andPage 38F-Waste Constituents Reported to TRI: 1988 to 1990				
Constituent	c/Waste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
 F005		1,320	0	0	
	Waste Total	1,320	6,685	0	
Facility 0(	00387			·	
F-Waste TETR	Constituents Reported to TRI ACHLOROETHYLENE	: 0	0	0	
	Constituent Total	0	0	0	
F-Waster F001	s Reported to NCAR:	16,821	0	0	
	Waste Total	16,821	0	0	
Facility 0	00388				
F-Waste METH	Constituents Reported to TRI YL ETHYL KETONE (MEK)	: 12,429	9,882	2,637	
•	Constituent Total	12,429	9,882	2,637	
F-Waste: D0011 D0011 F0011 F003 F003 F005	s Reported to NCAR: F003 F005 F003F005 F005	0 0 12,667 0 100,620	0 3,720 3,520 0 16,000 25,640	2,752 0 0 3,211 0	
	Waste Total	113,287	48,880	5,963	
Facility 0	00394				
F-Waste F003	Constituents Reported to TRI	: 19,800	0	0	
	Waste Total	19,800	0	0	
Facility 0	00397				
F-Waste ACET METH	Constituents Reported to TRI ONE ANOL	: 11,680 0	0 0	0 0	
	Constituent Total	11,680		0	
F-Waste: F001 F002 F003 F003	s Reported to NCAR: F003 F005	0 0 43,318 0	1,342 1,282 0 38,966	952 1,924 2,519 24,416	
	Waste Total	43,318	41,590	29,811	
Facility 0	00398				
F-Waste ACETC METH TOLUI XYLEI	Constituents Reported to TRI ONE YL ETHYL KETONE (MEK) ENE NE	: 3,479 10,707 4,296 1,500	4,122 9,309 0	0 6,989 0 0	
	Constituent Total	19,982	13,431	6,989	

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07/23/93 Page 39	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs Tra 1988	ansferred Off 1989	f-site 1990
F-Wastes Repo D001F003F0 D001F004F0 F003 F005	rted to NCAR: 05 05	0 0 66,67 <b>4</b> 151,653	142,590 10,638 0 0	74,822 0 0 0
	- Waste Total	218,327	153,228	74,822
Facility 000399				
F-Waste Const D001F001 D001F002	ituents Reported to TRI:	0 0	0 0	36,000 5,900
	Waste Total	0	0	41,900
Facility 000400				
F-Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL TOLUENE XYLENE	ituents Reported to TRI: COHOL YL KETONE (MEK)	0 0 12,642 9,453 0	0 0 0 0 0 0	0 0 0 0 0 0
	Constituent Total	22,095	0	0
F-Wastes Repo D001D035F0 F003F005 F005	rted to NCAR: 03F005	0 0 <b>4</b> 5,150	0 85,386 0	31,535 0 0
	Waste Total	45,150	85,386	31,535
Facility 000404				
F-Waste Const METHYL ETH TOLUENE XYLENE	ituents Reported to TRI: YL KETONE (MEK)	0 0 0	0 0 0	0 0 0
	Constituent Total	0	0	0
F-Wastes Repo F005	rted to NCAR:	539,512	750,000	816,160
	Waste Total	539,512	750,000	816,160
Facility 000405				
F-Waste Const ACETONE n-BUTYL AL ETHYLBENZE METHYL ETH METHYL ISO TOLUENE XYLENE	ituents Reported to TRI: COHOL NE YL KETONE (MEK) BUTYL KETONE	0 0 0 533 2,100	0 0 0 31,306 0	0 0 0 11,484 0
	Constituent Total	2,633	31,306	11,484
F-Wastes Repo F003	rted to NCAR:	37,250	27,199	28,945

07/23/93 Page 40	Comparison of F-Wastes R F-Waste Constituents Rep	eported to ported to TR	NCAR and NI: 1988 to 1	990
Constituent/	- Waste Code	Lbs Tra 1988	ansferred Off 1989	-site 1990
			17 EE7	12 114
FUUS				12,114
	Waste Total	37,250	44,756	41,059
Facility 000	406			
F-Waste C METHYL 1,1,1- XYLENE	onstituents Reported to TRI: ETHYL KETONE (MEK) TRICHLOROETHANE	0 0 0	37,000 0 0	0 0 0
	Constituent Total	0	37,000	0
F-Wastes D001D0 D001F0 F003F0 F005	Reported to NCAR: 02F003F005 05 05 Waste Total	0 0 38,530 38,530	0 2,503 37,132 0 	3,575 0 36,749 40,324
Facility 000	407	,	,	
F-Waste Co TRICHL	ONSTITUENTS Reported to TRI: OROETHYLENE	3,250	4,631	0
	 Constituent Total	3,250	4,631	0
F-Wastes F001 F002 F003	Reported to NCAR:	6,628 0 3,390	12,187 1,218 4,870	6,468 0 5,485
	 Waste Total	10,018	18,275	11,953
Facility 000	410			
F-Waste C METHAN 1,1,1-	onstituents Reported to TRI: OL TRICHLOROETHANE	0 0	0 0	0 0
	 Constituent Total	0		0
F-Wastes F004	Reported to NCAR:	6,000	20,000	5,450
	Waste Total	6,000	20,000	5,450
Facility 000	411			
F-Waste Co XYLENE	onstituents Reported to TRI:	0	0	0
	Constituent Total	0	0	0
F-Wastes F002	Reported to NCAR:	45	0	0
	 Waste Total	45		0
Facility 000	414			
F-Waste C	onstituents Reported to TRI:			
ACETON METHYL	E ETHYL KETONE (MEK)	0 0	0 0	0 0

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# Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

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	Lbs Ti	ansferred Of	f-site
Constituent/Waste Code	1988	1989	1990
METHANOL TOLUENE XYLENE	0 0 0	0 0 0	0 0 0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: D001F003F005 F003F005	0 0	0 232,127	341,013 0
Waste Total	0	232,127	341,013
Facility 000417			
F-Waste Constituents Reported to TRI	:		
ACETONE	368,488	301,975	0
n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK)	0 72,864	59.713	0
METHANOL	27,065	21,598	õ
METHYL ISOBUTYL KETONE	0	0	0
TOLUENE	218,595	179,138	0
XILENE			
Constituent Total	687,012	562,424	0
F-Wastes Reported to NCAR:			
D001F003F005		0	1,604,863
F003 F003F005	266,400 ∩	1.706.075	0
F005	1,775,850	0	Ō
		1 706 075	
Waste Total	2,042,250	1,/06,0/5	1,604,863
Facility 000426			
F-Waste Constituents Reported to TRI	:		
ACETONE	0	0	0
n-BUTYL ALCOHOL	0	0	0
METHYL ETHYL KETONE (MEK) METHANOL	0	0	0
METHYL ISOBUTYL KETONE	ŏ	õ	Õ
TOLUENE	0	0	0
XYLENE	0	0	0
Constituent Total	0	0	0
E-Wastas Perorted to NCAR.			
D001F003F005	0	0	291,540
F003F005	0	344,824	0
F005	460,224	0	0
Waste Total	460,224	344,824	291,540
Facility 000427			
F-Waste Constituents Reported to TRI METHANOL	: 0	0	0
Constituent Total	0	0	0
E-Waston Ponented to MORE			
F003	330	0	0
F005	435	0	0
Waste Total	765	0	0
nabee rotar		-	-

07/23/93Comparison of F-Wastes Reported to NCAR andPage42F-Waste Constituents Reported to TRI: 1988 to 1990				1990	
Constituent (Waste	-	Lbs Transferred Off-site			
waste	·	1900			
Facility 000428					
F-Waste Consti	tuents Reported to TRI:				
METHANOL TETRACHLORC	DETHYLENE	7,900 0	5,700 0	505 0	
	 Constituent Total	7,900	5,700	505	
F-Wastes Repor	ted to NCAR:				
D001F002		0	400	0	
D001F005 F001		6.710	40	0	
F002		0	1,788	1,000	
F003		0	0	400	
	Waste Total	6,710	2,228	1,400	
Facility 000432					
F-Waste Consti	tuents Reported to TRI:				
1,1,1-TRICH	ILOROETHANE	0	0	10,350	
	Constituent Total	0	0	10,350	
F-Wastes Repor	ted to NCAR:				
F001		17,020	18,516	10,783	
F002 F005		25,185	0 1,191	0 3,172	
	 Waste Total	42,205	19.707	13,955	
				- • - •	
Facility 000439					
F-Waste Consti	tuents Reported to TRI:				
n-BUTYL ALC	COHOL	0	0	0	
TOLUENE	HLORIDE	0	15.091	0	
XYLENE		õ	14,592	õ	
	 Constituent Total	0	29,683	0	
E-Wastes Benor	ted to NCAR.				
F003	teu to mean.	66,469	40,500	50,600	
F005		0	66,000	60,000	
	Waste Total	66,469	106,500	110,600	
Facility 000441					
F-Waste Consti	tuents Reported to TRI:				
ACETONE		9,900	5,900	1,900	
TOLUENE	·	650	1,800	9,200	
	Constituent Total	10,550	7,700	11,100	
F-Wastes Repor	ted to NCAR:				
F003 F003F005		41,250	990 36 450	0 36 420	
rvusrvus			JU,4JU	50,420	
	Waste Total	41,250	37,440	36,420	

### Facility 000442

F-Waste Constituents Reported to TRI:

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Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

Constituent/Waste Code	Lbs T 1988	ransferred C 1989	)ff-site 1990
ACETONE			
ACEIONE -			
Constituent Total	U	U	U
F-Wastes Reported to NCAR: F001	3,573	2,928	5,422
F003	11,414	5,105	6,287
Waste Total	14,987	8,033	11,709
Facility 000445			
F-Waste Constituents Reported to TRI:			
ACETONE CRESOL(S)	2,592	3,478	5,034
METHYL ETHYL KETONE (MEK)	25,807	1,105	1,091
METHYLENE CHLORIDE TOLUENE	0 36,834	0 24,433	286 31,683
- Constituent Total	65,233	29,016	38,094
F-Wastes Reported to NCAR:			
D001F005K086	0	0	189,489
Waste Total	0	0	189,489
Facility 000447			
F-Waste Constituents Reported to TRI: TOLUENE	: 19,800	19,145	13,805
Constituent Total	19,800	19,145	13,805
E-Wastes Reported to NCAR.			
F005	9,262	10,883	10,560
Waste Total	9,262	10,883	10,560
Facility 000451			
F-Waste Constituents Reported to TRI:	:		
ACETONE METHANOL	0	750 750	0
Constituent Total	0	1 500	0
	v	1,000	· · · ·
F-wastes Reported to NCAR: D001F001F002F003	0	0	29,296
F002	1,370	459	0
F003 F003F005	4,565	3,211 0	3,577
Waste Total	5,935	3,670	34,243
Facility 000452			
F-Waste Constituents Reported to TRI	:		
CYCLOHEXANE	0	0 5 410	0
TOLUENE	0	20,704	19,855
Constituent Total	0	26,114	 19,855
F-Wastes Reported to NCAR:			
D001F005	0	28,700	28,300

07/23/93Comparison of F-wastes Reported to NCAR andPage44F-Waste Constituents Reported to TRI: 1988 to 1990				990
Constituent/Wast	- e Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
F005		68,656	0	20,182
	 Waste Total	68,656	28,700	48,482
Facility 000456				
F-Waste Const 1,1,1-TRIC TRICHLOROE	ituents Reported to TRI: HLOROETHANE THYLENE	0 0	0 0	0 0
	 Constituent Total		0	0
F-Wastes Repo F001	rted to NCAR:	12,332	11,871	5,107
	 Waste Total	12,332	11,871	5,107
Facility 000457				
F-Waste Const TOLUENE	ituents Reported to TRI:	0	0	0
•	 Constituent Total	0	0	0
F-Wastes Repo D001D002F0 D001F002	rted to NCAR: 03	0 0	0 0	14,098 12,640
	 Waste Total	0	0	26,738
Facility 000463				
F-Waste Const D001F003 F002	ituents Reported to TRI:	0 0	0 0	2,950 2,327
	 Waste Total	0	0	5,277
Facility 000465				
F-Waste Const FREON 113 TOLUENE XYLENE	ituents Reported to TRI:	0 0 0	1,650 0 0	1,6 <b>40</b> 0 0
	 Constituent Total	0	1,650	1,640
F-Wastes Repo F001 F001F002 F001F003 F002 F002F003 F003	rted to NCAR:	7,103 0 1,365 0 11,565	1,084 0 2,328 12,932 816 50,068	0 1,640 2,328 1,720 816 22,162
	Waste Total	20,033	67,228	28,666
Facility 000470				
F-Waste Const D001D008F0 D001F003F0 F001	ituents Reported to TRI: 05 05	0 0 2,600	0 650 3,055	1,995 0 4,355
	Waste Total	2,600	3,705	6,350

Comparison of F-Wastes Reported to NCAR and

07/23/93

07/23/93 0 Page 45 1	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Waste	Code	Lbs Tr 1988	ansferred O 1989	ff-site 1990
Facility 000472				
F-Waste Constit	tuents Reported to TRI:	_		_
ACETONE		0	0	0
METHANOL		ŏ	Ő	Ő
TOLUENE		0	0	0
	Constituent Total	0	- 0	0
F-Wastes Report	ted to NCAR:			
D001D005F003	3F005	0	0	1,140
D001D007D000	8F003	. 0	3.700	1,000
1002	-			
	Waste Total	0	3,700	2,140
Facility 000475				
F-Waste Constit	tuents Reported to TRI:	1		
D001F001F00	3	0	0	5,974
D001F001F00. D004D005F00	3F005 1F006	0	0	4,800
D007D008D03	5F005	Ő	õ	880
D007D008F00	3F005	0	0	400
F002		0	2,949	U O
F003F005		ŏ	<b>4</b> 58	3,000
	Washa Mahal			
Facility 000479	waste local	U	5,40/	25,111
Facility 0004/9				
F-Waste Consti	tuents Reported to TRI:		C 000	0
FREON 113 METHYLENE CI	HLORIDE	2,700	6,000	0
1,1,1-TRICH	LOROETHANE	40,000	11,000	Õ
TRICHLOROET	HYLENE	29,000	38,000	0
	Constituent Total	71,700	55,000	0
E Wastes Banan	tod to NCNP.			
Fool	Led to NCAR:	2,720	5,510	0
F002	_	72,500	50,400	0
F002F003F00	5	0	2,050	0
	Waste Total	75,220	57,960	0
Facility 000481				
F-Waste Constit	tuents Reported to TRI:	:		
ACETONE		6,377	0	0
TOLUENE	L RETONE (MER)	1.307	0	0
1,1,1-TRICH	LOROETHANE	0	Ō	Ō
XYLENE		0	0	0
	Constituent Total	7,684	0	0
F-Wastes Report	ted to NCAR:			
D001F003		0	0	57,450
F001 F002F003F00	5	4,125	3-677	000
F003	-	10,805	0	0

07/23/93 Page 46	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs Tra 1988	ansferred Off 1989	-site 1990
F003F005 F005		0 3,575	50,411 0	32,105 0
	- Waste Total	18,505	54,088	90,455
Facility 000485				
F-Waste Const	ituents Reported to TRI:	627	٥	0
ACETUNE	COHOL	027	0	0
METHYL ETH	YL KETONE (MEK)	7.224	ŏ	ŏ
METHANOL		18,424	0	0
TOLUENE		9,482	0	0
XYLENE	-	0	0	0
	Constituent Total	35,757	0	0
F-Wastes Repo	rted to NCAR:		•	40 700
D001D035F0	03F005	0	E1 000	49,700
F005		65,800	51,800	0
•	- Waste Total	65,800	51,800	49,700
Facility 000486				
- E Wasta Canat	ituante Denerted to MDT.			
r-waste Const	ituents Reported to IRI:	٥	0	0
METHANOL		ő	õ	0 0
	Constituent Total	0	0	0
F-Wastes Repo	rted to NCAR:	<b>500</b>		400
F001		500	200	400
F002		8 400	0	Ő
F003F005		0,400	8.000	8.301
F005		35,000	0	0
	Waste Total	44,000	8,700	8,701
Facility 000491				
F-Waste Const	ituents Reported to TRI:			
METHYLENE	CHLORIDE	18,936	158,860	24,286
	Constituent Total	18,936	158,860	24,286
F-Wastes Repo	rted to NCAR:			
D001F002		0	0	51,400
D022F002F0	03F005	0	0	28,300
F002		154,200	54,500	0
F002F003	•	0	492,400	545,431
	Waste Total	154,200	546,900	625,131
Facility 000492				
F-Waste Const	ituents Reported to TRI:			
TETRACHLOR	OETHYLENE	38,388	0	0
	Constituent Total	38,388	0	0
F-Wastes Repo	orted to NCAR:			
D001F001		0	0	2,500

07/23/93 Page 47 Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Tra	ansferred Off	-site
Constituent/Waste Code	1988	1989	1990
F001	38,388	0	5,400
F001F002	0	0	16,700
F002	1,400	0	0
F005 -	3,490		
Waste Total	43,284	0	24,600
Facility 000493			
F-Waste Constituents Reported to TRI: XYLENE	12,840	9,230	11,165
Constituent Total	12,840	9,230	11,165
F-Wastes Reported to NCAR: F003	22,950	27,450	18,480
- Waste Total	22.950	27.450	18,480
Facility 000496	22,550		
F-Waste Constituents Reported to TKI: FREON 113	6.485	5,503	14,190
METHANOL	0	0	0
1,1,1-TRICHLOROETHANE	0	0	0
Constituent Total	6,485	5,503	14,190
F-Wastes Reported to NCAR:			
F001	7,500	6,109	14,191
- Waste Total	7,500	6,109	14,191
Facility 000497			
F-Waste Constituents Reported to TRI:			
ACETONE	2,186	6,068	5,389
METHYL ETHYL KETONE (MEK)	13	28	33 836
TOLUENE	0	1,102	71
-			
Constituent Total	2,575	7,262	6,329
F-Wastes Reported to NCAR:	11 900	23 348	0
F003F005	0	17,904	õ
F005	29,325	0	0
- Waste Total	41,225	41,252	0
Facility 000499			
F-Waste Constituents Reported to TRI:			
ACETONE	0	750	250
n-BUTYL ALCOHOL FTHYLBENZENE	0	750	250
METHYL ETHYL KETONE (MEK)	ŏ	750	250
METHANOL	0	750	250
METHYL ISOBUTYL KETONE	0	750	1 000
XYLENE	0	1,600	750
		7 400	2 000
Constituent Total	0	/,400	5,000

07/23 Page	/93 48	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Tr 1988	ansferred Of 1989	f-site 1990
F-1	Wastes Repo: F003 F005	rted to NCAR:	0 988,797	519,575 222,675	574,233 197,783
		Waste Total	988,797	742,250	772,016
Facil	ity 000506				
F-1	Waste Const METHYL ETH 1,1,1-TRIC	ituents Reported to TRI: YL KETONE (MEK) HLOROETHANE	3,609 0	3,926 0	2,756 0
		Constituent Total	3,609	3,926	2,756
F-'	Wastes Repo F001 F002 F003F005 F005	rted to NCAR: - Waste Total	0 13,870 0 24,000 37,870	0 22,500 31,500 0 54,000	4,122 0 23,000 0 27,122
Do at 14	000E07		ŗ	·	
Facil	1CY 000507				
F-1	Waste Const F002 F003	ituents Reported to TRI:	1,089 10,200	0 8,600	0 0
		Waste Total	11,289	8,600	0
Facil	ity 000508				
F-'	Waste Const F003 F005	ituents Reported to TRI:	0 0	0 0	0 0
		Waste Total	0	0	0
Facil	ity 000512				
F-'	Waste Const	ituents Reported to TRI:	:		
	ACETONE n-BUTYL AL	COHOL	65,700 10,800	0 0	0 0
	ETHYLBENZE		0	0	0
	METHYL ETH METHANOL	IL RETONE (MER)	92,100	0	0
	METHYL ISO	BUTYL KETONE	0	0	0
	TOLUENE XYLENE		53,200 34,800	0	0
		- Constituent Total	257,500	0	0
F-1	Wastes Repo F003F005 F005	rted to NCAR:	0 322,712	512,671 0	512,671 0
		- Waste Total	322,712	512,671	512,671
Facil	ity 000514				
F-1	Waste Const	ituents Reported to TRI:	:		
-	CRESOL(S)	····	2,600	4,200	17,350
	ETHYLBENZE	NE	0 2.100	1,500 3,500	5,205
		-	_,		

07/23/93 Page 49 Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

Constituent ///sta Cada	Lbs T:	ransferred O:	ff-site
Constituent/waste Code	1988	1909	1990
Constituent Total	4,700	9,200	39,510
F-Wastes Reported to NCAR: F002	1,260	1,215	630
Waste Total	1,260	1,215	630
Facility 000517			
F-Waste Constituents Reported to TRI:	:	19 980	0
1,1,1-TRICHLOROETHANE TRICHLOROETHYLENE	21,274 8,404	1,815 8,780	0
Constituent Total	29,678	30,575	0
F-Wastes Reported to NCAR:			
F001 F002	51,040 0	1,000	0
F003	0	5,200	0
Waste Total	51,040	31,400	0
Facility 000519			
F-Waste Constituents Reported to TRI:	: 	750	750
METHYLENE CHLORIDE	38,000	/ 50	0
METHANOL	12,000	750	0
TOLUENE	27,000	750	. 10
Constituent Total	102,900	4,600	765
F-Wastes Reported to NCAR:			
D001F005 F001F002F003F004	0	13,534	0
F003	265,135	20,000	ŏ
F005 F005P030P077P098	417,973 0	120	91,466 2,725
Waste Total	683,108	40,314	94,191
Facility 000522			
F-Waste Constituents Reported to TRI	:		
ACETONE	0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F003	41,145	71,000	83,350
Waste Total	41,145	71,000	83,350
Facility 000524			
F-Waste Constituents Reported to TRI	:		
ACETONE METHANOL	15,040 1,003,000	15,400 1,050,000	15,000 22,000
Constituent Total	1,018,040	1,065,400	37,000
F-Wastes Reported to NCAR: F005	0	5,632	4,327

07/23 Page	/93 50	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs T 1988	ransferred C 1989	off-site 1990
		Waste Total	0	5,632	4,327
Facil	ity 000529				
F-	Waste Const	tituents Reported to TRI:	0	•	1 400
	F001F003F0	105	· 0	0	83,100
	F003		85,014	0	0
	F003F004 F003F005		U O	0 106.256	64,520 0
	10031003	- Waste Total	85,014	106,256	149,020
Facil	ity 000532				·
- 7	Waste Const	ituents Reported to TRI:			
•	METHYL ETH	YL KETONE (MEK)	637	1,500	0
	METHANOL		855	1 500	0
	XYLENE		2,202	1,500	õ
		- Constituent Total	3,774	3,000	0
F-	Wastes Repo	orted to NCAR:			
	D001D001F0	005	0	23,513	0
	D001F001F0 D001F003F0	)05 )05	0	ں 35,488	15,137
	D001F005F0	006	Ō	0	2,752
	F003		22,193	0	0
		Waste Total	22,193	59,001	33,365
Facil	ity 000533				
F-	Waste Const METHANOL	tituents Reported to TRI:	0	750	5
		- Constituent Total	0	750	5
F-	Wastes Repo	orted to NCAR:			
	F003		0	420	0
		- Waste Total	0	420	0
Facil	ity 000538				
F-	Waste Const METHYLENE	ituents Reported to TRI: CHLORIDE	0	0	0
		Constituent Total	0	0	0
F-	Wastes Repo	orted to NCAR:			
	F002		0	1,200	0
	F003 F005		0	2,000	376
		- Waste Total	0	3,200	376
Facil	ity 000539				
F-	Waste Const	ituents Reported to TRI:			
•	METHYL ETH	IYL KETONE (MEK)	0	0	0
	TOLUENE	-	0	0	0

