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FINAL SUMMARY OF RECOMMENDATIONS: BILL 167 – THE TOXICS REDUCTION ACT, 2009*

Prepared for:

Standing Committee on General Government

Prepared by:

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May 28, 2009

* Projects prepared by the Legislative Research Service are designed in accordance with the requirements and instructions of the Committee making the request. The views expressed should not be regarded as those of the Legislative Research Service or of the individual preparing the project.

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INTRODUCTION

This Final Summary sets out recommendations and concerns presented to the Standing Committee on General Government during its public hearings on Bill 167, the *Toxics Reduction Act, 2009*, held in Toronto on May 13 and 25, 2009.

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This Final Summary is intended to assist Members in their deliberations on Bill 167. It is not a complete record of all the evidence that the Committee heard, nor is it a comprehensive review of all arguments made by witnesses. Accordingly, comments have been abbreviated and recommendations summarized. For a full account of the evidence presented to the Committee, reference should be made to *Hansard*, and to the written submissions themselves.

Where possible, witnesses' comments have been assigned to the most appropriate provision of the Bill. Concerns or recommendations of witnesses on other matters have been added under appropriate topical headings at the end of the document.

An alphabetical list of the witnesses and the abbreviations used to identify them appears at the end of the document. Where more than one abbreviation appears in connection with a specific recommendation, it indicates that the witnesses identified expressed substantially the same viewpoint, or recommendation.

RECOMMENDATIONS RELATED TO SPECIFIC SECTIONS OF THE BILL

Global Change

In all places where Bill 167 speaks of toxic substances being used or created, the words "release," "released" or "releases" should be added as appropriate. Note: In addition, please see the marked-up copy of Bill 167 submitted by the CCPA, which contains a complete record of the CCPA's recommended amendments.

(CCPA)

Section 1 Purpose

Add to the purposes of the Bill the following:

- promotion of safer alternatives to toxics; and
- application of the precautionary principle and principles of sustainable development to the above goals.

(CELA, RNAO)

Add to the purposes of the Bill the following:

- the reduction or elimination of toxic releases (and not just their use and creation); and
- recognition of Ontarians' right to know the identity and amounts of toxics that are used, created, occur in consumer products or are released into the environment or workplace.

(RNAO)

Amend section 1(a) to read as follows:

(a) to prevent pollution and protect human health and the environment by reducing the use and creation of toxic questions.

(CCSPA)

Section 2 Definitions

"facility"

Define the term "facility"; a definition should at minimum discuss the nature and size of establishments that will be subject to Bill 167.

(CCSPA)

"substance of concern"

Bill 167 should provide at least a minimal description as to what is the basis for a "concern."

(CCTFA)

If the purpose of defining substances as "substances of concern" is to report on these substances because they are not on the federal National Pollutant Release Inventory (NPRI), then instead of setting up a separate reporting regime, Ontario should seek to have these substances added to the NPRI.

(CCPA)

Delete the term "substance of concern" as it will be redundant if an appropriate definition of "toxic substance" is included.

(CCSPA)

"toxic substance"

Harmonize the federal/provincial management of toxic substances by amending the definition of "toxic substance" to read as follows:

"toxic substance" means a substance prescribed by the regulations and is consistent with Federal *Canadian Environmental Protection Act, 1999* Schedule 1 list of toxic substances, or has been determined as toxic through applying an equivalent process and criteria as contained under the Federal Chemicals Management Plan, for the purposes of this Act.

(CPPI, SLEA)

Amend the definition of "toxic substance" to read as follows:

"toxic substance" means a substance on Schedule 1 of the *Canadian Environmental Protection Act, 1999* and prescribed by the regulations as a toxic substance for the purposes of this Act.

(CCSPA, CCPA, CME, CPCA, CPIA)

Amend the definition of "toxic substance" to state as follows:

"toxic substance" means:

A substance is toxic if it is entering or may enter the environment in a quantity or concentration or under conditions that

- have or may have an immediate or long-term harmful effect on the environment or its biological diversity;
- constitute or may constitute a danger to the environment on which life depends; or
- constitute or may constitute a danger in Canada to human life or health.