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07/2 Page	/23/93Comparison of F-Wastes Reported to NCAR andge 51F-Waste Constituents Reported to TRI: 1988 to 1990				1990
Cons	stituent/Wa	ste Code	Lbs Tra 1988	nsferred Off 1989	 1990
		Constituent Total	0	0	0
Ŧ	F-Wastes Re F003F005 F005	ported to NCAR:	0 4,800	2,752 0	5,504 0
		Waste Total	4,800	2,752	5,504
Faci	lity 00054	0			
F	F-Waste Con ACETONE n-BUTYL METHYL E	stituents Reported to TRI: ALCOHOL THYL KETONE (MEK)	: 0 0	0 0 0	0 0
	METHANOL TOLUENE		0	0	0
	XYLENE		0	0	0
		Constituent Total	0	0	0
F	F-Wastes Re F003F005 F005	ported to NCAR:	0 78,360	31,926 0	54,510 0
		Waste Total	78,360	31,926	54,510
Faci	ility 00054	1			
F	F-Waste Con ACETONE n-BUTYL METHYL E METHANOL TOLUENE XYLENE	stituents Reported to TRI ALCOHOL THYL KETONE (MEK)	0 0 0 0 0 0	0 0 0 0 0 0	0 0 0 0 0 0
		Constituent Total	0	0	0
I	F-Wastes Re F003F005 F005	ported to NCAR:	0 166,800	197,946 0	113,218
		Waste Total	166,800	197,946	113,218
Faci	ility 00054	2			
H	F-Waste Con ACETONE n-BUTYL METHYL E METHANOL METHYL I TOLUENE XYLENE	stituents Reported to TRI ALCOHOL THYL KETONE (MEK) SOBUTYL KETONE	: 0 0 0 0 0 0	0 0 0 0 0 0	. 0 0 0 0 0
		Constituent Total	0	0	0
F	F-Wastes Re F003F005 F005	ported to NCAR:	0 109,200	111,281	80,431 0
		Waste Total	109,200	111,281	80,431

07/23/93 Page 52	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Waste	- Code	Lbs Tra 1988	ansferred Off 1989	E-site 1990
Facility 000543				
F-Waste Constit	tuents Reported to TRI:	0	0	n.
METHYL ETHY	L KETONE (MEK)	0	õ	0
METHANOL		0	0	0
TOLUENE		0	U O	0
	Constituent Total	0	0	0
F-Wastes Repor	ted to NCAR:	•		226 740
F003F005 F005		0 442.168	306,537	226,748
1005				
	Waste Total	442,108	306,537	220,/48
Facility 000544				
F-Waste Consti	tuents Reported to TRI:	•		<u> </u>
ACETONE	OHOT.	0	0	0
METHYL ETHY	L KETONE (MEK)	0	0	0
METHANOL		0	0	0
XYLENE		0 0	0	0
	 Constituent Total	0	0	0
F-Wastes Repor	ted to NCAR:			
F003F005		0	43,890	70,882
F005		54,000		0
	Waste Total	54,000	43,890	70,882
Facility 000545				
F-Waste Consti	tuents Reported to TRI:			
FREON 113	HLORIDE	6,600 12 893	0 11 550	0
	Constituent Total	19,493	11,550	0
F-Wastes Repor	ted to NCAR:		2 600	
F002 F002F003F00	5	26,400	3,600	3,600
F003	5	32,400	31,800	0
	 Waste Total	58,800	64,200	38,800
Facility 000547				
T. Masha Carati	turnt - Devented to MDT.			
METHYL ETHY	L KETONE (MEK)	350,000	451,760	200,000
	Constituent Total	350,000	451,760	200,000
F-Wastes Repor	ted to NCAR:			
F003		0	451,760	0
F003F005 F005		0 452 360	0 976	303,050 A
	Waste Total	452,360	452,696	303,050

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07/23/93 Page 53	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
Facility 000548				
F-Waste Const n-BUTYL AL METHANOL TETRACHLOR XYLENE	ituents Reported to TRI: COHOL OETHYLENE	0 0 1,748 0	0 0 0 0	0 0 0 0
	Constituent Total	1,748	0	0
F-Wastes Repc D001F002F0 F002 F002F003 F002F003F0	orted to NCAR: 03F005 05	0 38,750 0 0	0 4,400 0	2,000 600 4,400 36,660
	waste lotal	38,750	4,400	43,000
Facility 000550 F-Waste Const ACETONE METHYL ETH METHANOL TOLUENE	ituents Reported to TRI: YL KETONE (MEK)	2,520 1,196 0 2,520	6,189 627 157 0	0 0 0 0
	Constituent Total	6,236	6,973	0
F-Wastes Repc F003 F005	rted to NCAR: - Waste Total	0 10,312 10,312	8,565 0 8,565	14,025 0 14,025
Facility 000552				
- F-Waste Const ACETONE METHANOL	ituents Reported to TRI: - Constituent Total	0 1,301 1,301	9,213 233 9,446	0 0 
F-Wastes Repo F005	rted to NCAR:	7,923	0	0
	- Waste Total	7,923	0	0
Facility 000553				
F-Waste Const TETRACHLOR 1,1,1-TRIC	ituents Reported to TRI: OETHYLENE HLOROETHANE	0 0	0 0	0 0
	Constituent Total	0	0	0
F-Wastes Repo F001 F003 F005	orted to NCAR:	18,800 3,100 0	16,050 0 200	20, <b>4</b> 75 0 0
	Waste Total	21,900	16,250	20,475

#### Facility 000557

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F-Waste Constituents Reported to TRI:

07/23 Page	/93 5 <b>4</b>	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Tr 1988	ansferred Of 1989	ff-site 1990
	ACETONE METHYL ETH METHANOL TOLUENE	YL KETONE (MEK)	0 0 1,450 4,000	0 414 126 754	0 1,390 0 6 <b>4</b> 0
		Constituent Total	5,450	1,294	2,030
F-'	Wastes Repo F003	rted to NCAR:	21,113	0	9,632
		Waste Total	21,113	0	9,632
Facil	ity 000563				
F-	Waste Const n-BUTYL AL METHYL ETH	ituents Reported to TRI COHOL YL KETONE (MEK)	: 8,595	0 23,100	0 16,640
		Constituent Total	8,595	23,100	16,640
F-`	Wastes Repo D007F005 F005	rted to NCAR:	0 49,764	54,000 62,400	42,657 30,972
		Waste Total	49,764	116,400	73,629
Facil	ity 000568				
F-	Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL TOLUENE XYLENE	ituents Reported to TRI COHOL YL KETONE (MEK)	: 21,934 0 6,580 6,142 6,580 0	2,800 0 0 0 0 0	0 0 0 0 0
		Constituent Total	41,236	2,800	0
F-	Wastes Repo F001 F003 F005	rted to NCAR:	0 0 82,350	0 21,150 0	3,875 0 4,998
		Waste Total	82,350	21,150	8,873
Facil	ity 000574				
F-	Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL METHYL ISO TOLUENE XYLENE	ituents Reported to TRI COHOL YL KETONE (MEK) BUTYL KETONE	: 0 0 0 0 0 0 0	• 0 0 0 0 0 0 0	0 0 0 0 0 0 0
		Constituent Total	0	0	0
F-	Wastes Repo F003 F005	rted to NCAR:	0 0	0 0	0 0
		Waste Total	0	0	0

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07/23/93 Comparison of F-Was Page 55 F-Waste Constituent	'93 Comparison of F-Wastes Reported to NCAR and 55 F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Waste Code	Lbs 7 1988	Transferred C 1989	off-site 1990	
Facility 000575				
F-Waste Constituents Reported to	TRI:			
ACETONE	0	0	0	
n-BUTYL ALCOHOL	0	0	0	
METHYL ETHYL KETONE (MEK)	0	0	0	
METHANOL	. 0	0	0	
TETRACHLOROETHYLENE	0	0	0	
	0	0	0	
I, I, I-TRICHLOROEINANE VVLENE	ŏ	Ő	Ő	
Constituent Moto				
		Ū	v	
F-Wastes Reported to NCAR:	20 912	21 458	20 636	
	20,012	1,308	4,814	
F003F005	õ	1,500	7,522	
10031003				
Waste Tota	al 20,812	22,766	32,972	
Facility 000576				
F-Waste Constituents Reported to	TRI:			
1,1,1-TRICHLOROETHANE	0	0	32,984	
Constituent Tota	a1 0	0	32,984	
F-Wastes Reported to NCAR:				
D001F003	0	5,280	1,765	
F001	26,548	37,408	32,984	
F002	0	0	481	
Waste Tota	al 26,548	42,688	35,230	
Facility 000579				
F-Waste Constituents Reported to	TRI:			
n-BUTYL ALCOHOL	0	0	0	
METHYL ISOBUTYL KETONE	0	0	0	
TOLUENE	0	0	0	
1, 1, 1-TRICHLOROETHANE	0	0	0	
XYLENE		U 		
Constituent Tota	al 0	0	0	
F-Wastes Reported to NCAR:				
F001	0	0	1,000	
F002	17,456	12,316	10,392	
F003	3,194	24,073	41,416	
F005	/,634		10,972	
Waste Tota	al 28,284	36,389	63,780	
Facility 000580				
F-Waste Constituents Reported to	TRI:			
F003	0	0	1,200	
Waste Tota	al 0	0	1,200	
Facility 000583				
- P-Wasta Constituents Banarted to	TTR T •			
METHYL ETHYL KETONE (MEK)	0	0	16,397	

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07/23 Page	/93 56	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Tra 1988	nsferred Of 1989	f-site 1990
		Constituent Total	0	0	16,397
F-	Wastes Repo D001F003 F003	rted to NCAR:	0 37,440	13, <b>4</b> 55 26,197	19,724 16,221
		Waste Total	37,440	39,652	35,945
Facil	ity 000586				
F-	Waste Const METHYL ETH	ituents Reported to TRI: YL KETONE (MEK)	5,227	3,378	0
		Constituent Total	5,227	3,378	0
F-	Wastes Repo	rted to NCAR:			
	F003F005 F005	-	0 109,416	42,000 46,328	39,093 77,062
		Waste Total	109,416	88,328	116,155
Facil	ity 000590				
F-	Waste Const	ituents Reported to TRI:			
-	FREON 113		0	1,424	8,059
	METHYLENE	CHLORIDE OFTHVI ENE	0	1,424	0
	1,1,1-TRIC	HLOROETHANE	õ	0 0	7,591
		Constituent Total	0	2,848	15,650
F-	Wastes Repo F001	rted to NCAR:	3,350	5,798	8,059
		Waste Total	3,350	5,798	8,059
Facil	lity 000599				
F-	Waste Const	ituents Reported to TRI:	•		
	TOLUENE	HLOROFTHANE	0	0	0
	XYLENE		2,964	Ő	3,000
		Constituent Total	2,964	0	3,000
F-	Wastes Repo	rted to NCAR:	•	10 000	1 401
	F001 F003		0	10,200	1,401
	F005		õ	õ	6,359
		Waste Total	0	10,200	17,785
Facil	lity 000601				
F-	Waste Const	ituents Reported to TRI:		-	-
n-BUTYL AL	n-BUTYL AL METHYL ETH	COHOL YL KETONE (MEK)	2,252 0	0 0	0
		Constituent Total	2,252	0	0
F.	Wastes Pero	orted to NCAR:			
L.	F003 F005		4,120 4,565	3,887 10,229	3,215 12,400
		•			

07/23/93 Page 57	Comparison of F-Wastes F-Waste Constituents Re	omparison of F-Wastes Reported to NCAR and Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs Ti 1988	ransferred O 1989	ff-site 1990	
	Waste Total	8,685	14,116	15,615	
Facility 000603					
F-Waste Const ACETONE	ituents Reported to TRI:	0	0	0	
	- Constituent Total	0	0	0	
F-Wastes Repo F003	rted to NCAR:	0	0	0	
	- Waste Total	0	0	0	
Facility 000604					
F-Waste Const 1,1,1-TRIC	ituents Reported to TRI: HLOROETHANE	74,000	0	0	
	Constituent Total	74,000	0	0	
F-Wastes Repo F001	rted to NCAR:	70,470	58,971	57,924	
	- Waste Total	70,470	58,971	57,924	
Facility 000608					
F-Waste Const METHYL ETH TOLUENE	ituents Reported to TRI: YL KETONE (MEK)	0 0	0 1,500	0 500	
	- Constituent Total	0	1,500	500	
F-Wastes Repc D001D007F0 D007D008F0 F003F005 F003F006 F005	rted to NCAR: 03F005 03F005 - Waste Total	0 0 118,387 	0 0 184,144 	12,750 67,020 0 0 	
Facility 000610					
F-Waste Const ACETONE METHYL ETH METHANOL TOLUENE	ituents Reported to TRI: YL KETONE (MEK)	5,900 15,700 0 14,600	11,300 46,000 2,700 47,000	10,500 49,505 1,750 25,803	
	constituent iotai	50,200	107,000	87,558	
F-Wastes Repo D001F003F0 F003 F003F005 F005	orted to NCAR: 05 	0 294,206 0 0	0 185,321 0 266,789	300,030 0 39,410 0	
Facility 000611	maste IUtai	274,200	<del>4</del> 72,11V	<i>JJJ</i> , <del>1</del> 10	
F-Waste Const METHYL ETH	ituents Reported to TRI: YL KETONE (MEK)	7,320	10,800	0	

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07/23 Page	/93 58	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Const	ituent/Wast	e Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
	1 1 1-TRTC	THLOROETHANE	76.840	12.040	0	
	1,1,1 1111	Constituent Total	84,160	22,840	0	
F-	Wastes Repo F001 F002	orted to NCAR:	71,170	560 11,480	905 8,520	
	F005	Waste Total	6,000 	3,045	9,425	
Facil	ity 000613		·		·	
F-	Waste Const ACETONE	tituents Reported to TRI	C: 0	0	0	
	n-BUTYL AI METHYL ETH METHANOL TOLUENE	LCOHOL HYL KETONE (MEK)	0 3,909 0 0	0 0 0	0 0 0 0	
	AIDENE	Constituent Total			<u>-</u> 0	
F-	Wastes Repo D001D035F( F003F005 F005	orted to NCAR: 003F005 Waste Total	0 0 24,430 	0 41,650 0 	33,600 456 0 34,056	
Facil	ity 000614					
F-	Waste Const ACETONE n-BUTYL AI METHYL ETH METHANOL TOLUENE XYLENE	tituents Reported to TR LCOHOL HYL KETONE (MEK)	L: 9,039 0 2,280 2,382	0 8,304 3,061 1,500 4,176 3,000	0 4,129 9,539 10 3,524 0	
		Constituent Total	13,701	20,041	17,202	
F-	Wastes Repo F005	orted to NCAR:	77,859	80,109	100,800	
		Waste Total	77,859	80,109	100,800	
Facil	ity 000615					
F-	Waste Const ACETONE n-BUTYL AI METHYL ETH METHANOL TOLUENE XYLENE	tituents Reported to TR LCOHOL HYL KETONE (MEK)	I: 2,168 1,024 0 2,130 0	750 1,500 954 3,999 2,250	0 255 0 255 255 255	
		Constituent Total	5,322	9,453	765	
F-	Wastes Repo D001D035F( F003F005 F005	orted to NCAR: 003F005	0 0 <b>44</b> ,225	0 21,025 0	21,886 0 0	
		Waste Total	44,225	21,025	21,886	

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07/23/93 Page 59	Comparison of F-Wastes F-Waste Constituents R	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wa	aste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990	
Facility 00061					
F-Waste Cor 1,1,1-TF TRICHLOF	nstituents Reported to TRI RICHLOROETHANE ROETHYLENE	:: 16,200	0 0	1,005 7,704	
	Constituent Total	16,200	0	8,709	
F-Wastes Re F001 F002 F006	eported to NCAR:	0 16,200 100,055	9,575 0	7,704 0 0	
	Waste Total	116,255	9,575	7,704	
Facility 00062	22				
F-Waste Cor n-BUTYL METHYL I	ALCOHOL ALCOHOL THYL KETONE (MEK)	0	0 0	0 0	
•	Constituent Total	0	0	0	
F-Wastes Reporte D001F005 F005	eported to NCAR:	0 14,640	<b>44</b> ,202 19,230	0 22,620	
	Waste Total	14,640	63,432	22,620	
Facility 00062	23				
F-Waste Con n-BUTYL TOLUENE	nstituents Reported to TRI ALCOHOL	[: 0 0	0 0	0 0	
	Constituent Total	0	0	0	
F-Wastes R F003 F003F00!	eported to NCAR: 5	0 0	800 0	0 12,800	
	Waste Total		800	12,800	
Facility 00062	24				
F-Waste Con ACETONE METHYL 1 METHYL 1 TOLUENE XYLENE	nstituents Reported to TR ETHYL KETONE (MEK) ISOBUTYL KETONE	I: 2,323 7,251 2,323 20,847 3,485	4,350 13,131 6,526 41,658 7,005	0 0 0 0	
	Constituent Total	36,229	72,670	0	
F-Wastes R D001F00 F001 F005	eported to NCAR: 3F005	0 0 48,255	0 4,000 94,336	173,394 0 0	
	Waste Total	48,255	98,336	173,394	

## Facility 000626

F-Waste Constituents Reported to TRI:

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07/23 Page	/93 60	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Tra 1988	ansferred Off 1989	E-site 1990
	FREON 113 1,1,1-TRIC	HLOROETHANE	0 0	0 0	0 0
		 Constituent Total	0	0	0
F-	Wastes Repo F001 F002 F003 F003F005 F005	rted to NCAR:	0 11,912 0 198	10,092 41,283 0 459 459	1,376 11,492 229 0 0
		Waste Total	12,110	52,293	13,097
Facil F-	ity 000631 Waste Const ACETONE	ituents Reported to TRI:	0	0	0
F-	Wastes Repo	constituent Total	0	U	0
•	F003		144,111	54,087	111,500
Es ei 1	iter 000622	Waste Total	144,111	54,087	111,500
Facil	Waste Const ACETONE	ituents Reported to TRI:	7,155	0	0
		Constituent Total	7,155	0	0
F-	Wastes Repo F003	rted to NCAR:	35,285	16,246	0
		Waste Total	35,285	16,246	0
Facil F-	ity 000634 Waste Const ACETONE n-BUTYL AL METHYL ETH METHANOL METHYL ISO TOLUENE XYLENE	ituents Reported to TRI: COHOL YL KETONE (MEK) BUTYL KETONE	1,340 1,673 1,231 3,362 8,830 11,060 3,360	3,474 3,744 3,914 9,230 29,100 29,331 6,537	1,489 2,395 1,786 3,468 18,655 19,206 4,697
		Constituent Total	30,856	85,330	51,696
F-	Wastes Repo F001 F003 F003F005 F005	rted to NCAR:	750 3,150 0 139,727	0 167,482 359,065 0	0 43,196 162,266 0
		Waste Total	143,627	526,547	205,462
Facil	ity 000637	ituante Dosautad to MPT			
F	TRICHLOROE	THYLENE	0	0	0
		Constituent Total	0	0	0

Same A

07/23/93Comparison of F-Wastes Reported to NCAR andPage 61F-Waste Constituents Reported to TRI: 1988 to 1990				1990
Constituent/Waste	- Code	Lbs Transferred Off-site 1988 1989 1990		
F-Wastes Report F001	ted to NCAR:	4,123	4,128	9,174
	Waste Total	4,123	4,128	9,174
Facility 000638				
F-Waste Constit	tuents Reported to TRI:			
METHYLENE CI 1,1,1-TRICHI	HLORIDE LOROETHANE	20,300	0 0	0 0
	Constituent Total	20,300	0	0
F-Wastes Report	ted to NCAR:	10 154	22.252	15 200
F001 F002		42,154 340	29,850	17,320
F002F004 F003F005		0	312 9,080	1,176 6,479
	 Waste Total	42,494	39,242	24,975
Facility 000639				
F-Waste Constit	tuents Reported to TRI:			
METHYLENE CI TOLUENE	HLORIDE	0 6,140	0 10,698	0 10,600
1,1,1-TRICH	LOROETHANE	0	0	30
	Constituent Total	6,140	10,698	10,630
F-Wastes Report	ted to NCAR:	•		
D001D002F003 D001D006D003	1 8F002	0	3,300 22,694	0
D001F002F00	5 8 F 0 0 1	0	0	25,540 540
F001	8F001	19,250	19,449	9,686
F005		19,113	0	0
	Waste Total	38,363	45,443	35,766
Facility 000640				
F-Waste Constit	tuents Reported to TRI:	٥	100	600
BENZENE		Ő	0	0
METHANOL TOLUENE		0 0	0 100	0 660
	 Constituent Total	0	200	1,260
F-Wastes Report	ted to NCAR:			
D001F005		0	0 1,600	138,900 6,600
F005		0	22,400	0
	Waste Total	0	24,000	145,500
Facility 000643				
F-Waste Constit	tuents Reported to TRI:	10 429	7 627	E E01
TETRACHLORO	ETHYLENE	4,980	7,515	4,141
1,1,1-TRICH	LOROETHANE	7,415	10,035	10,199

Comparison of F-Wastes Reported to NCAR and

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07/23 Page	/93 62	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
		Constituent Total	22,824	25,187	21,034
F-1	Wastes Repo: F001 F002 F003 F005	rted to NCAR: - Waste Total	1,600 19,790 2,790 0 	11,950 10,496 9,803 1,600 	18,334 3,640 6,746 1,505 30,225
Facil	ity 000644				
F-	Waste Const METHYL ETH XYLENE	ituents Reported to TRI: YL KETONE (MEK)	8,000 6,520	0	4,620
_		Constituent Total	14,520	U	4,020
F'-	Wastes Repo F003F005 F005	rted to NCAR:	0 46,569	<b>49,7</b> 00 0	92,400 0
•		Waste Total	46,569	49,700	92,400
Facil	ity 000647				
F-	Waste Const n-BUTYL AL METHYL ETH METHANOL TOLUENE XYLENE	ituents Reported to TRI: COHOL YL KETONE (MEK)	0 0 0 0	0 0 0 0	0 0 0 0
_			0	0	U
F-	Wastes Repo F003 F003F005	rted to NCAR:	68,923 0	101,469 3,691	79,826 27,619
		Waste Total	68,923	105,160	107,445
Facil	ity 000648				
F-	Waste Const n-BUTYL AL METHANOL TOLUENE	ituents Reported to TRI: COHOL - Constituent Total	6,479 0 2,905 9,384	2,860 7,328 3,604 13,792	7,745 1,949 6,296 15,990
F-	Wastes Repo	rted to NCAR:			
	F003F005 F005		0 54,100	61,950 0	72,450 0
		Waste Total	54,100	61,950	72,450
Facil	ity 000649				
F-	Waste Const METHYLENE METHANOL	ituents Reported to TRI: CHLORIDE	5,454 1,432	15,430 334	2,741 188
		Constituent Total	6,886	15,764	2,929

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07/23/93 Page 63 Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Tr	ansferred Of	f-site
Constituent/Waste Code	1988 	1989	1990
F-Wastes Reported to NCAR: D001F002F003F005 F002	0 5,464	0 15,430	456 2,741
Waste Total	5,464	15,430	3,197
Facility 000650			
F-Waste Constituents Reported to TRI: ACETONE TOLUENE	962 1,846	2,064 2,540	3,984 3,710
Constituent Total	2,808	4,604	7,694
F-Wastes Reported to NCAR: F003	0	0	6,497
- Waste Total	0	0	6,497
Facility 000651			
F-Waste Constituents Reported to TRI: METHYLENE CHLORIDE TOLUENE	2,657 16,920	0 3,750	0 0
- Constituent Total	19,577	3,750	0
F-Wastes Reported to NCAR: D001F003 D001F003F005 D001F005 F001 F002 F002F003 F003F005 F005	0 0 12,750 8,855 0 0 12,825	733 367 367 0 0 550 0 1,467 64,065	0 0 2,200 0 367 2,434 12,031
Waste Total	34,430	67,549	17,032
Facility 000660 F-Waste Constituents Reported to TRI: ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE 1,1,1-TRICHLOROETHANE XYLENE	6,949 0 0 34,423 7,805 0	1,478 14,473 0 70 22,828 4,556 2,963	0 0 0 0 0 0 0
Constituent Total	49,177	46,368	0
F-Wastes Reported to NCAR: D001D035F003F005 D001F003F005 F001 F003 F003F005 F005	0 0 0 0 622,882	0 235,802 39,496 42,467 16,300 445,593 0	285,704 167,752 0 89,698 0 287,732 0
Waste Total	622,882	779,658	830,886

07/23/93 Page 64	omparison of F-Wastes Reported to NCAR and -Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs Tra 1988	nsferred Of: 1989	f-site 1990
Facility 000668				
F-Waste Const 1,1,1-TRIC	ituents Reported to TRI: HLOROETHANE	4,025	0	0
	- Constituent Total	4,025	0	0
F-Wastes Repo F001 F005	rted to NCAR:	11,297	8,360	0
roos	- Waste Total	11,297	9,491	0
Facility 000670				
F-Waste Const METHYLENE 1,1,1-TRIC	ituents Reported to TRI: CHLORIDE CHLOROETHANE	62,370 48,510	2,250 0	250 250
	- Constituent Total	110,880	2,250	500
F-Wastes Repo 'F002 F005	orted to NCAR:	<b>4</b> 1,580 0	44,000 8,865	0 0
	- Waste Total	41,580	52,865	0
Facility 000672				
F-Waste Const ACETONE n-BUTYL AL METHANOL TOLUENE	ituents Reported to TRI:	0 0 30,000 6,391	50,906 0 42,549 100	0 0 0
	Constituent Total	36,391	93,555	0
F-Wastes Repo F002 F003 F003F005 F005	orted to NCAR: - Waste Total	6,050 0 55,250 61,300	0 62,871 0 0 62,871	0 0 35,762 0 35,762
Facility 000674				
F-Waste Const XYLENE	ituents Reported to TRI:	36,135	19,609	14,906
	- Constituent Total	36,135	19,609	14,906
F-Wastes Repo D001F003 F001 F003 F005	orted to NCAR:	0 1,200 9,600 9,200	0 0 19,358 0	14,906 0 0
	- Waste Total	20,000	19,358	14,906
Facility 000675				
F-Waste Const ACETONE FREON 113	ituents Reported to TRI:	5,404 0	0 7,057	0
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# Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Tr	ansferred Of	f-site
Constituent/Waste Code	1988		
METHYLENE CHLORIDE 1,1,1-TRICHLOROETHANE	0 0	2,610 4,689	0 0
Constituent Total	5,404	14,356	0
F-Wastes Reported to NCAR: D001F003F005 F001 F002 F003 F005	0 0 0 27,720	33,522 0 14,904 1,140 0	0 0 0 0
Waste Total	27,720	49,566	0
Facility 000682			
F-Waste Constituents Reported to TRI TOLUENE 1,1,1-TRICHLOROETHANE	: 0 37,476	0 48,852	0 0
Constituent Total	37,476	48,852	0
F-Wastes Reported to NCAR: D001F003 F001 F002 F003	0 153,777 0 0	0 127,460 31,801 3,300	1,376 63,301 0 0
Waste Total	153,777	162,561	64,677
Facility 000683			
F-Waste Constituents Reported to TRI 1,1,1-TRICHLOROETHANE	: 0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F001	0	1,560	3,120
Waste Total	0	1,560	3,120
Facility 000684			
F-Waste Constituents Reported to TRI 1,1,1-TRICHLOROETHANE	: 0	0	0
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F001	12,320	0	0
Waste Total	12,320	0	0
Facility 000686			
F-Waste Constituents Reported to TRI 1,1,1-TRICHLOROETHANE	: 0	2,400	21,700
Constituent Total	0	2,400	21,700
F-Wastes Reported to NCAR: F001	23,791	20,740	18,341
Waste Total	23,791	20,740	18,341

07/23/93Comparison of F-Wastes Reported to NCAR andPage 66F-Waste Constituents Reported to TRI: 1988 to 1990				1990
Constituent/Waste	e Code	Lbs T 1988	ransferred O 1989	ff-site 1990
Facility 000694				
F-Waste Consti METHANOL	tuents Reported to TRI:	0	0	0
	- Constituent Total			
		v	Ŭ	Ŭ
F-wastes Repor D001F003F00 F003	15 15	0 126,778	0	122,036
F003F002	- Waste Total	126,778	142,793	122,036
Facility 000697				
F-Waste Consti	tuents Reported to TRI:			
METHANOL TRICHLOROET	'HYLENE -	0	0 0	0 0
•	Constituent Total	0	0	0
F-Wastes Repor	ted to NCAR:	25 100	22 529	31 657
F001 F003F005 F005		0 1,700	800 0	0 0
	- Waste Total	36,800	33,338	34,657
Facility 000700				
F-Waste Consti ETHYLENE GL FREON 113 TETRACHLORC 1,1,1-TRICH	tuents Reported to TRI: YCOL MONOETHYL ETHER DETHYLENE ILOROETHANE	8,900 14,000 81,000 100	0 24,000 29,800 0	0 0 0 0
	- Constituent Total	104,000	53,800	0
F-Wastes Repor F001 F002 F003 F005	ted to NCAR:	84,646 98,374 0 0	160 64,501 1,860 3,720	1,120 88,224 1,260 60
	Waste Total	183,020	70,241	90,664
Facility 000704				
F-Waste Consti METHANOL TOLUENE	tuents Reported to TRI:	106,280 868,923	0 807,454	0 305,100
	- Constituent Total	975,203	807,454	305,100
F-Wastes Repor	ted to NCAR:	·	·	
F005	-	891,760	833,880	385,430
	Waste Total	891,760	833,880	385,430
Facility 000707				
F-Waste Consti METHYL ETHY	tuents Reported to TRI: 'L KETONE (MEK)	4,972	0	0

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07/23/93 Page 67	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
	- Constituent Total	4,972		0
F-Wastes Repo F003F005 F005	orted to NCAR:	0 27,475	17,960 0	0 0
	Waste Total	27,475	17,960	0
Facility 000710				
F-Waste Const	ituents Reported to TRI:			
n-BUTYL AL	COHOL	0	0	0
METHYL ETH	IYL KETONE (MEK)	0	0	0
METHANOL		0	0	0
TOLUENE		0	0	0
	- Constituent Total	0	0	0
F-Wastes Repo	orted to NCAR:	917	0	0
1005	-			
•	Waste Total	917	0	0
Facility 000712				
F-Waste Const 1,1,1-TRIC	ituents Reported to TRI: CHLOROETHANE	32,000	44,000	0
	- Constituent Total	32,000	44,000	0
F-Wastes Repo F001	orted to NCAR:	45,628	52,137	63,422
	- Waste Total	45,628	52,137	63,422
Facility 000713				
F-Waste Const	ituents Reported to TRI:			
ACETONE METHYLENE	CHLORIDE	0 0	0 0	0 0
	- Constituent Total		0	0
E Wester Bong	anted to NCAR.			
F-Wastes Repo F002		654	0	0
	Waste Total	654	0	0
Facility 000717				
F-Waste Const METHANOL	tituents Reported to TRI:	0	0	0
	- Constituent Total	0	0	0
F-wastes Repo F003F005 F005	DITED TO NCAK:	0 12,705	21,945 0	20,691 0
	Waste Total	12,705	21,945	20,691

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### Facility 000721

F-Waste Constituents Reported to TRI:

07/23/93 Comparison of F-Wast Page 68 F-Waste Constituents	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Waste Code	Lbs Tr 1988	ansferred Off 1989	-site 1990	
ACETONE METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	0 0 0 0 0	0 0 0 0 0	0 0 0 0 0	
Constituent Total	0	0	0	
F-Wastes Reported to NCAR: F003F005 F005	0 20, <b>4</b> 05	21,9 <b>4</b> 5 0	21,054 0	
Waste Total	20,405	21,945	21,054	
Facility 000722				
F-Waste Constituents Reported to 1 ACETONE TOLUENE	TRI: 12,570 12,570	6,589 6,589	0 0	
Constituent Total	25,140	13,178	0	
F-Wastes Reported to NCAR: F002 F003	0 47,083	0 26,985	495 12,052	
Waste Total	47,083	26,985	12,547	
Facility 000725				
F-Waste Constituents Reported to 7 CRESOL(S) METHYLENE CHLORIDE TETRACHLOROETHYLENE 1,1,1-TRICHLOROETHANE	TRI: 0 0 0 0	0 0 0 0	5 5 5 5	
Constituent Total	L 0	0	20	
F-Wastes Reported to NCAR: F001	186,000	0	0	
Waste Total	186,000	0	0	
Facility 000730				
F-Waste Constituents Reported to 7 ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHYLENE CHLORIDE METHANOL TOLUENE TRICHLOROETHYLENE	TRI: 0 400 80,000 22,500 400 40,000 2,000	0 6,482 30,197 3,412 5,413 23,210 0	1,071 7,489 70,786 0 16,161 175,259 0	
Constituent Total	L 145,300	68,714	270,766	
F-Wastes Reported to NCAR: D001D005F001F003 D001D010F003F005 D001F001F003F005 D006F001F003F005 F001 F001F003F005 F005	0 0 0 44,170 0 158,983	0 0 0 10,850 0 189.660	305 3,100 25,275 0 181,815 0 97,820 0	

07/23/93 Cor Page 69 F-V	mparison of F-Wastes F Waste Constituents Rep	eported to ported to T	NCAR and RI: 1988 to 1	1990
Constituent/Waste Co	- ode	Lbs Tra 1988	ansferred Of: 1989	f-site 1990
	Waste Total	203,153	200,510	308,315
Facility 000733				
F-Waste Constitue	ents Reported to TRI:			
ACETONE		0	0	0
METHYL ETHYL I	KETONE (MEK)	0	0	0
TOLIENE		0	0	0
XYLENE		0	õ	Ō
	 Constituent Total	0	0	0
F-Wastes Reported	d to NCAR:			
F003F005		0	54,800	0
F005		24,000	0	36,000
	Waste Total	24,000	54,800	36,000
Facility 000734				
F-Wasté Constitue	ents Reported to TRI:	-		
FREON 113 METHYLENE CHLO	ORIDE	0 5,600	8,422 8,423	9,652 9,652
	Constituent Total	5,600	16,845	19,304
F-Wastes Reported	d to NCAR:			
F001		10,250	16,691	7,912
F002		1,600	0	2,140
F003		1,600	0	0
	Waste Total	13,450	16,691	10,052
Facility 000735				
F-Waste Constitue	ents Reported to TRI:			
METHYL ETHYL	KETONE (MEK)	9,203	5,518	12,800
TOLUENE	I. ENIE	14,/8/	3,652	9,020
INICHBOROEINI				
	Constituent Total	23,990	9,170	21,820
F-Wastes Reported	d to NCAR:	^	C 700	•
D001D007D008F	005	0	6,/20 N	U 7.350
F001		11,205	8,600	5,746
F002		600	300	1,768
F003F005		0	67,400	57,463
F005		72,100	U 	U 
	Waste Total	83,905	83,020	72,327
Facility 000737				
F-Waste Constitue TETRACHLOROET	ents Reported to TRI: HYLENE	750	750	0
	 Constituent Total	750	750	0
F-Wastes Reported	d to NCAR:			
F001	a oo morain	800	0	1,485
F003		0	3,200	4,268
F005		2,100	2,400	0

07/23/93 Page 70	Comparison of F-Wastes F-Waste Constituents Re	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent	/Waste Code	Lbs Tr 1988	ansferred Of 1989	f-site 1990	
	 Waste Total	2,900	5,600	5,753	
Facility 00	00740				
F-Waste TOLUE 1,1,1	Constituents Reported to TRI: NE L-TRICHLOROETHANE	0 0	0 0	0 0	
	- Constituent Total	0	0	0	
F-Wastes F001 F003I F005	Reported to NCAR: 7005	23,595 0 18,920	28,435 34,760 0	0 0 0	
	Waste Total	42,515	63,195	0	
Facility 00	00743				
F-Waste METHA	Constituents Reported to TRI: NOL	0	0	0	
	- Constituent Total	0	0	0	
F-Wastes F002 F003	Reported to NCAR:	0 0	651 3,204	1,317 2,852	
	- Waste Total	0	3,855	4,169	
Facility 00	00744				
F-Waste METHY METHY	Constituents Reported to TRI: (L ETHYL KETONE (MEK) ANOL	70,000 15,000	0 0	0 0	
	- Constituent Total	85,000	0	0	
F-Wastes F0031 F005	s Reported to NCAR: F005	0 70,397	64,911 0	<b>4</b> 3,679 0	
	- Waste Total	70,397	64,911	43,679	
Facility 00	00750				
F-Waste ACETC	Constituents Reported to TRI: DNE	0	0	0	
	Constituent Total	0	0	0	
F-Wastes D0011 D0011 F002 F003	Reported to NCAR: 7003 7003F005	0 0 2,293 4,617	9,160 80 126 80	0 0 220 9,590	
	Waste Total	6,910	9,446	9,810	
Facility 00	00751				
F-Waste METHY METHY	Constituents Reported to TRI: YL ETHYL KETONE (MEK) YLENE CHLORIDE	68,200 6,040	29,600 21,700	75, <b>494</b> 11,211	

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## Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

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Constituent/Waste Code	Lbs Ti 1988	ransferred O 1989	ff-site 1990
TOLUENE	14.679	7.410	18.874
Constituent Total	88 919	58 710	105 579
	00,515	50,710	103,373
F-Wastes Reported to NCAR: F002	5,995	19,470	13,189
F003F005 F005	0 101.248	0 82,567	73,100 147,576
Waste Total	107 243	102 037	233 865
maste iotai	107,245	102,037	255,005
Facility 000/8/			
F-Waste Constituents Reported to TRI: METHYLENE CHLORIDE	: 3,621	1,799	0
Constituent Total	3,621	1,799	0
P-Wastes Peperted to NCAP.		·	
F002	3,597	3,175	0
F005		1,3/6	
• Waste Total	3,597	4,551	0
Facility 000768			
F-Waste Constituents Reported to TRI	:		1 550
D001D008F005	0	0	1,550
Waste Total	0	0	1,550
Facility 000769			
F-Waste Constituents Reported to TRI: 1,1,1-TRICHLOROETHANE	: 23,100	0	26,092
Constituent Total	23,100	0	26,092
F-Wastes Reported to NCAR:			
F001	3,600	0	0
F002 F003	19,500	21, 555	390
Waste Total	23,100	21,533	26,482
Facility 000771			
ACETONE	: 0	0	0
CYCLOHEXANE TOLLIENE	0	0	0
1,1,1-TRICHLOROETHANE	Ő	0	0
TRICHLOROETHYLENE			
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F002	0	0	9,010
Waste Total	0	0	9,010
Facility 000774			-
T Waste Constituents Described to mot	_		
n-BUTYL ALCOHOL	969	676	1,224

- onstituent/Waste Code	Lbs Transferred Off-site 1988 1989 1990			
METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	1,130 2,504 2,422 735	862 1,315 1,663 484	1,560 2,380 3,010 876	
Constituent Total	7,760	5,000	9,050	
F-Wastes Reported to NCAR: D001F003F005 F003F005 F005	0 0 26,652	22,000 18,098 0	50,055 32,751 0	
Waste Total	26,652	40,098	82,806	
acility 000775				
F-Waste Constituents Reported to TRI: ACETONE	18,456	42,205	10,807	
Constituent Total	18,456	42,205	10,807	
F-Wastes Reported to NCAR: *F001 F003 F005 Waste Total	0 12,989 39,062 52,051	0 33,748 12,536 46,284	3,610 10,807 40,690 55,107	
acility 000778				
F-Waste Constituents Reported to TRI: ACETONE	0	0	0	
Constituent Total	0	0	0	
F-Wastes Reported to NCAR: F002	39,920	0	0	
Waste Total	39,920	0	0	
acility 000794				
F-Waste Constituents Reported to TRI: ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	3,320 1,930 1,366 2,948 8,310 2,831	3,601 1,800 1,351 2,851 8,103 2,426	1,657 1,059 1,081 1,594 4,767 1,427	
Constituent Total	20,705	20,132	11,585	
F-Wastes Reported to NCAR: F003F005 F005	0 121,408	195,073 0	102,132 0	
 Waste Total	121,408	195,073	102,132	
acility 000798				
F-Waste Constituents Reported to TRI: n-BUTYL ALCOHOL METHANOL	0 0	0 0	0	

07/23/93 Comparison of F-Wastes Reported to NCAR and

07/23/93 Page 73	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Waste	e Code	Lbs Tra 1988	ansferred Of 1989	f-site 1990
	Constituent Total	0	0	0
F-Wastes Repo D001F003F00 F003F005 F005	rted to NCAR: 05	0 0 21,945	0 7,590 0	11,385 0 0
	Waste Total	21,945	7,590	11,385
Facility 000801				
F-Waste Const: ACETONE	ituents Reported to TRI	: 1,688	0	0
	Constituent Total	1,688	0	0
F-Wastes Report D001F003 F003	rted to NCAR:	0 16,110	0 12,300	11,070
	Waste Total	16,110	12,300	11,070
Facility 000808 F-Waste Consti ACETONE	ituents Reported to TRI	: 0	0	0
METHYL ETHY 1,1,1-TRICH	YL KETONE (MEK) HLOROETHANE	0 0	0 0	0
	Constituent Total	0	0	0
F-Wastes Report D001F003 F003	rted to NCAR:	0 3,280	0 17,220	12,300 0
	Waste Total	3,280	17,220	12,300
Facility 000809				
F-Waste Const ACETONE METHANOL	ituents Reported to TRI	: 0 0	0 0	0 0
	Constituent Total	0	0	0
F-Wastes Repor F003	rted to NCAR:	0	0	0
	Waste Total	0	0	0
Facility 000812				
F-Waste Consti METHANOL 1,1,1-TRICH	ituents Reported to TRI HLOROETHANE	: 0 592	6,539 11,490	0 0
	Constituent Total	592	18,029	0
F-Wastes Repor F001 F002 F002F003 F003 F005	ted to NCAR:	300 1,100 0 4,458 0	9,676 0 610 6,538 1,113	14,098 0 4,368 371

07/23 Page	/93 74	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Const	ituent/Wast	e Code	Lbs Tr 1988	ansferred O 1989	ff-site 1990
		Waste Total	5,858	17,937	18,837
Facil	ity 000816				
F-	Waste Const	ituents Reported to TRI:			
	METHYL ETH	YL KETONE (MEK)	0	0 71 136	2,772
	1.1.1-TRIC	HLOROETHANE	· 0	,1,150	4,299
	1,1,1 1	Constituent Total	0	71,136	24,351
F-	Wastes Repo	rted to NCAR:	0	0	6.000
	F002		õ	õ	12,600
	F005		0	0	2,800
		Waste Total	0	0	21,400
Facil	ity 000823				
- 4	Waste Const	ituents Reported to TRI	•		
-	METHYL ETH	YL KETONE (MEK)	0	20,208	39,295
TOLUENE		0	0	250	
	1,1,1-TRIC	HLOROETHANE	0	6,218	26,373
		Constituent Total	0	26,426	65,918
F-	Wastes Repo	rted to NCAR:			
	F003		23,240	0	0
	F003F005			40,584	62,812
		Waste Total	23,240	40,584	62,812
Facil	ity 000829.				
F-	Waste Const	ituents Reported to TRI	:		
	ACETONE		0	0	0
	METHYL ETH METHANOL	IL RETONE (MER)	0	0	0
		n			
		Constituent Total	U	0	0
F-	Wastes Repo	rted to NCAR:	-		
	D001D035F0	03F005	0		22,692
	F003F005 F005		20.540	22,575	0
	1000				
		Waste Total	20,540	22,575	22,692
Facil	ity 000834				
F-	Waste Const	ituents Reported to TRI	:		
	F003	- · · ·	0	240	0
		Waste Total	0	240	0
Facil	ity 000836				
F-	Waste Const	ituents Reported to TRT	:		
<b>L</b>	ACETONE	Tranto Heber con co INT	0	0	0
	METHYL ETH	YL KETONE (MEK)	0	0	0
	METHANOL		0	0	0
	XYLENE		õ	õ	ŏ

Page 75	F-Waste Constituents Rep	ported to TH	RI: 1988 to 1	990
Constituent/Waste	e Code	Lbs Tra 1988	ansferred Off 1989	-site 1990
	Constituent Total	0	0	0
F-Wastes Repo F005	rted to NCAR:	0	0	23,600
	Waste Total	0	0	23,600
Facility 000841				
F-Waste Const n-BUTYL AL ETHYLBENZE METHANOL 2-NITROPRO TOLUENE XYLENE	ituents Reported to TRI: COHOL NE PANE	1,020 385 24 225 5,493 0	0 0 0 0 0 0	0 0 0 0 0
	Constituent Total	7,147	0	0
F-Wastes Repo F005	rted to NCAR:	3,935	14,508	1,300
•	Waste Total	3,935	14,508	1,300
Facility 000842				
F-Waste Const ACETONE TOLUENE XYLENE	ituents Reported to TRI:	0 0 0	0 0 0	0 0 0
	Constituent Total	0	0	0
F-Wastes Repo F001	rted to NCAR:	0	5,400	1,439
	Waste Total	0	5,400	1,439
Facility 000844				
F-Waste Const METHANOL TOLUENE	ituents Reported to TRI:	0 0	0 0	0 0
	Constituent Total	0	0	0
F-Wastes Repo D035F003F0 F003F005 F005	rted to NCAR: 05	0 0 26,843	0 87,360 0	5,310 0 0
	Waste Total	26,843	87,360	5,310
Facility 000845				
F-Waste Const FREON 113 TETRACHLOR	ituents Reported to TRI: OETHYLENE	0 105,2 <b>4</b> 7	0 0	0 0
	- Constituent Total	105,247		0
F-Wastes Repo F001	rted to NCAR:	131,428	126,445	42,918
	- Waste Total	131,428	126,445	42,918

Comparison of F-Wastes Reported to NCAR and

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07/23/ Page	93 76	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990				
Constit	tuent/Waste	Code	Lbs Tra 1988	ansferred Off 1989	-site 1990	
Facili	ty 000847					
F-W	aste Consti ACETONE 1,1,1-TRICH	tuents Reported to TR	I: 79,200 0	36,900 0	29,700 0	
		Constituent Total	79,200	36,900	29,700	
F-W	astes Repor F003	ted to NCAR:	0	36,900	29,700	
		Waste Total	0	36,900	29,700	
Facili	ty 000850					
F-W	aste Consti METHANOL	ituents Reported to TR	I: 412	316	0	
		Constituent Total	412	316	0	
F-W	astes Repor F003F005 F005	ted to NCAR:	0 3,680	7,035 0	8,267 0	
		Waste Total	3,680	7,035	8,267	
Facili	ty 000852					
F-W	aste Consti METHYL ETHY TOLUENE XYLENE	tuents Reported to TR L KETONE (MEK)	I: 0 0 0	0 0 0	0 0 0	
		Constituent Total	0	0	0	
F-W	astes Report D001D035F00 D001F003F00 F005	rted to NCAR: 03F005 05	0 0 6,947	23,340 0	31,517 0 0	
- 11/		waste lotal	0,947	23,340	31,317	
Facili F-W	aste Consti ACETONE n-BUTYL ALC METHANOL TOLUENE	ituents Reported to TR COHOL	EI: 9,265 0 0 0	8,336 0 0 0	6,160 0 0	
		Constituent Total	9,265	8,336	6,160	
F-W	astes Repor F003	ted to NCAR:	1,050	15,378	6,160	
		Waste Total	1,050	15,378	6,160	
Facili	ty 000856					
F-W	aste Consti ACETONE TOLUENE	ituents Reported to TR	CI: 0 1,300	7,800 120,000	60,000 89,000	
		Constituent Total	1,300	127,800	149,000	