(CVMA)

Section 3 Requirement for toxic substance reduction plans

Threshold of Toxics Used or Created

Reduce the thresholds with respect to amounts used or created to approximately 1% of the National Pollutant Release Inventory thresholds (100kg or lower), and lower the threshold to 50kg for high hazard priority chemicals within five years of implementing the regulations made under Bill 167.

(CCS, ED/USW, OPHA)

Consider lower thresholds than those used by the National Pollutant Release Inventory for substances that are carcinogens, reproductive toxins, persistent or bioaccumulative.

(CELA)

Commit to the goal of comprehensive coverage of toxic production and use, and not just to coverage of a limited number of industries, or to the very large users and creators of toxics. All emitters and users who reach threshold levels must report. Reporting thresholds must be significantly lower than those of the National Pollutant Release Inventory, and they must be lowered over time.

(RNAO)

Provide for the same reporting exemptions as are provided for with respect to the calculation of mass reporting thresholds, as set out in section 3 of Schedule 2 of the *Notice with respect to substances in the National Pollutant Release Inventory for 2008*, dated 16 February 2008.

(CVMA)

Applicable Sectors

Make the requirement of toxic substance reduction plans applicable to all sectors, not just manufacturing and mineral processing.

(CCS, CPIA, OPHA)

Consider having Bill 167 apply to all sectors that report to the National Pollutant Release Inventory, or in the alternative, to any industrial facility that has an approval to emit contaminants to air or deposit them on land under the *Environmental Protection Act* or discharge contaminants to water under the *Ontario Water Resources Act*.

(CELA)

Bill 167 should apply to sewage treatment plants, which are responsible for 87% of mercury, 37% of arsenic, 71% of lead and nearly all chlorine releases into Ontario waterways.

(ED/USW)

Employee Threshold

Reduce the proposed ten-employee standard to a five-employee threshold.

(CCS, OPHA)

Small- and medium-sized businesses should not be exempted from the application of Bill 167. There should be no employee threshold.

(ED/USW)

General

Bill 167 should include provisions for reviewing and lowering reporting thresholds over time.

(TPH)

Section 4(1) Contents of plan

Amend subsection (1), prior to paragraph 1, to read as follows:

4.(1) A toxic substance reduction plan will consider, in accordance with the regulations, the following:

Subsection (1) should be amended to require a description of material processes at the facility collectively, rather than each substance separately.

(CME)

Recommendations re Specific Paragraphs

Paragraph 1

Add the words "on a risk prioritized basis" to the end of paragraph 1, prior to subparagraph i, to read:

1. Subject to paragraph 2, a statement that the owner or the operator of the facility intends, on a risk prioritized basis

(CPPI)

After subparagraph ii, add a new requirement which will read as follows:

iii. To reduce the level of emissions of toxic substances for the total facility, on a risk prioritized basis.

(CPPI)

Paragraph 4

The contents of toxic substance reduction plans should be facility-based, not process-based.

(SLEA)

This paragraph should be amended to require a description of material processes at the facility collectively, rather than each substance separately.

(CME)

The processes that are the most significant sources of substances should be considered when developing plans. Amend clause C of subparagraph ii by adding the words "significant processes."

(CME)

Amend this paragraph to read as follows:

4. A description of material processes at the facility that uses or creates the toxic substance, including,

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- i. a description of how, when, where and why the substance is used or created, and
- ii. quantifications that,
 - A. were made under section 9 before the plan was prepared,
 - B. were used to prepare the plan, and
 - C. show, as of the time the quantifications were made, how the substance entered the significant processes, whether it was created, destroyed or transformed during the process, and what happened to it after it left the significant processes at the facility.

(CVMA)

Amend this paragraph, prior to subparagraph i, to read as follows:

4. A description of the total facility that uses or creates the toxic substance, including,

(CPPI)

Amend subparagraph i to read as follows:

i. a description of how, when and where the substance is emitted from the total facility, and

(CPPI)

Paragraph 5

Amend this paragraph to read as follows:

5. a description and analysis of the options that were considered from an emission risk exposure basis, as warranting reducing the use and creation of the toxic substances at the facility.

(CPPI)

Proposed Paragraph 9

Add a new paragraph 9 to read as follows:

9. Facilities are exempt from complying with Section 4(1) paragraphs 4, 5 and 6 as it relates to the use and processing of raw material feedstocks from nature, such as crude oil, rock, trees. Plans to reduce toxic substances from facilities using raw material feedstocks from nature will be based on reducing the level of toxic substances that are emitted to the air, water and land.

(CPPI)

Other

Include in the contents of the plan a statement of intent to reduce the discharge of toxics to the air, land or water.

(RNAO)

The focus on emissions by toxic substance reduction plans should be based on a scientific evaluation of risk exposure.

(SLEA)

Clarify the definition of "creation" or develop differing requirements depending on the type of facility, focusing on the reduction of use of "toxic substances" in manufacturing articles (consumer products); creation of toxic by-products in the production of "value substances" (those for sale on the market); and release of "value substances" and intermediates into the environment and to humans from production of "value substances." Defined terms should align with those used under the REACH Regulation.

Bill 167 should include additional options for managing toxic chemicals other than reducing their use. These options could include environmental performance agreements and management plans developed under the *Canadian Environmental Protection Act*, 1999.

(PP)

(PP)

Section 8 Summary of Plan

Disclosure to the Public

Summaries should be made public through an easily-searchable and easilyaccessible format on the internet. A mechanism should also be in place to answer specific questions the public may not have about toxic substances in their environment.

Ensure the public's right to know about toxics in their environment, workplaces and products by collecting all necessary toxics data and making it available in a readily searchable format; and making available not only data collected under Bill 167, but also under all other environmental legislation.

The information made available to the public should be limited to the risk of exposure to emissions and not the use or presence of toxic substances in a facility.

Content of Plan Summaries

There should be a mandatory provision requiring employers to share information concerning the health consequences of exposure to toxic substances in the workplace with employees.

Bill 167 should require companies to disclose their actual use, creation and release of toxic substances.

(OPHA)

(CCS)

Section 9 Toxic substance accounting

Specific Amendments

Toxic substance accounting should be done on a total facility level, not on an individual process level. This section should be amended to read as follows:

9. The owner and the operator of a facility who are required under section 3 to ensure that a toxic substance reduction plan is prepared for a toxic substance shall ensure that, for the total facility, the net use and the total emissions of the substance from the total facility are quantified.

(RNAO)

(CCS, RNAO)

(SLEA)

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Add a new sentence under section 9 to read as follows:

Facilities are exempt from complying with Section 9 as it relates to the use and processing of raw material feedstocks from nature, such as crude oil, rock, trees. Plans to reduce toxic substances from facilities using raw material feedstocks from nature will be based on reducing the level of toxic substances that are emitted to the air, water and land.

Amend this section to read as follows:

9. The owner and the operator of a facility who are required under section 3 to ensure that a toxic substance reduction plan is prepared for a toxic substance shall ensure that, for processes at the facility which are significant that uses or creates the substance, the substance is tracked and quantified, in accordance with the regulations.

(CVMA)

(CPPI)

Method of Toxic Substance Accounting

Do not dictate what type of toxic substance accounting system a company uses. If a company is already using a recognized accounting method, it should not be required to change.

(CME)

Limit the prescriptive requirement for materials accounting and reporting and allow for other methods (*e.g.*, direct monitoring of emissions, engineering calculations, risk assessments, use of emission factors) as determined by professional judgment.

(SLEA)

Bill 167 needs to utilize the full suite of tools available for reducing the use of toxics. In addition to toxic substance accounting, Bill 167 should provide for the use of root cause analysis and business case for implementation.

(OCETA)

Account Only for Significant Sources

Ensure that only processes which provide significant sources of substances be tracked and quantified.