07/23/93 Page 77	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
	- Codo	Lbs Tra	ansferred Off	-site
was		1900		
F-Wastes Rep D001F003F F003 F003F005	orted to NCAR: 005	0 252,016 0	174,812 0 0	0 24,043 193,147
	- Waste Total	252,016	174,812	217,190
Facility 000857				
F-Waste Cons METHYL ET TOLUENE	tituents Reported to TRI: HYL KETONE (MEK)	7,543 7,985	11,120 11,120	15,423 3,027
	Constituent Total	15,528	22,240	18,450
F-Wastes Rep F003 F003F005 F005	orted to NCAR:	0 0 21,190	10,700 30,366 0	0 36,933 0
	Waste Total	21,190	41,066	36,933
Facility 000858				
F-Waste Cons 1,1,1-TRI	tituents Reported to TRI: CHLOROETHANE	0	0	0
	Constituent Total	0	0	0
F-Wastes Rep F001	orted to NCAR:	86,000	12,468	0
	Waste Total	86,000	12,468	0
Facility 000859				
F-Waste Cons ACETONE METHYLENE TOLUENE 1,1,1-TRI	tituents Reported to TRI: CHLORIDE CHLOROETHANE	0 0 0 0	0 0 0	0 0 0
	- Constituent Total	0	0	0
F-Wastes Rep D001F003 D001F005 D008F001 F001 F002 F003 F005	orted to NCAR:	0 0 182,268 0 224,774 43,228	0 0 111,996 10,368 149,974 33,532	145,536 21,624 35,046 0 0 0
	Waste Total	450,270	305,870	202,206
Facility 000860				
F-Waste Cons TRICHLORO	tituents Reported to TRI: ETHYLENE	18,871	0	0
	Constituent Total	18,871	0	0
F-Wastes Rep F002	orted to NCAR:	23,411	10,513	5,740

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07/23, Page	7/23/93 Comparison of F-Wastes Reported to NCAR and age 78 F-Waste Constituents Reported to TRI: 1988 to 1990				990
Const:	ituent/W	Naste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
			23,411	10,513	5,740
Facili	ity 0008	367			
F-V	Waste Co ACETONI METHYLI	onstituents Reported to TRI: E ENE CHLORIDE	0	0 0	0 0
		 Constituent Total	0	0	0
F-V	Wastes H F003	Reported to NCAR:	0	0	1,394
	FUUSFUU	 Waste Total		0	4,494
Facil:	ity 0008	377			
F-1	Waste Co METHANO	onstituents Reported to TRI: DL	0	0	0
•		 Constituent Total		0	0
F-V	Wastes F D001F00 D007D03 D022F00 F001 F002 F005	Reported to NCAR: 02F003F005 03FF005 01F002	0 0 0 0 0	0 0 61,110 0 860	2,508 856 424 10,668 813 0
		Waste Total	0	61,970	15,269
Facil	ity 0008	884			
F-1	Waste Co METHYL METHANO TOLUEN	onstituents Reported to TRI: ETHYL KETONE (MEK) DL E	0 0 0	0 0 0	0 0 0
		Constituent Total	0	0	0
F-1	Wastes I F003F00 F005	Reported to NCAR: 05	0 30,030	13,838	11,594
Facil	· +	waste Iotai	50,050	13,030	11,394
Facil.	Waste Co	onstituents Reported to TRI:			
	METHYLI	ENE CHLORIDE	0	0	0
		Constituent Total	0	0	0
F-1	Wastes H F001 F001F00 F002	Reported to NCAR: 02F003	2,442 0 2,416	0 3,670 0	11,859 0 0
		Waste Total	4,858	3,670	11,859

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07/23/93 Page 79	Comparison of F-wastes F F-Waste Constituents Rep	orted to TR	NCAR and I: 1988 to 1	990
Constituent/Waste	- Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
Facility 000888				
F-Waste Consti 1,1,1-TRICH	tuents Reported to TRI: LOROETHANE	0	0	1,000
	Constituent Total	0	0	1,000
F-Wastes Repor F001	ted to NCAR:	500	0	1,000
	Waste Total	500	0	1,000
Facility 000892				
F-Waste Consti ACETONE	tuents Reported to TRI:	45,000	0	0
METHYLENE C.		49,000		
	constituent Total	94,000	U U	U
F002 F003 ·	Led LO NCAR:	26,395 70,352	28,236 45,019	20,150 46,351
	Waste Total	96,747	73,255	66,501
Facility 000901				
F-Waste Consti ACETONE n-BUTYL ALC METHYL ETHY METHANOL METHYL ISOB TOLUENE XYLENE	tuents Reported to TRI: OHOL L KETONE (MEK) UTYL KETONE	0 0 0 0 0 0 0	0 0 0 0 0 0 0	0 0 0 0 0 0
	Constituent Total	0	0	0
F-Wastes Repor D001F003F00 F003F005 F005	ted to NCAR: 5  Waste Total	0 0 15,300 15,300	4,090 0 4,090	33,100 0 0 33,100
Facility 000903				
F-Waste Consti XYLENE	tuents Reported to TRI:	0	0	0
	 Constituent Total	0	0	0
F-Wastes Repor F003 F005	ted to NCAR:	8, <b>4</b> 70 1,155	10,550 0	18,700 0
	 Waste Total	9,625	10,550	18,700
Facility 000911				
F-Waste Consti ETHYLBENZEN TOLUENE XYLENE	tuents Reported to TRI: E	0 0 0	0 0 0	0 0 0

07/23/93 Page 80	Comparison of F-Wastes F-Waste Constituents Re	Reported to ported to T	NCAR and RI: 1988 to	1990
Constituent/Wast	te Code	Lbs Tra 1988	ansferred Of 1989	f-site 1990
	Constituent Total	0	0	0
F-Wastes Repo F002	orted to NCAR:	0	0	6,010
	- Waste Total	0	0	6,010
Facility 000914				
F-Waste Const	ituents Reported to TRI:			
ACETONE		0	0	0
n-BUTYL AL	COHOL	0	0	0
METHYL ETH	AYL KETONE (MEK)	0	0	0
METHANOL		0	0	0
XYLENE	_	0	0 	0
	Constituent Total	0	0	0
F-Wastes Repo	orted to NCAR:			
F003F005		26 125	59,622	66 007
FUUS	-	30,135		
	Waste Total	36,135	59,622	66,997
Facility 000917				
F-Waste Const	tituents Reported to TRI:			
ACETONE		0	0	0
n-BUTYL AI	LCOHOL	0	0	0
METHYL ETH	HYL KETONE (MEK)	0	0	0
METHANOL		0	0	0
TOLUENE		0	0	0
XYLENE		0	U 	0
	Constituent Total	0	0	0
F-Wastes Repo	orted to NCAR:			
F003F005		0	268,111	187,992
F005		128,720	0	0
	Waste Total	128,720	268,111	187,992
Facility 000921				
F-Waste Const	tituents Reported to TRI:			
D001F003	-	0	0	324,543
D005F003		0	0	2,500
F003		0	411,836	1,432,236
	Waste Total	0	411,836	1,759,279
Facility 000939	·			
F-Waste Const	tituents Reported to TRI:	:		
n-BUTYL AI	LCOHOL	0	0	750
METHYL ETH	HYL KETONE (MEK)	0	0	750
TOLUENE		0	0	750
XYLENE		0	0	1,500
	- Constituent Total	0	0	3,750
F-Wastes Repo	orted to NCAR:			
F002		0	0	500

07/23/ Page	07/23/93Comparison of F-Wastes Reported to NCAR andPage81F-Waste Constituents Reported to TRI: 1988 to 1990			<b>1990</b>	
Consti	ituent/Wast	e Code	Lbs 1 1988	Transferred ( 1989	off-site 1990
	F003 F005		0 0	0 139,920	17,585 176,485
		Waste Total	0	139,920	194,570
Facili	ity 000955				
F-W	Naste Const METHYL ISO	ituents Reported to TRI BUTYL KETONE	: 0	0	0
		Constituent Total	0	0	0
F-V	Nastes Repo F005	rted to NCAR:	0	90,000	164,520
		Waste Total	0	90,000	164,520
Facili	ity 000957				
F-W	Naste Const 1,1,1-TRIC	ituents Reported to TRI HLOROETHANE	: 0	13,537	10,192
		Constituent Total	0	13,537	10,192
F-W	Nastes Repo D001F003F0 F001 F002 F005	rted to NCAR: 05	0 0 0 0	0 14,541 601 2,428	2,833 12,619 0 0
		Waste Total	0	17,570	15,452
Facili	ity 000970				
F-V	- Naste Const XYLENE	ituents Reported to TRI	: 0	1,500	0
		Constituent Total	0	1,500	0
F-V	Vastes Repo D001F003F0 F001F002F0 F003F005	rted to NCAR: 05 03F005	0 0 0	4,800 0 10,500	0 3,650 7,900
		Waste Total	0	15,300	11,550
Facili	ity 000984				
F-V	Vaste Const ACETONE n-BUTYL AL METHYL ETH METHANOL METHYL ISO TOLUENE XYLENE	ituents Reported to TRI COHOL YL KETONE (MEK) BUTYL KETONE	: 0 0 0 0 0 0 0	0 0 0 0 0 0	0 0 0 0 0 0
		Constituent Total	0	0	0
F-V	Wastes Repo D001F003F0	rted to NCAR: 05	0	0	9,600
		Waste Total	0	0	9,600

Page 82 F-Waste Constituents Repo	ported to T	RI: 1988 to 3	1990
 Constituent/Waste Code	Lbs Tra 1988	ansferred Of: 1989	f-site 1990
Facility 000987			
F-Waste Constituents Reported to TRI: METHYL ETHYL KETONE (MEK)	0	0	0
METHANOL TETRACHLOROETHYLENE TOLUENE	0 0 0	0 0 0	0 0 0
 Constituent Total	0	0	0
F-Wastes Reported to NCAR: F001	0	7,256	12,093
F003	0	103,583	105,576
Waste Total	0	110,839	117,669
acility 000990			
F-Waste Constituents Reported to TRI: FREON 113	0	0	0
. Constituent Total	0	0	0
F-Wastes Reported to NCAR: F001 F002	0 0	3,527 2,520	950 7,250
 Waste Total	0	6,047	8,200
Facility 001008			
F-Waste Constituents Reported to TRI: TOLUENE XYLENE	0 0	18,758 0	0 11,727
 Constituent Total	0	18,758	11,727
F-Wastes Reported to NCAR: D001F003F005 F005	0	46,457	81,000 28 500
 Waste Total		46.457	109.500
Facility 001010	·		,
F-Waste Constituents Reported to TRI: METHYL ETHYL KETONE (MEK) TOLUENE	0	18,750 0	3,350
Constituent Total		18,750	
F-Wastes Perorted to NCAP.	·	207,00	0,000
D001D007F005 D001F005 D007F005 F002	0 0 0 0	9,500 3,750 14,500 1,000	27,000 0 22,100 1,250
 Waste Total	0	28,750	50,350
Facility 001024			
F-Waste Constituents Reported to TRI: METHYL ETHYL KETONE (MEK) TOLUENE	0	0	0
XYLENE	õ	õ	ŏ

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07/23/93 Page 83	7/23/93 Comparison of F-Wastes Reported to NCAR and age 83 F-Waste Constituents Reported to TRI: 1988 to 1990			1990
Constituent/W	aste Code	Lbs Tra 1988	nsferred Off 1989	-site 1990
	 Constituent Total	0	0	0
F-Wastes R D035F00 F003F00	eported to NCAR: 3F005 5	0 0	0 168,990	110,625 0
	Waste Total	0	168,990	110,625
Facility 0010	33			
F-Waste Co TETRACH	nstituents Reported to TRI: LOROETHYLENE	0	1,040	7,920
	Constituent Total	0	1,040	7,920
F-Wastes R F001	eported to NCAR:	0	6,930	7,920
	Waste Total	0	6,930	7,920
Facility 0010	40			
F-Waste Co F003	nstituents Reported to TRI:	0	0	37.947
	Waste Total		0	37,947
Facility 0010	43			
F-Waste Co TETRACH	nstituents Reported to TRI: LOROETHYLENE	0	40,256	55,740
	 Constituent Total	0	40,256	55,740
F-Wastes R F001	eported to NCAR:	0	41,999	55,740
	 Waste Total	0	41,999	55,740
Facility 0010	46			
F-Waste Co	nstituents Reported to TRI:	•	•	<u>^</u>
ACETONE METHYL	ETHYL KETONE (MEK)	0	0	0
METHANO	L	0	0	0
XYLENE		Ō	0	õ
	Constituent Total	0	0	0
F-Wastes R F003F00	eported to NCAR: 5	0	23,590	39,200
	Waste Total	0	23,590	39,200
Facility 0010	50			
F-Waste Co FREON 1	nstituents Reported to TRI: 13	0	0	0
	Constituent Total	0	0	0
F-Wastes R F001	eported to NCAR:	0	26	0

Page	84	F-waste Constituents	Reported to 1	IRI: 1988 CO	1990
Const	ituent/Wast	e Code	Lbs Tr 1988	ansferred O 1989	ff-site 1990
	F002 F003 F003F005 F005		0 0 0 0	1,592 25,307 0 338	0 20,051 18,619 0'
		Waste Total	0	27,263	38,670
Facil	ity 001059		•		
F-	Waste Const F001	ituents Reported to T	RI: 1,435	1,152	1,363
	F003	Waste Total	1,435	1,235	1,363
Facil	ity 001073				
F-	Waste Const ACETONE	ituents Reported to T	RI: 0	0	0
		Constituent Total	0	0	0
F-	Wastes Repo F003	orted to NCAR:	68,046	72,567	26,080
		Waste Total	68,046	72,567	26,080
Facil	ity 001074				
F-	Waste Const D035F001F0 F001 F001F002	ituents Reported to T 003F005	RI: 0 0 0	0 0 0	15,750 1,260 15,080
		Waste Total	0	0	32,090
Facil	ity 001094				
F-	Waste Const F001 F002	ituents Reported to T	RI: 0 0	0 0	400 400
	F005		0	0	6,400
		Waste Total	0	0	7,200
Facil	ity 001095.				
F-	Waste Const D007F005	ituents Reported to T	RI: 0	0	3,303
		Waste Total	0	0	3,303
Facil	ity 001107.				
F-	Waste Const F001 F003 F005	ituents Reported to T	RI: 0 0 0	0 0 0	2,289 1,977 2,550
		Waste Total	0	0	6,816
Facil	ity 001127.				
F-	Waste Const FREON 113	ituents Reported to T	RI: O	0	0

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Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990 -

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# Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990

	Lbs Trans	ferred Off	-site
Constituent/Waste Code	1988	1989	1990
Constituent Total	0	0	0
F-Wastes Reported to NCAR: F002	0	0	53,654
Waste Total	0	0	53,654
Facility 001129			
F-Waste Constituents Reported to TRI: 1,1,1-TRICHLOROETHANE	0	0	0
- Constituent Total	0	0	0
F-Wastes Reported to NCAR: F002	0	0	9,600
Waste Total	0	0	9,600
Facility 001132			
F-Waste Constituents Reported to TRI: n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) TOLUENE XYLENE	0 0 0 0	0 0 0 0	0 0 0 0
Constituent Total			
	Ū	Ū	Ŭ
F-Wastes Reported to NCAR: D001F003 D001F005 F002	0 0 0	0 0 0	11,550 82,756 593
Waste Total	0	0	94,899
Facility 001134			
F-Waste Constituents Reported to TRI: METHYL ETHYL KETONE (MEK) TOLUENE	: 0 0	0 0	11,585 11,786
Constituent Total	0	0	23,371
F-Wastes Reported to NCAR: D001F003 F003 F003F005	0 0 0	0 0 0	6,900 38,300 3,200
Waste Total	0	0	48,400
Facility 001142			
F-Waste Constituents Reported to TRI: n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK)	: 0 0	0 0	11,700 6,900
Constituent Total	0	0	18,600
F-Wastes Reported to NCAR: F003F005	0	0	821,770
Waste Total	0	0	821,770

07/2 Page	23/93 = 86	Comparison of F-Wastes F-Waste Constituents Re	Reported to NCA ported to TRI:	R and 1988 to 1	990
Cons	stituent/Was	te Code	Lbs Transf 1988	erred Off 1989	-site 1990
rac.					
1	F-Waste Const METHYL ET	tituents Reported to TRI: HYL KETONE (MEK)	0	0	·0
		Constituent Total	0	0	0
]	F-Wastes Repo F003	orted to NCAR:	0	0	388,519
		Waste Total	0	0	388,519
Fac	ility 001156				
]	F-Waste Cons	tituents Reported to TRI:	_		
	n-BUTYL A METHYL ET	LCOHOL HYL KETONE (MEK)	0	0	0
		- Constituent Total	0	0	0
1	F-Wastes Rep	orted to NCAR:			
	F001 • F005		0 0	0	2,107 15,170
		- Waste Total	0	0	17,277
Fac:	ility 001162				
]	F-Waste Const 1,1,1-TRI	tituents Reported to TRI: CHLOROETHANE	0	0	7,250
		- Constituent Total	0	0	7,250
1	F-Wastes Ren	orted to NCAR:			
	F001 F003		0 0	0 0	5,294 200
		- Waste Total	0		5,494
Fac	ility 001165				-
]	F-Waste Const METHANOL	tituents Reported to TRI:	0	0	0
		- Constituent Total	0	0	0
1	F-Wastes Rep	orted to NCAR:			
•	F003F005		0	0	21,054
		Waste Total	0	0	21,054
Fac	ility 001174				
1	F-Waste Const ETHYLENE ( METHYL ETH TOLUENE 1 1.1-TRI(	tituents Reported to TRI: GLYCOL MONOETHYL ETHER HYL KETONE (MEK) CHLOROETHANE	0 0 0	0 0 0	750 750 750 0
	_, _,		0		2,250
]	F-Wastes Rep F001 F003	orted to NCAR:	0 0	0 0	15, <b>4</b> 97 3,017

07/23/93 Page 87	Comparison of F-Wastes Reported to NCAR and F-Waste Constituents Reported to TRI: 1988 to 1990			
Constituent/Wast	e Code	Lbs T 1988	ransferred O 1989	ff-site 1990
	Waste Total	0	0	18,514
Facility 001177				
F-Waste Const 1,1,1-TRIC	ituents Reported to TRI HLOROETHANE	0	0	0
	Constituent Total	0	0	0
F-Wastes Repo F001 F003	rted to NCAR:	0	0	1,716 788
F005		Ő	0	1,718
	Waste Total	0	0	4,222
Facility 001190				
F-Waste Const METHANOL TOLUENE	ituents Reported to TRI	[: 0 0	0 0	250 250
	Constituent Total	0	0	500
F-Wastes Repo F005	rted to NCAR:	0	0	61,414
	Waste Total	0	0	61,414
Facility 001193				
F-Waste Const	ituents Reported to TRI	:	0	0
METHYL ETH	YL KETONE (MEK)		0 	0 
	Constituent Total	0	0	0
F-Wastes Repo F003 F005	rted to NCAR:	0 0	0 0	94,469 116,899
	Waste Total	0	0	211,368
Facility 001197				
F-Waste Const n-BUTYL AL	ituents Reported to TRI COHOL	C: 0	0	750
	Constituent Total	0	0	750
F-Wastes Repo D001F002F0	rted to NCAR: 03F005	0	0	41,900
	Waste Total	0	0	41,900
TOTAL FOR ALL FA	CILITIES:	1988	1989	1990
F-Waste Constit F-Wa	uents Reported to TRI stes Reported to NCAR	8,104,871 29,309,689	6,424,963 29,146,745	2,911,618 34,022,672

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APPENDIX F RESULTS OF COMPARISON OF SINGLE-MEDIA DATA AND TRI THRESHOLDS

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Qty (lbs)	
Chemical	
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Facilities Exceeding the Processed or Manufactured Threshold \*

2262 ZINC 2262 COPPER

21,385 208,206

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sIC Chemical Qty (lbs)

\* Facilities Exceeding the Processed or Manufactured Threshold \*

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None

sIC Chemical Qty (lbs)

 $^{\star}$  Facilities Exceeding the Processed or Manufactured Threshold  $^{\star}$ 

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None

-Page Facilities Reporting Water Releases to NPDES over the TRI Reporting Threshold for 1990 07/22/93

sic Chemical Qty (lbs)

 $^{\star}$  Facilities Exceeding the Processed or Manufactured Threshold  $^{\star}$ 

4952 COPPER 4952 ZINC

14,497 32,926

Qty (lbs)	
Chemical	
SIC	

 $\star$  Facilities Exceeding the Processed or Manufactured Threshold  $\star$ 

13,654 234,725 319,170,240	42,700 30,500 36,600 109,800	765,005	118,798	50,857 10,680 245,625,132 115,132 115,138 115,138 13,817 13,817 101,321 18,422 18,422 18,422	39,483 43,870 26,322 114,062 144,771 78,966	28,060 1,220,000,000	38,033,500 32,788 22,296 15,738 151,150 78,690 94,008	16,242 10,258 76,937	352,331	222,470,050 70,786 80,898 30,337 141,572 141,572 30,337 505,613,750 10,112 10,112
									(FUME OR DUST)	
COPPER ARSENIC PHENOL	TOLUENE SELENIUM MERCURY SILVER	PHENOL	TOLUENE	ANTIMONY SILVER PHENOL BARIUM TOLUENE SELENIUM LEAD MERCURY COPPER SILVER	TOLUENE SELENIUM LEAD MERCURY COPPER SILVER	TOLUENE PHENOL	BARIUM TOLUENE LEAD MERCURY COPPER ARSENIC SILVER TOLUENE	MERCURY COPPER ARSENIC	ALUMINUM	BARIUM TOLUENE SELENIUM LEAD MERCURY COPPER COPPER COPALT COBALT CADMIUM
N/A N/A N/A	N/A N/A N/A N/A	2011	2111	2221 2221 22221 22221 22221 22221 22221 22221 22221 22221 22221 22221 22221	2221 2221 2221 2221 2221 2221	2221 2221	55555555555555555555555555555555555555	2252 2252 2252	2253	2258 225558 22558 22558 2255757 22558 2255757 2255757 2255757 22557577 22557577777777

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sic	Chemical	Qty (1bs)
2258 2258	ALUMINUM (FUME OR DUST) SILVER	10,112 556,175
2262	COBALT	116,053
22269 22269 22269 22269 22269	PHENOL TOLUENE SELENIUM LEAD MERCURY CADMIUM	17, 306, 707 16, 013 26, 688 21, 350 64, 050 64, 050
2269	PHENOL	351,717
2269 2269 2269	TOLUENE MERCURY PHENOL	11,895 14,640 2,745,000,000
2298	SILVER	977,353
2322 2322 2322 2322 2322 2322 2322 232	PHENOL TOLUENE SELENIUM LEAD MERCURY COPPER COPPER COBALT CADMINUM FUME OR DUST) SILVER	85,044 120,032 87,296 87,296 87,296 120,032 10,912 10,912 10,912 578,336
3321 3321	Toluene Selenium	74,291 134,633
3339	CYANIDE COMPOUNDS	497,394
3471 3471	SELENIUM COBALT	132,844 1,195,592
3519 3519 3519 3519 3519	TOLUENE MERCURY COPPER COBALT ARSENIC	58,772 16,456 423,157 23,509 2,565,977
3536	SILVER	89,213
33714 37714 37714 37714 37714 4172 4172 4172 4172	TOLUENE SELENIUM LEAD MERCURY COPPER COPPER CADMIUM ALUMINUM (FUME OR DUST) SILVER	383,690 115,107 172,661 57,554 57,554 19,185 19,185 19,185 19,266

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Qty (lbs)	
iIC Chemical	

\* Facilities Exceeding the Otherwise Used Threshold \*

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50,349 20,979 12,587 12,587	62,049	33,947 46,757	47,183	37,988 34,823 12,353	64,050 10,169 30,500	62,296	44,041 73,353	20,714	58,721	14,647 54,900	20,506	12,189 18,284 12,525	57,6 <b>4</b> 5 17,080	14,234	13,366 13,366 28,960 33,643 13,364 14,601 12,161 12,161 12,161 12,161	
TOLUENE SELENIUM MERCURY COPPER	BARIUM	TOLUENE Arsenic	ARSENIC	SILVER PHENOL TOLUENE	LEAD COPPER SILVER	PHENOL	COPPER SILVER	MERCURY	SILVER	COPPER SILVER	SILVER	MERCURY PHENOL SILVER	TOLUENE MERCURY	PHENOL	TOLUENE SELENIUM LEAD MERCURY TOLUENE LEAD COPPER CHLORINE SELENIUM LEAD	
N/A N/A N/A	2086	2211 2211	2211	2231 2251 2252	2254 2254 2254	2257	2257 2257	2259	2261	2261 2261	2261	2262 2262 2262	2262 2262	2269	22281 22281 22281 22281 22281 22297 2297 2297 2451 2451 2451	

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SIC	Chemical	Qty (lbs)
2653	PHENOL	12,894
2834	MERCURY	21,960
3353	PHENOL	47,742
3411	SILVER	47,565
3425 3425 3425	TOLUENE MERCURY COPPER	10,120 10,795 62,072
3442	ARSENIC	31,959
3613 3613 3613	TOLUENE MERCURY COPPER	52,342 11,632 13,958
3621 3621	TOLUENE MERCURY	45,857 45,857
3676	LEAD	40,275
3679	TOLUENE	22,326
3829	MERCURY	10,614

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---Facilities Reporting POTW Releases to Pretreatment over the TRI Reporting Threshold for 1988 Page 07/22/93

Qty (lbs)

Chemical  .