(CME)

Section 10(3) Information available to public

Add the following sentences to the end of this subsection:

A facility is not required to disclose the use or presence of certain toxic substances, if providing that information publicly is disclosing company proprietary information that could advantage competitors in other jurisdictions, or for cases that would cause increased security concerns. A facility is not required to report the level of toxic substances in products manufactured in Ontario, where that requirement is not applied to products manufactured outside of Ontario and imported for resale in Ontario.

(CPPI)

Section 10(4) Information available to public

To better inform Ontarians about toxic substances, Bill 167 should

- stipulate that information will be available electronically via the internet, and specify a timeline for the launch of such an internet resource;
- require the director to ensure that progress reports are made publicly available via the internet;
- require that information be reported in a user-friendly manner (*i.e.*, geographically, linked with regional maps)—the information should be presented with consistent, interoperable standards such as Canadian Geospatial Data Infrastructure; and
- require that additional contextual information be provided for each toxic substance listed, including the rationale for inclusion on the list, and subsequent management measures taking place at the company—disclosure of management measures used by facilities should be shared amongst facilities to develop "best practices" throughout the industry.

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(PP)

The Director should not be permitted to link a toxic chemical to a specific consumer product unless that product has been assessed under the *Canadian Environmental Protection Act, 1999* and determined to be toxic.

Require the Director, prior to releasing any information publicly, to carry out an economic assessment of potential damage to another substance through misinterpretation by the public of information which the Director releases.

(CPIA)

(CPIA)

Section 22(1) Disposition of things seized

Provide scientific criteria for defining a "risk to health" that allow for the disposal of a thing, in accordance with this section, which was seized under section 19.

(CCSPA)

Section 29(6) Absolute liability

This provision should be amended to remove absolute liability and at least provide for a defence of due diligence as is used in the *Canadian Environmental Protection Act*, 1999.

(CCTFA).

Section 31 Appeal of order

Bill 167 should provide a simple procedure for appeals.

(EP)

Section 44 Document prepared for another purpose

Provide equivalency with other certified environmental management systems (EMS) such as ISO 14001 with no changes to the EMS. Provide powers to the Directors to recognize such plans under section 44.

(CME)

Amend section 44 to read as follows:

44. A document that was prepared for another government or for any other purpose may, if it deals with any of the requirements of this Act and the regulations, be used in the preparation of, or as part of, a document required under this Act, as long as the document meets the intended purpose of this Act and the regulations.

Section 49(1) Regulations

Prescribing Substances as Toxic Substances or Substances of Concern More substances should be subject to reduction under Bill 167.

Commit to the goal of comprehensive coverage of toxics, and not just to coverage of a limited number of toxics.

Refrain from classifying substances as toxic substances solely on a consideration of their inherent toxicity, without a disciplined consideration of exposure, which is a critical element of full risk evaluation and thoughtful management of chemical substances.

Additional effort should be directed towards listing specific substances. Consider eliminating from Phase 1 some high volume, but relatively less toxic chemicals, such as aluminum and compounds, copper and compounds, and zinc and compounds, which would significantly add to reporting facilities' challenges with relatively less toxics reduction impact. Consider the Massachusetts model, where metals and alloys were excluded from the state's toxic reduction legislation.

Conduct *a priori* assessments of alternatives before classifying substances as toxic.

Provide scientific criteria and processes for the placement or deletion of substances on the lists of toxic substances or substances of concern.

The following principles should be considered when defining the approach to developing a list of toxic substances:

There should be a transparent process that identifies specific criteria or considerations for inclusion on or removal from the list of toxic substances.

(OMA)

(RNAO)

(OPHA)

(OMA)

(CVMA)

IS

(OMA)

(CCSPA)

- Data and rationales used for determining whether a substance is toxic should be made publicly available.
- Analysis of alternatives to toxic substances should be evaluated using the same criteria as is used for determining whether a substance is toxic— consider using the methodologies developed by the European Union or the Massachusetts Toxics Use Reduction Institute.
- An analysis of the impact on the environment and human health from the use of an alternative substance or substitute for a toxic chemical should be required in addition to the feasibility study.
- Where assessment is pending for substances of concern, and to substitutes for toxic substances, the precautionary principle should be applied and careful consideration given to the potential impact on human health and the environment.
- The government should consider the approaches and decisions of other jurisdictions, such as those of the European Union under the REACH Regulation. Mechanisms should be included in Bill 167 to encourage the consideration of new data generated under the REACH Regulation.
- Consider an analysis of substance life-cycles for substances that, while inherently toxic, have a low risk of exposure (*e.g.*, nickel).