SIC	Chemical	Qty (lbs)
* Facilit	ies Exceeding	the Processed or Manufactured Threshold *
N/A	ARSENIC	71,039
N/A	PHENOL	1,323,022,282
2026	TOLUENE	64,604
2032	PHENOL	606,762
2221	PHENOL	42,122,025
2231	ARSENIC	223,769
2231	SILVER	11, 712
2251	MERCURY	10,361
2251	TOLUENE	51,806
2252	SELENIUM	12,414
2252	PHENOL TO LIENE	1,050,50 150,150
7677	VEDCIBV	21,02 20 758
2252	SILVER	10, 345
2257	PHENOL	59,870
2269	PHENOL	247,297
2281	PHENOL	6,717,625
2298	SILVER	89,222
2321	PHENOL	78,927
2843	PHENOL	61,268
3425	PHENOL	16,145,938
3519	ARSENIC	135,974
3519	PHENOL	2,948,806
3676	PHENOL	2,812,890
* Facilit	ties Exceeding	the Otherwise Used Threshold *
2016 2221	PHENOL PHENOL	16,144 37,028 10,201

016 121	PHENOL	16, 144 37, 028
52	PHENOL	10,201
52	MERCURY ARSENIC	10,190 10,190
57	SILVER	13,491
61	SILVER	25,639
62 69	PHENOL MERCURY PHENOL	14,903 13,334 14,889
69	PHENOL	14,870
39	CHLORINE PHENOL	26,873 11,011

٦ Facilities Reporting POTW Releases to Pretreatment over the TRI Reporting Threshold for 1989 Page 07/22/93

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\* Facilities Exceeding the Processed or Manufactured Threshold \*

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700,444	8,521,649 660,428 7,840,475 80,057 25,206 20,053,549	39,166	63,993	38,406	62,357	140,898	405,488	313,266,576	165,870	65,112 13,465 11,578 13,745 22,335	46,749 30,053 10,018 43,410	219,335 959,204	3,878,378	22, 355 132, 097	26,980,692	487,337 438,807 1,099,699 409,861 2,208,580	114,607 43,527 171,779 169,494	5,606,623
PHENOL	CYANIDE COMPOUNDS TOLUENE PHENOL ARSENIC TOLUENE PHENOL	ARSENIC	PHENOL	SILVER	TOLUENE	ARSENIC	PHENOL	PHENOL	CHLORINE	PHENOL TOLUENE TOLUENE MERCURY SELENIUM	COPPER TOLUENE SELENIUM MERCURY	SILVER Barium	TOLUENE	MERCURY TOLUENE	PHENOL	MERCURY SELENIUM MERCURY LEAD TOLUENE	MERCURY COPPER LEAD TOLUENE	PHENOL
N/A	2016 2016 2200 2200 2211	2211	2257	2258	2261	2262	2269	2281	2284	2321 2321 2328 2328 2328 2328	2396 2396 2396 2396	2641 2641	2843	2844 2844	3425	3471 3471 3471 3471	3471 3471 3471 3471	3519
2 Facilities Reporting POTW Releases to Pretreatment over the TRI Reporting Threshold for 1989 Page 07/22/93

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\* Facilities Exceeding the Otherwise Used Threshold \*

2221	CHROMIUM	11, 311
2253 2262 2269	SILVER PHENOL PHENOL	14,738 21,941 14,644
2269	PHENOL	18,321
2293	PHENOL	10,195
2821 2821 3281	ALUMINUM (FUME OR DUST) Toluene Phenol	17,723 10,502 16,592
3356	MANGANESE AND COMPOUNDS	21,781
3442	ARSENIC	10,331
3443	ARSENIC	16,894
3714	TOLUENE	10,093

-Facilities Reporting POTW Releases to Pretreatment over the TRI Reporting Threshold for 1990 Page 07/22/93

Qty (lbs)	****	Manufactured Threshold 🕈
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2024	PHENOL	1,492,663
2211	ARSENIC	55,992
2251 2251	The Phenol	664,365 230,844,779
2254	PHENOL	3,515,921
2257	PHENOL	53,577
2262	PHENOL	31,378
2262 2262	ZINC	35,556 76,191
2269	PHENOL	495,427
2269	PHENOL	8,400,459
2281	PHENOL	264,679,317
2284	CHLORINE	156, 388
2321 2321	MERCURY PHENOL	25,475 62,995
2843	PHENOL	65,801
3411	PHENOL	154,537
3425	PHENOL	28,328,606
3471 3471 3471 3471 3471 3471	LEAD SILVER COPPER MERCURY TOLUENE SELENIUM	20,681 14,218 23,266 16,803 19,388
3496	PHENOL	98,391
3519	PHENOL	3,506,294
3629 3629 3629 3629	LEAD SELENIUM MERCURY COPPER	20,601 13,734 68,670 13,734
3676	PHENOL	83,356
3692	PHENOL	19,666,082
3714	PHENOL	66,831,600

N Facilities Reporting POTW Releases to Pretreatment over the TRI Reporting Threshold for 1990 Page 07/22/93

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SIC	Chemical	Qty (1bs)
* Facilit	ies Exceeding the Otherwise Used Threshol	<b>ئ</b> •
2016	CYANIDE COMPOUNDS	12,067
2251	PHENOL	10,707
2252	PHENOL	11,332
2253	MERCURY	20,359
2262 2269	MERCURY PHENOL	15,500 17,658
2269	CADMIUM	10,162
2269	PHENOL	18,385
2439 2834	CHLORINE ARSENIC	10,289 14,736
3443	ARSENIC	15,981

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## APPENDIX G RESULTS OF COMPARISON OF TRI DATA ACROSS REPORTING YEARS

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07/22/	781 Facilities	ith Discrep	ancies	in Reporting	Status Over Time:
rage		Reporting	status 1080	in: 1990	
SIC	Cnemical	10LT			
20	COPPER	Х	×		
2013	HYDROCHLORIC ACID	х		Х	
2016 2016	AMMONIA CHLORINE	××	××	××	•
2016	AMMONIA	×	×		
2021	PHOSPHORIC ACID	×	×	Х	
2026	PHOSPHORIC ACID	ХХ		Х	
2026	PHOSPHORIC ACID	x		x	
2041	CHLOR INE	X	x	x	
2046 2046	HYDROCHLORIC ACID SULFURIC ACID	××	××	××	
2048	PHOSPHORIC ACID	x	х	Х	
2077	AMMONIA	Х		Х	
2077	CHLORINE	x		Х	
2077	AMMONIA	X		x	
2082	AMMONIA	x	x	х	
2086	PHOSPHORIC ACID	X		x	
2087 2087	METHANOL TOLUENE	××	×	××	
2111 2111 2111	ACETONE METHYL ETHYL KETONE (MEK) XYLENE	x x x	•	×××	
22	SULFURIC ACID	х х		x	
2200	AMMONIUM SULFATE (SOLUTION)	×	x	×	
2211	CHLORINE	х	X	Х	
2211	CHLORINE	x		Х	
2231	CHROMIUM COMPOUNDS	х	X	x	
2251	CHLORINE	X		x	
2252	AMMONIUM SULFATE (SOLUTION)	x	×		
2252	CHLORINE	x	X		
2257	SULFURIC ACID	X		x	
2257	BIPHENYL	ХХ		X	

X = reported

TRI Facilities with Discrepancies in Reporting Status Over	Time:	
TRI Facilities with Discrepancies in Reporting Status	Over	
TRI Facilities with Discrepancies in Reporting	Status	
TRI Facilities with Discrepancies in	Reporting	
TRI Facilities with Discrepancies	1n	
TRI Facilities with	Discrepancies	
TRI Facilities	with	
TRI 1901	Facilities	1000
	TRI	100

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07/22	/93 TRI Facilities	vith Discr	epan	cies	in Rep
sIC	chemical	Report1 1987 19	ng s 88	catus 1989	in: 1990
2257	BIPHENYL	x	×	8 1 1 1	×
2260	1, 1, 1-TRICHLOROETHANE	×		×	×
2261 2261	ETHYLENE GLYCOL GLYCOL ETHERS	××		××	××
2261	SULFURIC ACID	×			×
2262 2262	AMMONIA CHLORINE	××			××
2269	AMMONIUM SULFATE (SOLUTION)	x			×
2271 2271	AMMONIA CHLORINE	××	×		××
2272	ETHYLENE GLYCOL	×	×		x
2272	CHLORINE	×		×	
2282	ETHYLENE GLYCOL	×		×	×
2283 2283	AMMONIA AMMONIUM SULFATE (SOLUTION)	x	×		××
2283	AMMONIUM SULFATE (SOLUTION)	x		×	
2293 2293 2293	ACETONE CHLORINE SULFURIC ACID	×××	×××		×××
2295	SULFURIC ACID	x			×
2295	DECABROMODIPHENYL OXIDE	x		×	×
2296	SULFURIC ACID	x	×		x
2299 2299	CHLORINE COPPER COMPOUNDS	×	×		××
2421 2421	ARSENIC COPPER	××		××	××
2436 2436	FORMALDEHYDE PHENOL	x x			××
2491 2491 2491	ARSENIC COMPOUNDS CHROMIUM COMPOUNDS COPPER COMPOUNDS	×××		×××	×××
2491	COPPER COMPOUNDS	x		×	
2491	CHROMIUM COMPOUNDS	x		×	Х
2491 2491	ARSENIC COMPOUNDS CHROMIUM COMPOUNDS	××		××	××
2493	BARIUM COMPOUNDS	X			×
= X	reported				

07/22/	'93 TRI Facilities	with Discreps	ancies 1	in Reporting Status Over Time:	::
Page	3 TA8/ CO T220	Reporting	status	in: 1000	
SIC	Chemical	198/ 1988	T 967		
2493	SULFURIC ACID	x		х	
2499 2499	ACETONE TOLUENE	××	××	××	
2511	XYLENE	x	×	X	,
2511	TOLUENE	х х		Х	
2511 2511	ACETONE METHYL ETHYL KETONE (MEK)	x x x		×	
2511 2511	N-BUTYL ALCOHOL XYLENE	××		x	
2511 2511	ACETONE METHYL ETHYL KETONE (MEK)	x x x		x	
2511 2511 2511 2511 2511 2511 2511	ACETONE BIS (2-ETHYLHEXYL) PHTHALATE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	× ××××		*****	
2511	n-BUTYL ALCOHOL	х	х	Х	
2511 2511 2511 2511 2511 2511 2511	ACETONE BIS (2-ETHYLHEXYL) PHTHALATE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	* *****		× × × × × ×	
2511	METHYL ETHYL KETONE (MEK)	х х		Х	
2511	ACETONE	×	×	х	
2511	ACETONE	x		Х	
2511 2511 2511 2511 2511	ACETONE METHYL ETHYL KETONE (MEK) METHANOL TOLUENE	××××	× × × ×	X	
2511	XYLENE	ХХ		x	
2511	TOLUENE	x		х	
2511	n-BUTYL ALCOHOL	х	×	Х	
2511	GLYCOL ETHERS	Х		x	
2511	ACETONE	ХХ		Х	
2511	п-витур Аьсоноь	×	x	х	

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X = reported

07/22	/93 TRI Facilities v 1087 +0 1090	ith Discr	epanci	es İ	n Reporting St	atus (	Dver T	lme:
SIC	Chemical	Report1 1987 19	ng sta 88 19	tus 89	in: 1990			
2511	n-BUTYL ALCOHOL	x		×				•
2511	n-BUTYL ALCOHOL	×		x	х			
2511 2511	n-BUTYL ALCOHOL XYLENE	××		××				
2512	METHYL ETHYL KETONE (MEK)	×	x		x			
2512	TOLUENE	×	×		х			
2512	METHYL ETHYL KETONE (MEK)	x		×	Х			
2514 2514 2514	CHROMIUM CYANIDE COMPOUNDS NICKEL COMPOUNDS	×××		×××				
2514	XYLENE	x	×		×			
2517	XYLENE	×	×		X			
2519	METHYL ETHYL KETONE (MEK)	×		×				
2521 2521 2521	METHYL ETHYL KETONE (MEK) METHANOL TOLUENE	×××		×××				
2521 2521 2521 2521 2521 2521	ACETONE METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE		×××××		****			
2521	METHANOL	x	×		x			
2521	n-BUTYL ALCOHOL	x		×				
2521	GLYCOL ETHERS		×		X			
2621	POLYCHLORINATED BIPHENYLS (PC	вх		X				
2657 2657	DI-N-BUTYL PHTHALATE METHANOL	××		××	××			
2754	METHYL ETHYL KETONE (MEK)	×		×	X			
2815	TEREPHTHALIC ACID	x		×				
2818 2818 2818	ACRYLAMIDE AMMONIA SULFURIC ACID	×××		×××	×××			
2819 2819	ARSENIC COMPOUNDS CHROMIUM COMPOUNDS	×	××		××			
2819	SILVER COMPOUNDS	×		×	Х			
2819 2819	CHLORINE HYDROCHLORIC ACID	××		××	××			
×	reported							

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Time:		•												
Over														
atus														
lg St														
portir														
in Re	in: 1990	× × × × ×	× ××	* ***	×	×			××××	×××	*****	××	×××××	××××××
cies i	tatus 1989	×××××	×	×××	×	×	×	×××	×× ×	××	*****	××	*****	*****
with Discrepan	Reporting s 1987 1988	×××××	× ×	< ×××	x	x	х	×××	***	×××	*****	××	×××××	×××××
'93 TRI Facilities	chemical	ALUMINUM OXIDE AMMONIA ARSENIC ARRANUM COPPER	PHOSPHORIC ACID	SULFURIC ACID n-BUTYL ALCOHOL FORMALDEHYDE METHANOL	HYDROCHLORIC ACID	STYRENE	BIPHENYL	ACETONE METHYL ETHYL KETONE (MEK) STYRENE	ETHYLBENZENE Na Phthalene Toluene Xylene	AMMONIA DI-N-BUTYL PHTHALATE METHANOL	ACRYLIC ACID DI-N-OCTYL PHTHALATE 4,4'-METHYLENE DIANILINE 1-NAPHTHYLAMINE PHTHALIC ANHYDRIDE TOLUENE DIISOCYANATE TOLUENE DIISOCYANATE	ACRYLIC ACID AMMONIA	ACRYLONITRILE BUTYL ACRYLATE ETHYL ACRYLATE METHYL ACRYLATE METHYL ACRYLATE METHYL METHACRYLATE STYRENE	ACRYLONITRILE BUTYL ACRYLATE METHYL ACRYLATE METHYL METHACRYLATE STYRENE VINYL ACETATE
07/22/	SIC	2819 2819 2819 2819 2819	2819 2821	2821 2821 2821 2821	2821	2821	2821	2821 2821 2821	2821 2821 2821 2821 2821	2821 2821 2821	2821 2821 2821 2821 2821 2821 2821 2821	2821 2821	2821 2821 2821 2821 2821 2821 2821	2821 2821 2821 2821 2821 2821

X = reported

ng Status Over Time:			•													
in Reporti	in: 1990	1	×	****		×××	x	××	××	*****	××	**** ***	××	×××	××	×
ictes	status 1989	x	×	****	×	××	×	××			××	******	××	× ××		×
crepar	ting 1988									××× × ×				×	××	
th Dis	Repor 1987	×	×	×××××	×	×××	×	××	××	******	××	******	××	× ××		×
2/93 TRI Facilities wi	Chemical	TEREPHTHALIC ACID	SULFURIC ACID	ANTIMONY COMPOUNDS HYDROCHLORIC ACID MANGANESE COMPOUNDS 2-METHOXYETHANOL PHOSPHORIC ACID	TEREPHTHALIC ACID	AMMONIA ANTIMONY COMPOUNDS ETHYLENE OXIDE	CHLORINE	CHLORINE SULFURIC ACID	ETHYLENE GLYCOL HYDROCHLORIC ACID	n-BUTYL ALCOHOL DIETHYL PHTHALATE DIMETHYL PHTHALATE ETHYLENE GLYCOL METHANOL PHTHALIC ANHYDRIDE SULFURIC ACID TOLUENE	HYDROCHLORIC ACID NITRIC ACID	BENZYL CHLORIDE n-BUTYL ALCOHOL GLYCOL ETHERS NAPHTHALENE SULFURIC ACID TEREPHTHALIC ACID TEREPHTHALIC ACID 1,2,4-TRICHLOROBENZENE XYLENE	HYDROCHLORIC ACID	BIPHENYL DIETHANOLAMINE METHANOL 1,2,4-TRICHLOROBENZENE	MALEIC ANHYDRIDE PHOSPHORIC ACID	I METHANOL
07/22	SIC	2824	2824	2824 2824 2824 2824 2824 2824	2824	2824 2824 2824	2833	2834 2834	2834 2834	2834 2834 2834 2834 2834 2834 2834 2834	2834 2834	2841 2841 2841 2841 2841 2841 2841 2841	2842 2842	2843 2843 2843 2843 2843	2843 2843	2843

.

X = reported

07/22/	/93 TRI Facilities v 1087 to 1090	ith Discrep	ancies i	n Reporting Status Over Time:	
rage SIC	chemical	Reporting 1987 1988	status 1989	in: 1990	
2843	XYLENE	Х			
2843	SULFURIC ACID	×	×	Х	
2843 2843 2843 2843 2843 2843 2843 2843	ANTIMONY COMPOUNDS BENZYL CHLORIDE BIPHENYL BUTYL BENZYL PHTHALATE CHLORINE GLYCOL ETHERS 1,2,4-TRICHLOROBENZENE XYLENE	*****	******		
2843 2843 2843 2843 2843 2843	CHLOROACETIC ACID FORMALDEHYDE METHANOL SULFURIC ACID TETRACHLOROETHYLENE	****	×××××	××××	
2843	1,2,4-TRICHLOROBENZENE	×	×	х	
2851 2851 2851 2851 2851 2851	n-BUTYL ALCOHOL GLYCOL ETHERS METHYL ISOBUTYL KETONE 1,1,1-TRICHLOROETHANE XYLENE	×××××	****	****	
2851 2851 2851	ACETONE n-BUTYL ALCOHOL TOLUENE	×××	×××	×××	
2851	ETHYLENE GLYCOL	x	×	X	
2851 2851 2851	METHYL ETHYL KETONE (MEK) METHANOL TOLUENE	×××	x	×××	
2851 2851	METHYL ISOBUTYL KETONE XYLENE	x x x	×	X	
2851 2851 2851 2851 2851 2851	ACETONE n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL XYLENE	****	×× ××	××××	
2851 2851 2851 2851 2851	GLYCOL ETHERS METHANOL METHYL ISOBUTYL KETONE XYLENE	××××	x xx	××××	
2851 2851 2851	BARIUM COMPOUNDS DI-N-OCTYL PHTHALATE ETHYLBENZENE	×××	×××	X	
2851	BARIUM COMPOUNDS	x	×	Х	
2851 2851	GLYCOL ETHERS MANGANESE COMPOUNDS	××	××	Х	

.

X = reported

2851 GLYCOL ETHERS 2851 MANGANESE COMPOUNDS

17/22/	781 Facilities	vith Discrepa	incies i	n Reporting S	tatus Over Time:
age	B 1987 TO 1990	Reporting	status	in: 1000	
	Chemicai				
851	METHYL ISOBUTYL KETONE	X	×	×	-
865	GLYCOL ETHERS	X	×	х	
2865	FORMALDEHYDE	x	x		
2 Y C	CHI OBOBENZENE	×	×	X	
865	CHROMIUM COMPOUNDS	×	×	×:	
2865 2865	COBALT COMPOUNDS CODDER COMPOUNDS	< ×	< ×	< ×	
2865	HYDROCHLORIC ACID	×:	×	×	
2865 2865	NICKEL COMPOUNDS NITTRIC ACID	××	< ×	××	
2865	PHENOL	×	×	×	
2865 2865 2865	PICRIC ACID SULFURIC ACID o-TOLUIDINE	×××	×××	×××	
2869	ACRYLIC ACID	××	××	××	
2869	CHLORINE GLYCOL ETHERS	××	××	××	
2869	SULFURIC ACID	x		X	
2869 2869	MALEIC ANHYDRIDE SULFURIC ACID	x X X		××	
2869 2869 2869 2869	n-BUTYL ALCOHOL HYDROCHLORIC ACID SULFURIC ACID XYLENE	****	××××	× × ×	
2869	NITRIC ACID	×	×	x	
2869	ZINC COMPOUNDS	×		X	
2869	PHTHALIC ANHYDRIDE	х		X	
2869	TOLUENE	×	x		
2873 2873 2873	AMMONIUM SULFATE (SOLUTION) CHLORINE NITRIC ACID	×××	×××	××	
2874	CHLORINE	×		Х	
2875 2875 2875 2875 2875 2875 2875	AMMONIA AMMONIUM NITRATE (SOLUTION) MANGANESE COMPOUNDS PHOSPHORIC ACID SULFURIC ACID SULFURIC ACID ZINC COMPOUNDS	×××××	*****	*****	
2875	PHOSPHORIC ACID	Х		×	
2879	CHLORINE	X	×	X	
2891	DIMETHYL PHTHALATE	x	×		
Г = X	reported				

07/22 Page	/93 TRI Facilities v 9 1987 to 1990	vith Discr	epancies	in Reporting Status Over Time:
SIC	Chemical	Report11 1987 190	ng status 88 1989	in: 1990
2891	DI-N-BUTYL PHTHALATE	x	×	
2891	FORMALDEHYDE	×	×	
2893 2893 2893 2893 2893	ACETONE BARIUM COMPOUNDS COPPER COMPOUNDS LEAD COMPOUNDS	××××	××××	××
2893	ZINC	×		Х
2893	TOLUENE	~	~	Х
2899 2899 2899 2899	n-BUTYL ALCOHOL METHANOL PHOSPHORIC ACID SULFURIC ACID	××××	××	××××
2899 2899 2899 2899 2899	CRESOL(S) ETHYLENE GLYCOL METHYLENE CHLORIDE TETRACHLOROETHYLENE 1,1,1-TRICHLOROETHANE	****	****	****
2899	FORMALDEHYDE	×	×	×
3021	ZINC COMPOUNDS	×	ž	Х
3052 3052	TOLUENE 1,1,1-TRICHLOROETHANE	××	××	X
3052	METHYL ETHYL KETONE (MEK)	×	~	Х
3069	<b>BIS (2-ЕТНҮЦНЕХҮL) РНТНАLATE</b>	×	×	Х
3069	1, 1, 1-TRICHLOROETHANE	×	×	Х
3069	LEAD COMPOUNDS	x	×	Х
3069	1,1,1-TRICHLOROETHANE	×	×	×
3069 3069 3069 3069	DI-N-OCTYL PHTHALATE METHYL ETHYL KETONE (MEK) TOLUENE ZINC COMPOUNDS	××××	××××	××××
3079	STYRENE	x	х	Х
3079	METHYLENE CHLORIDE	×	x	х
3079	METHYL ETHYL KETONE (MEK)	x	x	х
3079	BIS (2-ЕТНҮЬНЕХҮЬ) РНТНАЬАТЕ	×	×	х
3079	SULFURIC ACID	x		Х
3079 3079	TOLUENE DIISOCYANATE TOLUENE DIISOCYANATE	××	××	

X = reported

07/22	/93 TRI Facilities wit 10 1987 to 1990	ch Discrep	ancles	in Reporting Status O	ver Time:
SIC	Chemical	Reporting 1987 1988	status 1989	in: 1990	
3079	METHYLENE BIS (PHENYLISOCYANATE	X	X	X	•
3079	STYRENE	×	x	X	
3079	1,1,1-TRICHLOROETHANE	×	×	X	
3079	XYLENE	×	×	X	
3087	ACETONE	x		X	
3111	CYCLOHEXANE	x		X	
3143	BIS(2-ETHYLHEXYL) ADIPATE	×		X	
3229 3229 3229 3229	ACETONE CHLOROPHENOLS HYDROFLUORIC ACID SULFURIC ACID	××××	×××	x	
3229 3229	CHLORINE PHOSPHORIC ACID	××	×	x x	
3231	SULFURIC ACID	x x		X	
3245	LEAD COMPOUNDS	x	X	Х	
3251	MANGANESE AND COMPOUNDS	×	х		
3291	TETRACHLOROETHYLENE	x		x	
3293	SULFURIC ACID	x		x	
32955 32955 32955 32995 32995 32995 32995	n-BUTYL ALCOHOL CHLOROBENZENE CYCLOHEXANE ETHYLBENZENE METHYLENE CHLORIDE TETRACHLOROETHYLENE TRICHLOROETHYLENE	× ×× ××××	× ×××	×××	
3298	PHOSPHORIC ACID	×		х	
3312	COPPER COMPOUNDS	x	х	х	
3315	SULFURIC ACID	×	x	х	
3351 3351	COPPER SULFURIC ACID	××	×	×	
3353	CHLORINE	х	×	х	
3354 3354	PHOSPHORIC ACID TOLUENE	××	××	××	
3356	LEAD COMPOUNDS	×	x	х	
3356	HYDROFLUORIC ACID	x		х	
3399	COBALT	×	×	х	
Г = X	reported				

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Over	
Status	
Reporting	;
in	~
Discrepancies	
with	1
TRI Facilities	1987 to 1990

Time:

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07/22/5	53	TRI Facilities wi	th Discrepa	incles 1	n Repo
Page	11	1981 CO 1990	Reporting	status 1989	in: 1990
SIC	chemical				
0000	AMMONITA		x	X	××
6622	COPPER		x	×	< :
34	ACETONE		××	××	××
34 94	FREON 113 METHANOL VVI. FNF		××	×	x
19 C	VIDENT ENE CH.	LORIDE	×	х	
34	MEINILENE		×		x
3411	GLYCOL ETHER	S	: :		>
3412	XYLENE		x		<b>x</b> :
3412 3412	ACETONE METHYL ETHYL	, KETONE (MEK)	××	x	××
3417	FREON 113		×	×	×
3423 3423 3423	ACETONE TOLUENE XYLENE		×××	×××	×××
3423	HYDROCHLORIC	c ACID	×	×	×
3425	HYDROCHLOR I(	C ACID	х	x	×
3429 3429	CHLORINE SULFURIC AC	ID	××	××	
3429	CYANIDE COM	SUNDO	x	X	×
3443	TOLUENE		x		×
3448 3448	n-BUTYL ALC TOLUENE	ОНОГ	××		××
3448	SULFURIC AC	ID	×	х	×
3449 3449	METHANOL XYLENE		××	××	×
3471	SULFURIC AC	CID	×	x	×
3471 3471 3471	HYDROCHLOR: NITRIC ACII SULFURIC AC	IC ACID D CID	×××	×××	×××
3479 3479 3479 3479	COPPER COM HYDROCHLOR SULFURIC A	POUNDS IC ACID CID UNDS	××××	××××	××××
3494	I NITRIC ACI TRICHLOROE	D THYLENE	××	~~	××
349(	6 НҮДКОСНГОК	LIC ACID	x	Ŷ	×

X = reported

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××

××××

07/22/	781 Facilities w	ith Discrep	ancies	in Reporting	Status Over Time:
Page		Reporting	status 1080	in: 1990	
sic	Chemical				
3496	HYDROCHLORIC ACID	X	x	x	
3496 3496	COPPER LEAD COMPOUNDS	××	××	x x	
3498 3498	METHANOL 1, 1, 1-TRICHLOROETHANE	× × × ×		××	
3499	AMMONIA	×	×	Х	
3499 3499	toluene Xy lene	××	××	×	
3499 3499	Toluene Xy lene	××	××	××	
3519 3519 3519 3519	DIETHANOLAMINE METHYL ISOBUTYL KETONE SULFURIC ACID XYLENE	× × ×	×	××××	
3545	AMMONIA	x	×	Х	
3546	AMMONIA	х		X	
3552 3552	SULFURIC ACID 1,1,1-TRICHLOROETHANE	××	×	××	
3562	METHANOL	×	×	Х	
3573 3573 3573 3573 3573 3573	COPPER HYDROCHLORIC ACID LEAD NITRIC ACID SULFURIC ACID SULFURIC ACID 1,1,1-TRICHLOROETHANE	× ××××××	×	*****	
3579	TETRACHLOROETHYLENE	×		x	
3579	METHANOL	×	×		
3600	1, 1, 1-TRICHLOROETHANE	x		×	
3613	COPPER	×	×	X	
3621	XYLENE	х	×	X	
3621 3621 3621	CHROMIUM COMPOUNDS NITRIC ACID PHOSPHORIC ACID	×××	×××	×××	
3622 3622	HYDROCHLORIC ACID SULFURIC ACID	××		××	
3622	COPPER	х	Х	Х	
3624	LEAD	×	х		
3629	SULFURIC ACID	×		х	
= X	reported				

07/22	/93 TRI Facilities wi	th Discr	epancies	in Reporting Status	Over Time:
Page	13 1987 to 1990	Reporti 1987 19	ng status 88 1989	in: 1990	
	Chemica I				
3634 3634	CHROMIUM COMPOUNDS HYDROCHLORIC ACID	××		X X	
3643	TETRACHLOROETHYLENE	x		×	•
3646 3646 3646 3646	CHLORINE NITRIC ACID PHOSPHORIC ACID SULFURIC ACID	××××	××××	××××	
3661 3661	NITRIC ACID SULFURIC ACID	××	××	××	
3674	SULFURIC ACID	x	×		
3674	HYDROFLUORIC ACID	x	х		
3674	SULFURIC ACID	x		х	
3675	XYLENE	х	x	x	
3676	1, 1, 1-TRICHLOROETHANE	×	×	×	
3678	SULFURIC ACID	X	X	х	
3691 3691	ANTIMONY ARSENIC	×	×	x	
3691	LEAD	×		Х	
3692	1,1,1-TRICHLOROETHANE		×	×	
3699 3699 3699 3699	COPPER NICKEL PHOSPHORIC ACID SULFURIC ACID	××××	××××	XXX	
3700 3700	n-BUTYL ALCOHOL XYLENE	××	××	XX	
3700	METHANOL	x	×	X	
3713	TOLUENE	×	×	X	
3714 3714	MANGANESE AND COMPOUNDS ZINC COMPOUNDS	××	x x	××	
3714	1, 1, 1-TRICHLOROETHANE		x	x	
3714	COPPER		x	Х	
3714	METHANOL	×	×	х	
3732 3732	ACETONE STYRENE	××	××	X X	
3732 3732	ACETONE STYRENE		××	××	

3732 ACETONE 3732 STYRENE X = reported

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07/22	781 781 781 781 781 781 793	with Discreps	incies	in Reporting Status Over Time:	
Page	14 1987 to 1990	Reporting	status		
SIC	Chemical	1987 1988	1989	1990	
3842	ETHYLENE OXIDE	×	×	•	
3873	LEAD	x	×	Х	
3964 3964	METHYL ETHYL KETONE (MEK) XYLENE	××	××		
666E	SULFURIC ACID	x	×	Х	
6666 6666 1	CHLORINE CHROMIUM COMPOUNDS HYDROCHLORIC ACID	×××	×× :	× ×:	
3999 5191	NITRIC ACID METHYL BROMIDE	××	××	XX	
0066	ACETONE	X	×	×	
0066	n-BUTYL ALCOHOL ETHYLBENZENE	××	×	××	
0066	FORMALDEHYDE	**	*	××	
0066	METHYL ETHYL KETONE (MEK)	~~>	<×>	~~~~	
0066	METHYLL ISOBUTYL KETONE Toluene	< × × :	< ×:	x x x :	
0066 VAN	XYLENE MANGANESE AND COMPOUNDS	× ×	××	× ×	

X = reported

MANGANESE AND COMPOUNDS

NA

Reported	
Not	
н	
NR	

07/22/5	3 TRI Facilities that S	topped Reporti	ng After Report	Ing significant <b>C</b>	)uantities:
SIC	chemical	Total 1987	l Quantity (lbs) 1988	Reported in: 1989	1990
1479	SODIUM HYDROXIDE	42,000	72,000	NR	NR
20	SODIUM HYDROXIDE	608,098	448,851	NR	NR
20	SODIUM HYDROXIDE	250,000	93,000	NR •	NR
2011	SODIUM HYDROXIDE	63,000	NR	NR	NR
2013	SODIUM HYDROXIDE	59,800	NR	NR	NR
2015	AMMONIA	17,000	17,000	30,300	NR
2024	SODIUM HYDROXIDE	37,191	44,627	NR	NR
2024	SODIUM HYDROXIDE	149,892	148,365	NR	NR
2046 2046	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	7,000,000 280,500	NR 1,200,000	NR NR	NR NR
2052	AMMONIA	NR	NR	53,000	NR
2086	SODIUM HYDROXIDE	73,841	45,674	NR	NR
2086	SODIUM HYDROXIDE	NR	37,363	NR	NR
2111 2111	METHYL ETHYL KETONE (MEK) Toluene	362,182 3,125,065	1,309,883 795,737	1,604,814 945,817	NR NR
2111	AMMONIA	180,883	NR	NR	NR
2111	AMMONIA	173,079	NR	NR	NR
2111	POLYCHLORINATED BIPHENYLS (P	NR	000'06	NR	NR
22	SODIUM HYDROXIDE	470,250	470,000	NR	NR
22 22	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	2,154,800 950,000	NR 757,470	NR NR	NR NR
2200 2200	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	3,000,000 517,000	NR 1,055,000	NR NR	NR NR
2200	ETHYLENE GLYCOL	75,000	NR	NR	NR
2200	SODIUM HYDROXIDE	1,422,000	552,779	NR	NR
2205	AMMONIA	6,619	45,183	46,683	NR
2211	SODIUM SULFATE (SOLUTION)	1,145,300	NR	NR	NR
2211	SODIUM SULFATE (SOLUTION)	4,971,650	NR	NR	NR
2211	SODIUM SULFATE (SOLUTION)	1,155,200	NR	NR	NR
2221 2221	SODIUM HYDROXIDE ETHYLENE GLYCOL	<b>4</b> 90,000 NR	711,000 NR	NR 53,200	NR NR
2221	SODIUM HYDROXIDE	18,250	101,150	NR	NR

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s)	
ODIUM SULFATE	Reported
03	Not
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225	NR

07/22/9 Page	93 TRI Facilities that 2 1987 to 1990	Stopped Reporti	ng After Reporti	ing significant Q	uantities:
SIC	Chemical	Total 1987	Quantity (Ibs) 1988	Reported in: 1989	1990
2221	SODIUM HYDROXIDE	NR	27,500	NR	NR
2221 2221	SODIUM HYDROXIDE SODIUM SULFATE (SOLUTION)	119,500 62,500	125,000 NR	NR NR	NR NR
2221	SODIUM HYDROXIDE	52,994	32,650	NR	NR
2221	SODIUM HYDROXIDE	143,050	115,065	NR	NR
2221	SODIUM SULFATE (SOLUTION)	420,000	NR	NR	NR
2231	SODIUM SULFATE (SOLUTION)	182,210	NR	NR	NR
2251	SODIUM SULFATE (SOLUTION)	105,000	NR	NR	NR
2252 2252	CHLORINE AMMONIUM SULFATE (SOLUTION)	NR 40,000	56,560 29,500	78,000 45,000	NR NR
2252	CHLORINE	171,787	252,670	323,184	NR
2252	CHLORINE	32,928	24,600	29,452	NR
2252 2252	SODIUM SULFATE (SOLUTION) BIPHENYL	854,200 39,560	NR 26,995	NR NR	NR NR
2252	CHLORINE	39,105	33,000	67,732	NR
2252	SODIUM SULFATE (SOLUTION)	60,000	NR	NR	NR
2253	SODIUM HYDROXIDE	111,345	NR	NR	NR
2253 2253	SODIUM HYDROXIDE SULFURIC ACID	NR NR	240,000 120,000	NR NR	NR
2253	SODIUM SULFATE (SOLUTION)	4,388,850	NR	NR	NR
2253	SODIUM SULFATE (SOLUTION)	8,976,700	NR	NR	NR
2257	SODIUM HYDROXIDE	50,700	58,825	NR	NR
2257	SODIUM SULFATE (SOLUTION)	1,508,433	NR	NR	NR
2257 2257	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	930,550 500	NR 47,500	NR NR	NR NR
2257	SODIUM SULFATE (SOLUTION)	2,300,039	NR	NR	NR
2257	SODIUM SULFATE (SOLUTION)	3,506,900	NR	NR	NR
2257	SODIUM SULFATE (SOLUTION)	5,300,200	NR	NR	NR
2257	SODIUM SULFATE (SOLUTION)	20,686,964	NR	NR	NR
2257 2257 2257 2257	GLYCOL ETHERS 1,2-DICHLOROBENZENE AMMONIUM SULFATE (SOLUTION) PSEUDOCUMENE PSEUDOCUMENE	NR NR 76,735 NR	NR NR NR 33, 220	100,000 100,000 NR NR	NN NN NN NN
2257	SODIUM SULFATE (SOLUTION)	332,640	NR	NR	NR

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Page SIC	3 1987 to 1990 Chemical	Total 1987	. Quantity (lbs) 1988	Reported in: 1989	1990
2258 2258	SODIUM HYDROXIDE TETRACHLOROETHYLENE	80,460	49,006 64,471	45,559	NR NR
2258	AMMONIUM SULFATE (SOLUTION)	19,400	19,950	25,850	NR
2258	SODIUM SULFATE (SOLUTION)	137,300	NR	NR	NR
2258 2258 2258	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE PSEUDOCUMENE	234,250 41,122 NR	NR 29, 885 NR	NR NR 28, 381	NR NR NR
2258 2258	AMMONIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	290,000 29,750	152,000 26,600	184,800 NR	NR NR
226	SULFURIC ACID	271,000	271,000	271,000	NR
2260 2260	TETRACHLOROETHYLENE 1,2,4-TRICHLOROBENZENE	NR NR	NR NR	39,650 30,800	NR NR
2261	SODIUM HYDROXIDE	NR	213,338	NR	NR
2261	SODIUM SULFATE (SOLUTION)	6,050,200	NR	NR	NR
2261	SODIUM SULFATE (SOLUTION)	1,000,000	NR	NR	NR
2261	SODIUM HYDROXIDE	167,000	NR	NR	NR
2261 2261	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	759,584 NR	NR 103,062	NR NR	NR NR
2261	SODIUM HYDROXIDE	NR	59,700	NR	NR
2262	AMMONIA	19,342	14,264	25,759	NR
2262	SODIUM SULFATE (SOLUTION)	68,911	NR	NR	NR
2262	SODIUM SULFATE (SOLUTION)	352,306	NR	NR	NR
2262	SODIUM SULFATE (SOLUTION)	970,000	NR	NR	NR
2262	SODIUM HYDROXIDE	75,687	91,397	NR	NR
2262	SODIUM HYDROXIDE	400,000	280,000	NR	NR
2262 2262	SODIUM SULFATE (SOLUTION) 1,1,1-TRICHLOROETHANE	680,000 63,000	NR 48,300	NR 66,000	NR NR
2262 2262	SODIUM HYDROXIDE SODIUM SULFATE (SOLUTION)	150,453 85,500	NR NR	NR NR	NR NR
2269 2269 2269	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE ANMONIUM SULFATE (SOLUTION)	637,910 981,532 45,862	NR 495,286 91,104	NR NR NR	NR NR NR
2269	SODIUM SULFATE (SOLUTION)	940,000	NR	NR	NR
2269	SODIUM HYDROXIDE	360,000	NR	NR	NR

NR = Not Reported

07/22/9 Page	3 TRI Facilities that 4 1990	Stopped Reporti	ng After Report	ing significant Q	uantities:
SIC	Chemical	Total 1987	Quantity (1bs) 1988	Reported in: 1989	1990
2269	SODIUM HYDROXIDE	38,475	34,500	NR	NR
2269 2269	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	880,000 390,000	NR NR	NR NR	NR NR
2269	SODIUM SULFATE (SOLUTION)	150,000	NR	NR	NR
2269 2269 2269	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE 1,2,4-TRICHLOROBENZENE	1,575,000 63,500 NR	NR 335,000 18,560	NR NR 33,000	NR NR NR
2269	SODIUM HYDROXIDE	12,500	36,500	NR	NR
2269	SODIUM HYDROXIDE	NR	290,000	NR	NR
2269	SODIUM HYDROXIDE	109,560	108,365	NR	NR
2269	SODIUM SULFATE (SOLUTION)	264,900	NR	NR	NR
2269	SODIUM HYDROXIDE	NR	26,600	NR	NR
2269 2269 2269	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE TETRACHLOROETHYLENE	259,950 68,700 43,902	NR 82,450 43,375	NR NR 38,625	NR NR NR
2272 2272	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	196,605 105,810	NR 114,908	NR NR	NR NR
2272 2272	AMMONIUM SULFATE (SOLUTION) AMMONIA	NR 155,250	52,650 47,680	NR NR	NR NR
2281	SODIUM SULFATE (SOLUTION)	669,550	NR	NR	NR
2281	SODIUM SULFATE (SOLUTION)	170,000	NR	NR	NR
2281	1,1,1-TRICHLOROETHANE	91,637	NR	NR	NR
2281	SODIUM HYDROXIDE	70,250	150,000	NR	NR
2282	SODIUM HYDROXIDE	34,000	157,971	NR	NR
2283	SODIUM SULFATE (SOLUTION)	360,000	NR	NR	NR
2293	SODIUM HYDROXIDE	1,858,724	1,857,734	NR	NR
2295	SODIUM HYDROXIDE	55,500	NR	NR	NR
2295	XYLENE	31,500	61,000	NR	NR
2297	BIS (2-ETHYLHEXYL) PHTHALATE	NR	39,000	41,800	NR
2298	SODIUM SULFATE (SOLUTION)	548,460	NR	NR	NR
2299	SODIUM SULFATE (SOLUTION)	155,823	NR	NR	NR
2299 2299	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	1,779,750 232,250	NR 348,070	NR NR	NR NR
2300	SODIUM HYDROXIDE	NR	305, 680	NR	NR

NR = Not Reported

07/22/9	33 TRI Facilities th 5 1987 to 1990	at Stopped Report1	ng After Report	ting Significant	Quantities:
SIC	Chemical	Total 1987	Quantity (lbs 1988	) Reported in: 1989	1990
2300	SODIUM HYDROXIDE	201,719	NR	NR	NR
2321 2321	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	620,000 55,000	NR 82,250	NR NR	NR NR
2322	SODIUM SULFATE (SOLUTION)	177,000	NR	NR	NR
2322 2322	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	236,000 500	NR 183,144	NR NR	NR NR
2322 2322 2322	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE BIPHENYL	768,450 500 39,650	NR 170,023 47,778	NR NR NR	NR NR NR
2322	SODIUM HYDROXIDE	NR	86,688	NR	NR
2341 2341	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	97,750 28,809	NR 26,606	NR NR	NR NR
2434	METHANOL	NR	NR	33,010	NR
2499 2499	GLYCOL ETHERS TOLUENE	NR NR	NR NR	52,012 34,018	NR NR
2511 2511	TOLUENE METHANOL	111,046 48,096	53,552 49,529	NR	NR NR
2511	ACETONE	60,312	82,535	93,284	NR
2511	METHANOL	47,234	27,717	NR	NR
2511 2511 2511 2511	TOLUENE METHYL ETHYL KETONE (MEK) n-BUTYL ALCOHOL XYLENE	105,319 61,000 55,314 39,482	100,669 58,252 49,607 38,028	110,643 69,034 57,593 38,573	NR NR NR
2511	ISOPROPYL ALCOHOL	NR	27,018	31,314	NR
2511 2511 2511	TOLUENE n-BUTYL ALCOHOL XYLENE	64,541 50,322 29,678	59,654 56,669 31,595	55,598 54,092 28,200	NR NR NR
2511	METHYL ETHYL KETONE (MEK)	26,079	31,139	36,848	NR
2511 2511	METHANOL TOLUENE	51,843 21,574	41,612 25,478	36,366 25,313	NR NR
2511	FORMALDEHYDE	NR	NR	30,504	NR
2511	XYLENE	14,191	19,839	30,346	NR
2511 2511	METHYL ETHYL KETONE (MEK) Toluene	NR NR	42,404 39,459	NR NR	NR NR
2511 2511	TOLUENE ACETONE	52,250 52,250	NR NR	NR NR	NR NR
2511	TOLUENE	NR	192,895	NR	NR
NR = N	ot Reported				

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07/22/9 Page	3 TRI Facilities that 6 1987 to 1990	Stopped Report	ing After Report	lng significant Q	uantities:
SIC	Chemical	Total 1987	l Quantity (lbs) 1988	Reported in: 1989	1990
2511	BIS (2-ЕТНҮЦНЕХҮL) РНТНАLATE	50,391	NR	NR •	NR
2512	METHANOL	NR	NR	27,773	NR
2517	m- XYLENE	NR	NR	44,509	NR
2519	XYLENE	NR	NR	26,000	NR
2521	GLYCOL ETHERS	65,991	NR	NR	NR
2600 2600	SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE	1,628,347 9,163,862	NR 1,125,300	NR NR	NR NR
2610	SODIUM SULFATE (SOLUTION)	60,655,582	NR	NR	NR
2611 2611	SODIUM SULFATE (SOLUTION) ALUMINUM OXIDE	32,820,000 2,950,000	NR NR	NR NR	NR NR
2611	SODIUM SULFATE (SOLUTION)	16,992,200	NR	NR	NR
2621 2621	SULFURIC ACID SODIUM HYDROXIDE	250 250	58,600 42,338	NR	NR NR
2641 2641 2641	SODIUM SULFATE (SOLUTION) ALUMINUM OXIDE SODIUM HYDROXIDE	24,015,000 16,930,000 1,500,000	NR NR 679, 800	NR NR NR	NR NR NR
2641	METHANOL	29,877	38,753	34,943	NR
2657 2657	METHYL ETHYL KETONE (MEK) DI-N-BUTYL PHTHALATE	49,050 NR	50,000 NR	NR 35,387	NN NN
2657	ACETONE	28,800	28,800	32,720	NR
2741	ISOPROPYL ALCOHOL	73,299	NR	NR	NR
2750	GLYCOL ETHERS	NR	NR	28,133	NR
2819	SODIUM HYDROXIDE	46,000	44,000	NR	NR
2819	SODIUM SULFATE (SOLUTION)	102,943	NR	NR	NR
2819	SODIUM SULFATE (SOLUTION)	855,130	NR	NR	NR
2819	ARSENIC COMPOUNDS	NR	550,000	NR	NR
2819 2819	SODIUM SULFATE (SOLUTION) ALUMINUM OXIDE	3,176,200 NR	NR 385,096	NR NR	NR NR
2821	SODIUM SULFATE (SOLUTION)	206,000	NR	NR	NR
2821	FORMALDEHYDE	NR	64,000	NR	NR
2824 2824	SODIUM SULFATE (SOLUTION) TEREPHTHALIC ACID	1,187,245 47,680	NR 3,103,540	NR 78, 650	NR NR
2824 2824	TEREPHTHALIC ACID ALUMINUM OXIDE	172,000 NR	190,000 68,000	170,750 NR	NR NR