Upon the initial receipt of data on substances of concern, the Minister should be required to take action upon assessing the information. Possible actions may include designating the substance as a toxic substance, removal of the substance from the list of substances of concern or developing an alternative strategy within a fixed time frame.

(PP)

(PP)

The government's proposed list of toxic substances should be expanded to include toxic chemicals under the *Canadian Environmental Protection Act, 1999*, even if they are not listed under the National Pollutant Release Inventory, and should automatically be updated to include chemicals recognized by the International Agency for Research on Cancer and toxics under California's *Safe Drinking Water and Toxic Enforcement Act* (Proposition 65).

(ED/USW)

Regulatory Approaches

Regulations made under Bill 167 should be

- science-based;
- collaborative and compatible with the federal government regulations under the *Canadian Environmental Protection Act, 1999* and Bill C-6 (the *Canada*

Consumer Product Safety Act), without duplication and extra burdens on business;

- based on an economic cost/benefit analysis conducted prior to the coming into force of regulations; and
- required to demonstrate that they will lead to an improvement in health or safety outcomes.

Any use of terms such as "green" or "better alternative" should be backed by science, not subjectively used to sound like an improvement.

(CCSPA)

Regulations should take into account the goals and objectives of other government initiatives such as the Open for Business campaign and budget commitments such as the 25% reduction in regulatory burden.

(CPCA)

Combining Facilities

Clarify or delete s. 49(1)(b), which permits the Minister to combine "related" facilities to define them as a "single facility," even if they have different owners or operators.

(CCSPA)

Sections 50 to 64

Delete these provisions of Bill 167.

(CCPA)

Section 64, amending s. 49(1) Regulations

Coming Into Force

Bring this section into force at the same time that s. 49(1) comes into force.

(RNAO)

Scope of Substances Regulated

Within one year of the passage of Bill 167, the government should identify priority substances and products for regulation, using a precautionary approach. Early action should focus on formaldehyde, lead, vinyl chloride, bis(2-ethylhexyl) phthalate and bisphenol A.

(ED/USW)

Consumer Product Labelling

All ingredients in consumer products should be fully disclosed on product labels. If carcinogens are present in products, they should be identified by a hazard symbol. The full ingredient list and hazard symbol should be visible to the consumer at point of sale and at point of use and presented in clear language.

(CCS, OPHA, RNAO)

The application of Bill 167 to consumer products should be clarified. At a minimum, Bill 167 should authorize labelling and warnings with respect to toxic substances in consumer products where the substances are capable of causing or have effects such as reproductive toxicity.

(CELA)

Exempt vehicles from the consumer protection provisions of Bill 167, as they are already covered by federal legislation. Add to section 64 a schedule similar to Schedule 1 of Bill C-6 (the *Canada Consumer Product Safety Act*) setting out those consumer products that are exempt from regulation under this provision.

(CVMA)

Bill 167 should include timelines for identifying priority substances and consumer products for regulation and labelling.

(TPH)

Require a risk-based approach to product regulation including consideration of both inherent toxicity and exposure pathways, and requiring consultation with affected parties prior to public notice of the intent to regulate.

(CAC/SMC)

Alternatives to Section 64

This provision should be deleted from Bill 167; in the alternative, consumer products and their ingredients already regulated under the federal *Food and Drugs Act* and *Canadian Environmental Protection Act, 1999* should be specifically exempted.

(CCTFA)

The authority to ban or restrict the manufacture, distribution or sale of a product known to contain a toxic substance should be vested in the federal government to avoid the Balkanization of the Canadian marketplace.