NR = Not Reported

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07/22/9	3 TRI Facilities that	Stopped Reporti	ng After Report	ing Significant (	Quantities:
Page	1 T38/ E0 T330	Total	Quantity (lbs)	Reported in:	1
SIC	Chemical	1987	1988	1989	
2833 2833	SODIUM SULFATE (SOLUTION) METHYL TERT-BUTYL ETHER	280,000 56,750	NR NR	NR NR	NR NR
2833	METHANOL	49,000	49,000		NR
2834 2834	SODIUM SULFATE (SOLUTION) FREON 113	369,000 27,800	NR 52,700	NR 66,400	NR NR
2834	SODIUM SULFATE (SOLUTION)	495,730	NR	NR	NR
2834	ISOPROPYL ALCOHOL	81,214	NR	NR	NR
2834 2834	AMMONIUM SULFATE (SOLUTION) n-BUTYL ALCOHOL	359,000 NR	560,210 29,094	489,810 NR	NR NR
2834	SODIUM SULFATE (SOLUTION)	124,800	NR	NR	NR
2843	SODIUM SULFATE (SOLUTION)	419,000	NR	NR	NR
2844 2844	SODIUM HYDROXIDE SULFURIC ACID	NR NR	71,525 25,517	NR NR	NR NR
2851 2851 2851 2851	METHYL ETHYL KETONE (MEK) ACETONE TOLUENE XYLENE	152,702 107,500 97,600 33,800	97,100 69,400 63,000 51,700	NR NR NR	NR NR NR
2851	TOLUENE	50,366	20,847	41,658	NR
2861	SODIUM SULFATE (SOLUTION)	7,300,000	NR	NR	NR
2865	SODIUM SULFATE (SOLUTION)	3,814,250	NR	NR	NR
2869	TEREPHTHALIC ACID	793,750	379,469	517,910	NR
2869	SODIUM SULFATE (SOLUTION)	2,055,470	NR	NR	NR
2869	AMMONIUM SULFATE (SOLUTION)	NR	121,533	59,298	NR
2869	ALUMINUM OXIDE	NR	83,000	NR	NR
2869	SODIUM SULFATE (SOLUTION)	137,500	NR	NR	NR
2874	SODIUM SULFATE (SOLUTION)	3,165,000	NR	NR	NR
2892 2892 2892	TETRACHLOROETHYLENE CYCLOHEXANE METHANOL	NR 192,000 40,850	411,081 105,278 52,205	N NR NR NR	NR NR NR
2893	CRESOL(S)	67,621	NR	NR	NR
2952	ALUMINUM OXIDE	87,201	32,080	NR	NR
3052 3052	TOLUENE SODIUM SULFATE (SOLUTION)	1,651,600 230,000	1,224,500 NR	377,237 NR	NR NR
3079 3079	TOLUENE ACETONE	23,493 25,734	28,637 23,191	NR 27,471	11R 11R

NR = Not Reported

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1/22/9	<pre>33 TRI Facilities that 8 1987 to 1990</pre>	Stopped Report	ing After Report	ing significant (	Quantities:
0	Chemical	1987	L QUANCICY (103) 1988	1989	1990
	SODIUM SULFATE (SOLUTION)	1,400,000	NR	NR .	NR
79	METHYLENE CHLORIDE	84,000	82,200	NR	NR
79	METHYLENE CHLORIDE	NR	15,694	32,620	NR
79	ISOPROPYL ALCOHOL	NR	NR	37,184	NR
79	ACETONE	12,528	173,922	NR	NR
86	1, 1, 1-TRICHLOROETHANE	NR	63,576	NR	NR
===	XYLENE m-XYLENE n-BUTYL ALCOHOL	54,337 NR 22,053	59,392 NR 22,950	NR 56, 385 51, 945	NR NR NR
31	SODIUM HYDROXIDE	166,000	900'06	NR	NR
45	ALUMINUM OXIDE	NR	2,198	32,376	NR
64	1,1,1-TRICHLOROETHANE	86,200	48,231	NR	NR
64	ALUMINUM OXIDE	147,400	60,000	60,750	NR
6	ASBESTOS	140,500	NR	NR	NR
6 6 6 0 0 0	SODIUM HYDROXIDE SULFURIC ACID HYDROCHLORIC ACID	160,000 130,000 120,000	NR NR NR	NR NR NR	NR NR NR
96	ALUMINUM OXIDE	NR	2,120,985	1,990,190	NR
97	ALUMINUM OXIDE	73,751	84,000	NR	NR
12	ALUMINUM OXIDE	1,350	305,655	266,070	NR
51	ALUMINUM OXIDE	NR	NR	122,750	NR
53	ALUMINUM OXIDE	163,799	182,918	NR	NR
53	ALUMINUM OXIDE FREON 113	836,050 NR	797,166 43,975	NR NR	NR NR
54	SODIUM HYDROXIDE	S	148,000	NR	NR
56	SODIUM SULFATE (SOLUTION) POLYCHLORINATED BIPHENYLS (P	170,000 7, <b>4</b> 00	NR 109,974	NR NR	NR NR
56	ALUMINUM OXIDE	180,620	513,640	NR	NR
161	ALUMINUM (FUME OR DUST)	NR	52	30,000	NR
161	ALUMINUM (FUME OR DUST)	180,000	180,000	NR	NR
862 862 862	COPPER 1,1,1-TRICHLOROETHANE PHENOL	580 93,250 80,086	122,000 76,000 74,081	123,500 90,750 74,829	NR NR NR

07/22/9 Pade	)3 TRI Facilities that Sto 9 1987 to 1990	pped Reporting	After Reporting	g Significant Ç	Quantities:
SIC	Chemical	Total Qu 1987	antity (lbs) Re 1988	eported in: 1989	1990
3411	CYCLOHEXANE	79,000	79,400	55,800	NR
3412	SODIUM HYDROXIDE	58,500	NR	NR	NR
3423	SODIUM HYDROXIDE	352,000	301,260	NR •	NR
3429 3429 3429	SODIUM SULFATE (SOLUTION) TETRACHLOROETHYLENE CVANIDE COMPOUNDS	110,000 38,500 800	NR 37,750 44,050	NR 37,750 28,700	NR NR NR
3429 3429 3429 3429 3429	COPPER SODIUM SULFATE (SOLUTION) CHROMIUM ZINC TETRACHLOROETHYLENE	205,900 205,900 NR NR 138,300	834,852 NR 115,940 106,399 59,772	n n n n n	N N N N N N N N N N N N N N N N N N N
3429 3429	BARIUM ALUMINUM OXIDE	252,340 NR	NR 68,530	NR NR	NR NR
3441	MANGANESE AND COMPOUNDS	NR	NR	331,344	NR
3441	MANGANESE AND COMPOUNDS	NR	NR	54,624	NR
3471	SODIUM SULFATE (SOLUTION)	57,000	NR	NR	NR
3471	SODIUM HYDROXIDE	48,020	31,452	NR	NR
3479	SODIUM HYDROXIDE	115,120	NR	NR	NR
3479 3479	SODIUM SULFATE (SOLUTION) ZINC	115,000 14,750	NR 107,900	NR NR	NR NR
3479	ALUMINUM OXIDE	29,920	34,614	NR	NR
3494	TR I CHLOROETHY LENE	33,058	28,174	29,795	NR
3499	SODIUM HYDROXIDE	21,875	33,191	NR	NR
3499	NA PHTHALENE	89,346	NR	NR	NR
3519	ALUMINUM (FUME OR DUST)	224,800	181,104	153,150	NR
3519	ALUMINUM (FUME OR DUST)	NR	46,500	48,000	NR
3531	DIETHANOLAMINE	NR	30,607	NR	NR
3542	1, 1, 1-TR ICHLOROETHANE	119,006	95,010	74,391	NR
3546 3546 3546	ALUMINUM (FUME OR DUST) Toluene Xylene	NR 56,298 51,750	NR 47,600 30,300	197, 725 Nr Nr	NR NR NR
3552	SODIUM HYDROXIDE	91,440	NR	NR	NR
3562	ALUMINUM OXIDE	NR	62,500	NR	NR
3562	ALUMINUM OXIDE	180,000	200,000	NR	NR
3585	1,1,1-TRICHLOROETHANE	322,720	182,952	308,112	NR
NR = N	ot Reported				

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07/22/9	3 TRI Facilities that :	stopped Reportin	ng After Report	ing significant (	Quantities:
Page I	U IVER CO IVEC	Total 1987	Quantity (lbs) 1988	Reported in: 1989	1990
					*
3599	ALUMINUM OXIDE	78,000	38,000	NR •	NR
3621	XYLENE	NR	NR	26,126	NR
3621	ALUMINUM OXIDE	NR	27,900	NR	NR
3622 3622	SODIUM HYDROXIDE FREON 113	26,304 35,312	39,839 32,604	NR NR	NR NR
3624 3624	COPPER METHANOL	92,418 Nr	121, 620 NR	40,500 27,250	NR NR
3634	METHYL ETHYL KETONE (MEK)	54,224	NR	NR	NR
3634	SODIUM HYDROXIDE	13,961	38,725	NR	NR
3645	1, 1, 1-TRICHLOROETHANE	45,600	15,015	26,697	NR
3661	TETRACHLOROETHYLENE	59,000	NR	NR	NR
3661 3661 3661	FREON 113 TETTRACHLOROETHYLLENE ETHYLENE GLYCOL MONOETHYL ET	53,200 318,300 38,200	87,700 341,000 37,300	80, 800 80, 200 NR	NR NR NR
3674	1, 1, 1-TRICHLOROETHANE	127,398	273,194	107,063	NR
3674 3674	HYDROCHLORIC ACID SODIUM HYDROXIDE	86,254 148,755	67,913 31,336	NR NR	NR NR
3675 3675	BARIUM FREON 113	101,078 86,791	159,019 77,562	35,000	NR NR
3675	ALUMINUM OXIDE	183,000	300,000	NR	NR
3675	ALUMINUM OXIDE	NR	NR	71,541	NR
3676 3676	METHOXYCHLOR METHYLENE CHLORIDE	NR 47,000	159,800 45,300	NR 26,750	NR NR
3678	SODIUM SULFATE (SOLUTION)	98,000	NR	NR	NR
3691	SODIUM SULFATE (SOLUTION)	163,000	NR	NR	NR
3695	METHYLENE CHLORIDE	NR	13,167	26,986	NR
3700 3700	SODIUM SULFATE (SOLUTION) TOLUENE	183,251 432,092	NR 91,272	NR 76,1 <b>4</b> 8	NR NR
3713	TOLUENE	3,900	25,800	30,950	NR
3714	1,1,1-TRICHLOROETHANE	422,300	NR	NR	NR
3714 3714	SODIUM HYDROXIDE TETRACHLOROETHYLENE	183,750 52,238	433,630 NR	NR NR	NR NR
3714	SODIUM HYDROXIDE	150,500	228,000	NR	NR
3714	SULFURIC ACID	92,140	90,200	NR	NR
NR = N	lot Reported				

Stopped Reporting After Reporting Significant Quantities: ÷. the .11111

07/22/	793 TRI Facilities t	lat Stopped Report	ing After Report	ing significant	Quantities:
Page SIC	li I987 CO 1990 Chemical	Tota 1987	l Quantity (lbs) 1988	Reported in: 1989	1990
3714	SODIUM HYDROXIDE	83,640	83,640	NR	NR
3714	ISOPROPYL ALCOHOL	NR	NR	40,200	NR
3714 3714	STYRENE ACETONE	96,970 83,400	50,299 39,200	NR NR	NR NR
3714 3714	ACETONE STYRENE	22,600 23,710	444,400 55,101	NR NR	NR NR
3714 3714	TETRACHLOROETHYLENE METHYLENE CHLORIDE	95,000 12,900	NR 37,651	NR 71,550	NR NR
3732	DIMETHYL PHTHALATE	60,808	NR	NR	NR
3732 3732	DIMETHYL PHTHALATE STYRENE	134,136 72,000	NR 32,736	NR 26,00 <b>4</b>	NR NR
3732	DIMETHYL PHTHALATE	78,693	NR .	NR	NR
3732	ACETONE	66,145	37,610	30,040	NR
3732 3732	STYRENE ACETONE	72,133 71,840	NR NR	NR NR	NR NR
3843	METHANOL	126,911	NR	NR	NR
3931	TR I CHLOROETHY LENE	266,448	NR	NR	NR
3931 3931	1, 1, 1-TRICHLOROETHANE COPPER	55,308 32,641	78,889 41,852	NR NR	NR NR
3999	ACETONE	51,700	24,600	26,572	NR
6662	COPPER	275,000	NR	NR	NR
4930	ALUMINUM OXIDE	5,140,000	5,196,000	NR	NR
7538	SODIUM HYDROXIDE	NR	80,783	NR	NR
8098 8098	SODIUM SULFATE (SOLUTION) SODTIM HYDROXIDE	501,500 101.500	NR NR	NR NR	NR NR

NR = Not Reported

NR NR

NR 501,500 101,500

SODIUM SULFATE (SOLUTION) SODIUM HYDROXIDE

8008 8098

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RI Facilities With Significant Quantity Discrepancies Between Reporting Years: 1987 to 1990

מספי	4	Total	Quantity (lbs)	Reported in:	1990	
SIC	Chemical	1987	TAR8			
2011	AMMONIA	NR	NR	100 .	6,300	
2015	AMMONIA	NR	NR	12,741	765	
2015	CHLORINE	NR	26	18,750	0	
2086 2086	AMMONIA PHOSPHORIC ACID	NR NR	NR NR	18,781 10,660	10 45	
2111	AMMONIA	551,891	476,710	255,905	250	
2111	1, 1, 1-TRICHLOROETHANE	27,040	20,450	22,561	185	
2221 2221	METHANOL SULFURIC ACID	15,000 280,000	<b>44</b> 0,000 350,000	355,474 16,500	293,311 250	
2221	1, 1, 1-TRICHLOROETHANE	36,000	41,000	21,000	0	
2251	BUTYL BENZYL PHTHALATE	NR	NR	17	5,400	
2253 2253 2253	BIPHENYL CHLORINE XYLENE	15,883 73,000 97,744	15,883 73,000 97,744	540 2,700 2,160	000	
2257	XYLENE	5,706	469	NR	NR	
2258	SULFURIC ACID	NR	NR	48,183	0	
2261	AMMONIA	500	21,473	22,043	28,539	
2261	AMMONIA	NR	18,085	14,253	154,703	
2262	XYLENE	10,876	117,995	174,784	174,784	
2262	TETRACHLOROETHYLENE	NR	1,300	970	10,500	
2262	AMMONIA	NR	NR	750	14,250	
2269	METHANOL	NR	NR	13,100	119,005	
2282	SULFURIC ACID	12,100	20,564	12,000	985	
2282	METHANOL	36,108	26,400	10,400	0	
2293	1,1,1-TRICHLOROETHANE	NR	NR	32,267	0	
2300	SODIUM HYDROXIDE	189,000	178	NR	NR	
2321	PHOSPHORIC ACID	NR	NR	46,640	0	
2389	ACETONE	NR	NR	16,020	1,620	
2434	TOLUENE	13,154	154,544	130,750	131,029	
2491 2491 2491	ARSENIC COMPOUNDS CHROMIUM COMPOUNDS COPPER COMPOUNDS	250 NR 250	86 101 58	10,750 10,750 10,750	22,710 10,000 10,000	
2491	NAPHTHALENE	22,724	2,470	807	401	

NR = Not Reported

1990	
to	
1987	
Years:	
Reporting	
Between	
Discrepancies	
Quantity	
Significant	
With	
Facilities	
TRI	
1/22/93	•
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Page	2	Total	Quantity (]bs)	Reported in:	
SIC	Chemical	1987	1988	1989	1990
2491 2491 2491	ANTHRACENE CHROMIUM COMPOUNDS COPPER COMPOUNDS NA PHTHALFNF	4 NR NR 1.152	30,000 1,900	4,055 1,100 820	NR 11,978 9,209 NR
2491	PENTACHLOROPHENOL	1	58	15,302	31, 830
2491 2491	CHROMIUM COMPOUNDS COPPER COMPOUNDS	1,187,625 7,036	1,864 1,065	3,884 2,541	00
2491	COPPER COMPOUNDS	NR	3,351	311	220
2492	FORMALDEHYDE	97,690	22,769	283,748	263,100
2493	FORMALDEHYDE	NR	NR	3,500	78,848
2511	METHYL ETHYL KETONE (MEK)	NR	3,585	361	NR
2511	n-BUTYL ALCOHOL	67,365	6,494	NR	NR
2511	METHYL ETHYL KETONE (MEK)	NR	19,209	31,664	0
2511	n-BUTYL ALCOHOL	32,872	35,070	28,215	0
2511	XYLENE	14,750	16,800	19,550	0
2511	ACETONE	NR	NR	12,100	0
2511 2511 2511 2511 2511	n-BUTYL ALCOHOL METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	14,787 17,978 55,932 35,930 12,705	18,031 20,065 103,008 46,169 14,792	21,490 24,637 55,588 44,976 16,857	00000
2511 2511 2511 2511	ACETONE n-BUTYL ALCOHOL TOLUENE XYLENE	NR NR 52,560 17,734	17,059 NR 63,261 20,789	11,245 15,720 54,000 14,746	0000
2511 2511 2511 2511 2511 2511 2511	ACETONE BIS (2-ETHYLHEXYL) PHTHALATE GLYCOL ETHERS METHYL ETHYL KETONE (MEK) METHANOL TOLUENE XYLENE	17,020 12,617 36,416 NR 155,016 85,423 35,667	46,907 13,528 13,528 49,759 166,9759 98,350 47,926	42,545 15,480 15,480 139,785 91,000 48,020	
2600	HYDROCHLORIC ACID	71,599	520	650	22,000
2611	HYDROCHLORIC ACID	37,000	1,400	1,400	1,300
2611	AMMONIA	6,050	620	1,350	725
2641 2641	AMMONIA CATECHOL	490,000 NR	8, 301 NR	8,850 1,510	340,280 17,600
2657	TOLUENE	153,900	12,300	28,800	30,300
2672	FORMALDEHYDE	002'6	69,011	1,926	NR
NR = NC	ot Reported				

TRI Facilities With Significant Quantity Discrepancies Between Reporting Years: 1987 to 1990

07/22/93 Pare 3

SEC         Chenkleal         1997         1996         1990         1991	טת	D	Total	Quantity (1bs)	Reported in:	
732         TEFRACHLORGETHYLENE         NR         NR         13         11, 211           2813         NETMANGL         8,1340         800         4,500         14,660         783         10,687           2813         NETMANGL         8,1340         800         4,500         14,600         780         14,650           2813         NETMANGL         13,960         13,960         12,000         750         1,0087           2813         NETMANGL         14,500         14,500         730         1,0087         2           2811         FERMANGL         NR         12,510         4,002         67         2         2           2821         RENTANGL         NR         12,510         14,500         11,750         15,5514         3,819           2823         FENTLALENE GLYCOL         27,400         8,510         14,600         11,750         15,5514           2824         FENTLANE         14,500         1,700         11,750         15,525         551           2824         FENTLALINE GLYCOL         27,400         8,510         11,750         15,525         551           2824         FENTLALACID         25,550         14,000         11,750         <	SIC	Chemical	1987	1988	1989	1990
2819         MEYNANGL         8,140         8,0         4,500         14,650         4,500         14,650         14,650         14,650         14,650         14,650         14,650         15,60         10,667         10,667           2813         WURDECHLORIC ACID         13,500         13,600         750         1,000         750         1,000           2813         SULEURAL CACID         14,500         14,500         20,000         750         1,000           2821         FORMALDENTOE         14,500         14,500         14,000         11,750         15,255           2821         FNETLENE CLID         21,400         8,500         11,500         15,255           2823         FNETLENE CLID         21,400         8,500         11,500         15,255           2824         FNETLENE CLID         21,400         11,500         15,255         15,121         15,255           2824 <td>152</td> <td><b>TETRACHLOROETHY LENE</b></td> <td>NR</td> <td>NR</td> <td>38</td> <td>13,211</td>	152	<b>TETRACHLOROETHY LENE</b>	NR	NR	38	13,211
2819         ANGNLA         NR         668         783         10,667           2813         HITSRCENLORIC ACID         13,560         12,000         756         1,000           2813         SULFURIC ACID         14,500         14,500         759         1,000           2821         FORMULENER         NR         14,500         750         1,000         750         1,000           2821         FORMULENER         NR         12,000         750         1,000         15,514         2           2821         BARUN CONDONDS         1,2500         34,000         11,750         15,255         500           2821         EPHVLENE GLYCOL         24,400         8,500         14,000         15,262         500           2823         PHORENCENDE         24,400         8,500         14,000         15,262         500           2824         PHYLENE GLUCOL         25,500         14,000         15,760         15,756         500           2824         PHYLENE GLUCOL         25,500         14,000         15,700         15,750         25,468           2824         PHYLENE GLUCOL         24,400         8,500         14,000         15,700         10,1111	619	METHANOL	8,340	800	4,500	14,850
2313         HYDROCHLORIC ACID         13,550         22,000         750         1,000           2813         SULFURIC ACID         1,500         2,000         750         1,000           2811         SULFURIC ACID         1,500         2,000         750         1,000           2821         FORMULENTOE         24,500         4,002         67         2           2821         BRTUM CONFORUDS         1,250         34,000         1,750         15,255           2821         ETENLEDE GLYCOL         24,400         8,500         14,000         1,500         500           2824         POSTHORIC ACID         25,500         14,000         1,500         50         500           2824         PENTULANCOROL         24,400         8,500         14,000         1,500         500           2824         PENTULANCOROL         24,400         8,500         14,000         1,500         55,500         10,111           2824         PENTULENE CHLORIDE         2,593         2,214         21,230         2,214           2831         PENTUL         ALCOL         14,000         1,500         10,111           2841         POLORIDE         2,500         14,000         1,500 <td>819</td> <td>AMMONIA</td> <td>NR</td> <td>668</td> <td>783</td> <td>10,867</td>	819	AMMONIA	NR	668	783	10,867
2819         GULFURIC ACID         14,500         4,002         67         2           2821         FORMALDERIDE         20,400         285,111         239,392         115,514         3,819           2821         METHANOL         NR         1,250         34,000         11,750         15,255           2821         ETHYLENE GLYCOL         24,400         8,500         1,500         500           2824         PHORANCHANCL         25,500         14,000         1,500         500           2824         PHETHALCANELOL         27,400         8,500         15,262         500           2824         PHETHALCANCHOL         25,500         14,000         1,500         500         2,214           2824         PHETHALCANTERE CHLORIDE         2,058         314,657         427,700         50,222         2,214           2834         PHULENE CHLORIDE         2,058         314,657         437,700         10,110           2834         PHETHALCANCHAL         18,950         16,200         10,110         25,400         10,110           2834         PHETHALLENE CLUCIL         14,857         414,657         429,7700         10,110           2844         PHOLIC         11,1-THICHLOROCTHANE	2819 2819	HYDROCHLORIC ACID NITRIC ACID	13,950 2,000	12,000 20,000	750 750	1,000 1,000
2211         FORMALDENTDE         20,400         285,131         239,392         115,514           2221         BATUN CONFOUNDS         1,250         34,000         11,750         15,255           2221         ETHYLENE GLYCOL         24,400         8,500         15,262         500           2231         FINYLENE GLYCOL         24,400         8,500         15,262         500           2323         FUNTULALCONDL         2,550         14,000         1,500         500           2323         PUNTULALCONDL         2,550         14,000         1,500         500           2323         PUNTULALCONDL         2,500         14,000         1,500         500           2323         PUNTULANE CHLORIDE         2,593         144,657         297,300         25,406           2333         PUNULANE         CHLORIDE         2,593         144,657         297,300         15,600         10,110           2334         PUNULANE         CHLORIDE         2,593         144,657         297,300         15,600         10,110           2333         PUNULANE         CHLORIDE         1,0,000         1,700         15,600         10,110           2343         JUNULANE         JUNULANE <t< td=""><td>2819</td><td>SULFURIC ACID</td><td>14,500</td><td>4,002</td><td>67</td><td>2</td></t<>	2819	SULFURIC ACID	14,500	4,002	67	2
2821         NETHANOL         NR         120         118         3,919           2821         BARTUM COMPONDS         1,250         34,000         11,750         15,255           2821         FTHYLENE GLYCOL         24,400         8,500         15,262         500           2823         FTHYLENE GLYCOL         24,400         8,500         15,262         500           2824         FHORTE ACID         25,500         14,000         1,500         500           2823         FTHYLENE GLYCOL         2,5978         1412         1716         97,20           2823         FTHYLENE CHLORIDE         2,5978         1412         171,500         15,200           2833         FTHYLENE CHLORIDE         2,5978         141,500         10,110           2834         FTHYLENE CHLORIDE         2,9958         14,557         29,550         27,240           2843         FTHYLENE GLYCOL         NR         100,000         1,500         10,110           2843         AMONIA         16,900         1,500         17,000         175,000           2843         MONIA         16,900         1,500         17,000         175,000           2843         MONIA         16,900	2821	FORMALDEHYDE	20,400	285,131	239,392	135,514
2821         BARTUM COMPONDS         1,250         34,000         11,750         15,255           2824         PHOSFHORTC ACID         24,400         8,500         15,262         500           2824         PHOSFHORTC ACID         25,500         14,000         1,500         500           2824         PHOTVLANC ANHYDRIDE         2,900         14,000         1,500         500           2824         PHTVLLIC ANHYDRIDE         2,900         2,112         10,120         9,320           2824         PETMYLLIC ANHYDRIDE         2,978         812,010         427,700         10,111           2824         PETMILIC ANHYDRIDE         3,058         144,657         427,700         188,155           2824         PETMILIC ANHYDRIDE         1,1,1-TRICHLORDETTANE         2,958         144,657         427,700         188,155           2834         PETMILE         1,1,1-TRICHLORDETTANE         5,000         1,1,1         25,405           2843         AMONIA         168,959         14,550         10,116         175,600           2843         TETRACHLORDETTANE         5,800         1,1,1         25,405         25,405           2843         AMONIA         1,1,1         1,1,1         2,146	2821	METHANOL	NR	120	118	3,819
2821         ETHYLENE GLYCOL         24,400         8,500         15,262         500           2824         PHOSFHORIC ACID         25,500         14,000         1,500         500           2824         P-BUTYL ALCOHOL         2,500         14,000         1,500         500           2824         P-BUTYL ALCOHOL         2,000         2,112         171,500         500           2824         PETHYLENE CHLORIDE         2,058         144,657         29,350         3,224           2824         METHYLENE CHLORIDE         3,058         144,657         21,283         3,2405           2834         FETHYLENE GLYCOL         118,950         16,200         15,600         10,110           2814         ETHYLENE GLYCOL         16,000         15,200         10,110         175,600           2813         PAMONIA         1,1,1-THATCHLOROETHANE         5,00         14,520         13,160         175,600           2813         TETRACHLOROETHANE         5,00         1,1,160         175,600         10,110           2813         TETRACHLOROETHANE         5,00         1,1,160         175,600         10,110           2813         TETRACHLOROETHANE         5,00         1,1,160         175,600 <t< td=""><td>2821</td><td>BARIUM COMPOUNDS</td><td>1,250</td><td>34,000</td><td>11,750</td><td>15,255</td></t<>	2821	BARIUM COMPOUNDS	1,250	34,000	11,750	15,255
2824         PHOSFHORIC ACID         25,500         14,000         1,500         500           2824         T-BUTYL ALCOHOL         2,000         2,110         19,120         9,320           2824         T-BUTYL ALCOHOL         2,000         2,000         2,110         19,120         9,320           2824         T-BUTYL ALCOHOL         2,070         2,000         2,110         110,120         9,320           2824         METHULIC ANHYDRIDE         2,978         912,010         27,100         10,110           2824         METHUALLINE CHLORIDE         2,978         912,010         10,110         25,405           2824         METHULINE CHLORIDE         2,978         144,657         99,350         10,110           2813         PMONIA         NR         100,000         76,000         175,600         175,600           2813         MENUL         RETRICHORDETHANE         5,000         1,1,1-TRICHLORDETHANE         5,000         1,75,600         175,600           2813         MONIA         1,1,1-TRICHLORDETHANE         5,000         1,740         175,600         175,600           2814         ADMONIA         1,1,1-TRICHLORDETHANE         5,000         1,114         20,25         21 <td>2821</td> <td>ETHYLENE GLYCOL</td> <td>24,400</td> <td>8,500</td> <td>15,262</td> <td>500</td>	2821	ETHYLENE GLYCOL	24,400	8,500	15,262	500
2824         n=BUTVL ALCOHOL         2,000         2,110         10,120         8,320         2,321         3,222         3,222         3,222         3,222         3,222         3,222         3,222         3,223         3,224         3,254         3,254         3,254         3,110         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         175,600         10,110         10,120         10,120         10,120         10,120         10,120         10,120         10,120         10,120         10,120         10,120         12,121         10,120         10,120	2824	PHOSPHORIC ACID	25,500	14,000	1,500	500
2824         METHYLENE CHLORIDE         3,058         144,657         99,350         25,405           2833         FVRIDINE         TOLUENE         2,978         812,010         427,700         186,150           2834         FWRIDINE         GLUCENE         18,950         16,200         15,800         10,110           2834         ETHYLENE GLYCOL         NR         100,000         76,000         14           2834         AMONIA         10,11-TRICHLORDETHANE         500         5,500         6,000         196,160           2843         TETRACHLORDETHANE         5,800         1,748         8,025         21           2843         AMONIA         31,014         29         5500         6,000         1,95,600           2843         AMONIA         31,014         29         5,500         6,000         1,75,600           2843         AMONIA         31,014         29         5,500         6,000         1,75,600           2843         AMONIA         31,014         20         30,25         21         41           2841         METHAUL         AMONIA         31,014         20         22         23           2841         METHAUL         AMONIA	2824 2824 2824	n-BUTYL ALCOHOL 2-METHOXYETHANOL PHTHALIC ANHYDRIDE	80 2,200 NR	710 2,412 NR	10,120 173,690 21,283	8,320 35,223 2,214
2833         PKRIDINE         418,950         16,200         15,800         10,110           2834         ETHYLENE GLYCOL         NR         100,000         76,000         14           2834         AMGONIA         10,110         75,000         13,160         175,600         14           2843         1,1,1-TRICHLORDETHANE         5,00         5,500         6,000         1,980         23           2843         TETRACHLORDETHYLENE         5,800         1,748         8,025         21           2843         AMOONIA         31,014         30         31,014         30         33         31           2843         AMOONIA         31,014         30         1,748         8,025         21         415           2843         SULFURIC ACID         31,014         30         33         32         33         31         415           2851         METHYL FETONE         5,800         1,148         8,025         33         35         415           2851         METHYL FETONE         1,014         302         32         33         32         35           2851         TOUDENE         TOUDENE         25,49         368,488         302         35	2824 2824	METHYLENE CHLORIDE TOLUENE	3,058 2,978	144,657 812,010	99,350 427,700	25,405 188,150
2834       ETHYLENE GLYCOL       NR       100,000       76,000       14         2834       AMONIA       5,500       5,000       13,160       175,600         2843       1,1,1-TRICHLOROETHANE       5,000       1,14       8,025       21         2843       TETRACHLOROETHALE       5,800       1,748       8,025       21         2843       TETRACHLOROETHALE       5,800       1,748       8,025       21         2843       AMONIA       31,014       30       33       31         2843       AMONIA       31,014       30       33       31         2843       AMONIA       31,014       30       33       31         2843       AMONIA       31,014       30       35       7,419         2843       AMONIA       31,014       30       33       37       37         2851       METHYL ETHYL KETONE       1,746       3,25       7,419       36       36         2851       METHYL ETHYL KETONE       MEK       23,56       1,466       5,755       1,466         2851       TOLUBNE       TOLUBNE       25,146       32,765       1,466       2,756         2861       TOLUBNE<	2833	PYRIDINE	418,950	16,200	15,800	10,110
2834         AMONIA         168,959         14,520         13,160         175,600           2843         1,1,1-TRICHLOROETHANE         500         5,500         6,000         1,880           2843         TETRACHLOROETHYLENE         5,800         1,748         8,025         21           2843         TETRACHLOROETHYLENE         5,800         1,748         8,025         21           2843         AMMONIA         31,014         30         37         31         33         31           2843         AMMONIA         31,014         3,549         368,488         8,025         21         31           2851         ACETONE         31,014         30         72,864         59,713         1,405           2851         METHYL ETHYL KETONE         88,292         221,565         181,577         33           2851         TOLUENE         25,146         72,864         302,725         1,405           2851         TOLUENE         25,809         1,548         3,506         1,405           2851         TOLUENE         25,809         1,648         2,351         3,26           2851         TOLUENE         NR         NR         18,758         2,36	2834	ETHYLENE GLYCOL	NR	100,000	76,000	14
2843       1,1,1-TRICHLOROETHANE       5,00       5,500       6,000       1,748       8,025       21         2843       TETRACHLOROETHYLENE       5,800       1,748       8,025       21         2843       AMONIA       5,800       1,748       8,025       21         2843       AMONIA       31,014       299       505       7,415         2843       SULFURIC ACID       31,014       30       33       37         2851       MCHAUL KETONE       1014       299       505       7,415         2851       ACETONE       31,014       23,549       369,488       302,725       1,000         2851       METHYL KETONE       EKK       25,146       72,516       5,713       1,400         2851       TOLUENE       27,065       13,460       2,755       1,400       2,755         2851       TOLUENE       NR       NR       NR       18,758       1,206         2851       TOLUENE       NR       NR       18,758       2,755       2,755         2851       TOLUENE       NR       NR       18,758       1,256         2869       INENE       NR       NR       18,758       1,256	2834	AMMONIA	168,959	14,520	13,160	175,600
2843         TETRACHLOROETHYLENE         5,800         1,748         8,025         21           2843         AMMONIA         31,014         299         505         7,415           2843         SULFURIC ACID         31,014         299         505         7,415           2843         SULFURIC ACID         31,014         302,725         1,000           2851         ACETONE         272,864         59,713         1,000           2851         METHYL ETHYL KETONE         (MEK)         272,864         59,713         1,400           2851         TOLUENE         281,000         277,864         59,713         1,400         2,756           2851         TOLUENE         METHANOL         86,292         221,363         181,577         3,366           2851         TOLUENE         NR         NR         NR         1,348         3,400         2,756           2851         TOLUENE         NR         NR         NR         1,348         3,400         2,756           2869         VILENE         NR         NR         NR         18,758         1,256           2869         DIMETHYL PHTHALATE         594,425         1,425,584         3,167         NR <td>2843</td> <td>1, 1, 1-TRICHLOROETHANE</td> <td>500</td> <td>5,500</td> <td>6,000</td> <td>1,880</td>	2843	1, 1, 1-TRICHLOROETHANE	500	5,500	6,000	1,880
2843       AMMONIA       30       505       7,415         2843       SULFURIC ACID       31,014       30       33       37,415         2851       ACETONE       METHYL ETHYL KETONE (MEK)       3,549       368,488       302,725       1,000         2851       ACETONE       ACETONE       25,146       72,864       59,713       1,450         2851       TOLUENE       272,864       322,725       1,000       27,755       1,000         2851       TOLUENE       88,292       221,363       181,577       3,350       5,351         2851       TOLUENE       92,778       1,686       22,348       3,400       2,755         2851       TOLUENE       NR       NR       NR       1,686       2,755       1,256         2851       TOLUENE       NR       NR       NR       1,686       2,755       1,256       1,256         2851       TOLUENE       S6,909       1,686       2,350       1,256       3,360       1,256         2869       DIMETHYL PHTHALATE       5       5       5,351       3,265       5,361       3,266         2869       METHANOL       284,425       142,584       3,167       N	2843	TETRACHLOROETHYLENE	5,800	1,748	8,025	21
2851ACETONE METHYL ETHYL ETHANOL3,549368,488302,7251,0002851TOLUENE25,1462,70627,06522,3485,515,3102851TOLUENE92,7781,3483,4002,7562851TOLUENE92,7781,3483,4002,7562851TOLUENE02,7781,6862,3501,2572851TOLUENENRNRNR18,758(2851TOLUENE59,7131,6862,3501,2562851TOLUENES65,5555,3501,2562869DIMETHYL PHTHALATE584,425142,5843,167NI2869XYLENE284,425142,5843,167NI2869XYLENE27626544,4495,542869ETHYLENE DICHLORIDE17,33522,800400,474412,82	2843 2843	AMMONIA SULFURIC ACID	NR 31,014	299 30	505 33	7,419 37
2851       TOLUENE       92,778       1,348       3,400       2,750         2851       XYLENE       25,809       1,686       2,350       1,250         2851       TOLUENE       NR       NR       NR       18,758       0         2859       DIMETHYL PHTHALATE       59       5,555       5,351       3,265         2869       METHAL       284,425       142,584       3,167       NI         2869       XYLENE       284,425       142,584       3,167       NI         2869       XYLENE       276       265       44,449       5,54'         2869       XYLENE       DICHLORIDE       17,335       22,800       400,474       412,82	2851 2851 2851 2851	ACETONE METHYL ETHYL KETONE (MEK) METHANOL TOLUENE	3,549 25,146 6,100 88,292	368,488 72,864 27,065 221,363	302,725 59,713 22,348 181,577	1,000 1,450 500 3,350
2851     TOLUENE     NR     18,758     (       2869     DIMETHYL     FHTHALATE     59     5,555     5,351     3,26       2869     METHANOL     284,425     142,584     3,167     N       2869     XYLENE     276     265     44,449     5,54       2869     ETHYLENE     DICHLORIDE     17,335     22,800     400,474     412,82	2851 2851	TOLUENE XYLENE	92,778 25,809	1,348 1,686	3,400 2,350	2,750 1,250
2869         DIMETHYL         PHTHALATE         59         5,555         5,351         3,26           2869         METHANOL         284,425         142,584         3,167         NI           2869         XYLENE         276         265         44,449         5,54           2869         ETHYLENE         17,335         22,800         400,474         412,82	2851	ToLUENE	NR	NR	18,758	0
2869         METHANOL         284,425         142,584         3,167         Ni           2869         XYLENE         276         265         44,449         5,54           2869         ETHYLENE DICHLORIDE         17,335         22,800         400,474         412,82	2869	DIMETHYL PHTHALATE	59	5,555	5,351	3,261
2869         XYLENE         276         265         44,449         5,54           2869         ETHYLENE DICHLORIDE         17,335         22,800         400,474         412,82	2869	METHANOL	284,425	142,584	3,167	NR
2869 ETHYLENE DICHLORIDE 17,335 22,800 400,474 412,82	2869	XYLENE	276	265	44,449	5,540
	2869	ETHYLENE DICHLORIDE	17,335	22,800	400,474	412,820

cs: 1987 to 1990 •

07/22/93	TRI Facilities With Sign	ifficant Quant	ity Discrepanc:	ies Between Repo	orting Years:
SIC	Chemical	Total 1 1987	Quantity (lbs) 1988	Reported in: 1989	1990
2869	XXLENE	21,583	954	630	207
2869 2869 2869	n-BUTYL ALCOHOL DIMETHYL PHTHALATE DI-N-BUTYL PHTHALATE	233 1,700 8,200	22,868 3,082 6,886	11,755 90,712 401 .	39,324 48,890 260
2869	METHANOL	1,250	826,605	563,375	551,086
2873	AMMONIUM NITRATE (SOLUTION)	1,000	270	230	29,080
2875	AMMONIA	250	6,042	637	605
2891	FORMALDEHYDE	6,469	195	2	2
2893 2893	METHANOL TOLUENE	14,100 83,165	1,215 5,525	<b>41</b> 0 1,830	NR 3,558
2893	METHANOL	1,000	250	3,750	NR
2893	METHYL ETHYL KETONE (MEK)	NR	4,014	99,600	147,135
2893	METHYL ETHYL KETONE (MEK)	4,200	29,115	2,975	2,858
2952	ZINC COMPOUNDS	175,279	2,693	NR	NR
3011	ZINC COMPOUNDS	1,630	14,460	8,000	150
3052	ZINC COMPOUNDS	10,150	40,839	35,184	3,500
3069	ACETONE	NR	NR	40,248	499,801
3078 3078	ACETONE STYRENE	NR NR	NR 4,136	12,045 3,283	00
3079	TOLUENE DIISOCYANATE	32,454	32,107	32,750	250
3079	1, 1, 1-TRICHLOROETHANE	106,919	68,415	48,400	4,950
3079 3079	METHYLLENE CHLORIDE METHANOL	NR NR	23,912 NR	19, <b>404</b> 11,750	00
3079 3079	ACETONITRILE HYDROCHLORIC ACID	26,400 200	22,249 79	34,435 79	30 37,463
3079	METHYLENE CHLORIDE	40,200	37,000	34,820	367,660
3079	METHYL ETHYL KETONE (MEK)	24,350	51,689	73,388	0
3081	METHYLENE CHLORIDE	251,500	237,500	18,449	NR
3089	LEAD	NR	6,145	6,895	0
3143	METHYL ETHYL KETONE (MEK)	86	111	1,600	38,300
3245	ZINC COMPOUNDS	8,250	3,102	64,151	56,593
329	PHENOL	1,250	31,000	48,190	71,250
3291	TOLUENE	200,000	186,376	18,000	NR

NR = Not Reported

1990
to
1987
Years:
Reporting
Between
Discrepancies
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TRI

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07/22/93 Page 5	TRI Facilities With Sign
SIC	Chemical
3291	XYLENE
2112	CHROMITIM COMPOUNDS

Page	ſ	Total	Ouantity (lbs)	Reported in:	
SIC	Chemical	1987	1988	1989	1990
3291	XYLENE	NR	47,502	1,000	NR
3312 3312	CHROMIUM COMPOUNDS MANGANESE COMPOUNDS	500 6,300	6,560 107,570	31,065 149,060	31,659 819,848
3351 3351	COPPER COMPOUNDS ZINC COMPOUNDS	467,138 50,052	563,234 47,162	5,271 2,050	NR NR
3353	METHANOL	NR	931	54,556	51,218
3354 3354 3354	NITRIC ACID SULFURIC ACID XYLENE	4 2 N R	4 4 500	14,750 12,750 16,092	500 122,635 500
3356 3356	HYDROFLUORIC ACID NITRIC ACID	32,300 410,500	<b>44,</b> 300 610,000	22,100 301,500	2,100 6,900
3356	CHROMIUM	22,497	16,411	12,411	633
3357	COPPER	250	5,000	6,550	250
3423 3423	HYDROCHLORIC ACID SULFURIC ACID	230,055 163,100	230,055 155,100	2,125 50	2,125 50
3429	ZINC COMPOUNDS	NR	775	8,764	500
3452	HYDROCHLORIC ACID	NR	NR	750	68,255
3471	COPPER	97,218	8	51	51
3471	SULFURIC ACID	NR	14,632	1,500	510
3471	SULFURIC ACID	65,354	36,283	18,513	250
3479	TOLUENE	3,950	533	31,306	11,484
3499 3499	n-BUTYL ALCOHOL GLYCOL ETHERS	NR NR	NR NR	3,000 6,551	00
3542	MANGANESE COMPOUNDS	34	24	7,012	NR
3546 3546	COPPER STYRENE	563,750 118,200	523,100 11,770	754 7,460	NR 12,150
3546	1, 1, 1-TRICHLOROETHANE	161,300	209,050	18,540	NR
3553	ETHYLENE GLYCOL	24,500	30,000	33,500	3,571
3573	AMMONIA	1,750	15,000	1,500	505
3586	XYLENE	16,300	1,600	NR	NR
3621	FREON 113	NR	34,992	46,117	0
3621	NICKEL COMPOUNDS	NR	NR	3,750	270
3622	ALUMINUM OXIDE	500	10,784	9,828	IIR
3622	1, 1, 1-TRICHLOROETHANE	12,000	10,800	1,200	NR
NR =	Not Reported				

Section 10
With Significant Quantity Discrepancies Between Reporting Years: 1987 to 1990 11444 ų L тRТ

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01/22/9 Page	J INI FACILICIES WICH	SIGNILLCAUL VUAIL	THÀ DISCLEMANC		OF CTILG I CATA:
SIC	Chemical	Total 1987	Quantity (lbs) 1988 	Reported in: 1989	1990
3624	ACETONE	81,000	72,000	96,000	0
3634	PHENOL	<i>TTT</i>	79,587	112,743	119,682
3646	ETHYLENE GLYCOL	NR	NR	83,000	0
3674	METHANOL	750	1,565	19,980	NR
3679 3679 3679 3679	COPPER FREON 113 NICKEL 1,1,1-TRICHLOROETHANE	NR 1,850 NR 475	14 26,859 22 443	4,102 68,194 9,710 5,361	NR 24,875 10 16,414
3692	TR I CHLOROETHYLENE	4,529	127,101	140,803	187,918
3692	1,1,1-TRICHLOROETHANE	1,400	13,500	NR	NR
3713 3713	n-BUTYL ALCOHOL GLYCOL ETHERS	105,700 NR	2,865 1,100	124,200 110,000	144,150 129,000
3713	XYLENE	500	14,500	NR	NR
3714	<b>TETRA</b> CHLOROETHY LENE	NR	105,247	304	500
3714	METHANOL	NR	NR	30,242	255
3714	ETHYLENE GLYCOL	NR	16,763	750	NR
3732 3732	STYRENE TOLUENE	222,750 43	20,070 12,612	5,060 6,591	15,270 NR
3825	FREON 113	28,000	27,000	36,000	0
3843	ETHYLENE OXIDE	15,862	17,604	16,950	0
3843	ACETONE	15,319	11,797	3,257	260
3942	ZINC	12,938	17,076	13,990	0
3963 3963 3963	ACETONE ALUMINUM OXIDE METHYLENE CHLORIDE	N N N N N N N N N N N N N N N N N N N	45,000 16,000 53,100	2,400 1,200 5,000	2,400 NR 5,250
3999	STYRENE	136,963	147,200	11,390	NR
6666 6666 9666	COPPER COMPOUNDS METHYL ETHYL KETONE (MEK) NICKEL COMPOUNDS	NR 6,050 4,403	3,000 75,750 7,500	2,250 32,000 2,250	22,700 93,875 22,737
3999	METHYLENE CHLORIDE	67,250	5,307	5,114	7,022
6666 6666 1000	ETHYLBENZENE Toluene Xylene	NR NR NR	10,010 35,751 35,751	13,079 30,977 44,747	000

NR = Not Reported

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