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(CPPI)

Delete subparagraph (ii) from the proposed amendment to s. 49(1).

(CCSPA)

OTHER RECOMMENDATIONS

Communication Along the Supply Chain

There should be communication from producers and importers to their downstream users (customers) and vice versa. Safe use guidelines, including risk management measures and operational conditions should be identified and shared amongst the entire supply chain.

Confidential Information

Bill 167 should clearly state that manufacturers' confidential information will not be disclosed to competitors.

Conflict with Other Laws

In the event of conflicts between Bill 167 and municipal bylaws or other provincial environmental legislation, allow the provision which is the most protective of health or the environment to prevail.

(CELA, RNAO)

Duplication Concerns

Provide equivalency with other certified environmental managements systems (EMS) such as ISO 14001 without requiring changes to the EMS and provide powers to the Ministry Directors to recognize such plans under Bill 167.

(CVMA)

Harmonize Bill 167 with the federal Chemical Management Plan.

(APMA, CPCA)

Harmonize Bill 167 with the Canadian Environmental Protection Act, 1999.

(CPIA)

(PP)

(CME)

Formally recognize the potential for overlap and duplication with federal efforts including the federal Chemical Management Plan and the *Canadian Environmental Protection Act, 1999*, and provide the Minister of the Environment with a specific directive to avoid overlap and duplication.

(CAC/SMC)

European Union *REACH* (*Registration, Evaluation, Authorisation and Restriction of Chemicals*) Regulation

Strengthen Bill 167 to account for the impact of the European Union's *REACH* regulation and other jurisdictions' chemical management plans so that Ontario's industries meet *REACH* standards and can export to Europe.

(CCS)

Funding for a Toxics Reduction Strategy

The toxics reduction strategy and its related institutions should be funded by fees levied on the regulated community.

(CCS, CELA, PP, RNAO)

The funding formula should be based on use and emission, with the weighting towards emissions.

(RNAO)

The \$24 million allocated by the Ministry of the Environment is not sufficient. Other ministries, especially the Ministry of Research and Innovation, need to be proactive in supporting the toxics reduction strategy. Support efforts by the Ministry of Research and Innovation to increase resources for university, government and industry research, including support to existing industries to accelerate the commercialization and marketing of new products that reduce the use of toxics.

(OBAC)

Independent Legal Review of Bill 167

There should be an independent legal review of the key provisions of Bill 167, as the rolling incorporation provision in s. 49(2) means that a change in a document outside of the regulations becomes incorporated into the regulations with no political oversight, and critical definitions and criteria related to the operation of the Bill have not been established.

(CCSPA)

Monitoring Program

Commit to an integrated monitoring program for toxic substances in the environment (*i.e.*, water, air and soil) and the human population (*i.e.*, biomonitoring) to identify where further action will be required to prevent pollution and protect human health and the environment.

Pre-Impact Safety Valve

Bill 167 should include a pre-impact safety valve to identify and notify those who may be adversely impacted by the designation of substances as being toxic substances or substances of concern.

(EP)

(PP)

Proposed Amendment to Occupational Health and Safety Act

Reinstate s. 36 of the Occupational Health and Safety Act (OSHA), which was repealed in 2001; this provision required employers to inventory the chemicals they used on an annual basis. Mandate that joint health and safety committees (established under OSHA) consider alternatives or substitutes to toxic chemicals (as is done under British Columbian and federal law) and require an employer to report to the joint committee and the Minister of Labour on the progress achieved with respect to the removal of toxic chemicals from the workplace.

(ED/USW)

Proposed Preamble

Include in the preamble an endorsement of the principles of the *Canadian Environmental Protection Act, 1999*—specifically the precautionary principle, pollution prevention, virtual elimination of persistent and bioaccumulative toxic substances, and the "polluter pays" principle.

(RNAO)

Reduction Targets

Bill 167 should include goals or targets for reducing the use and release of toxic substances in Ontario.

(CCS, ED/USW, PP, TPH)

Include aggressive targets for reductions in the use, creation and release of toxics, including:

- a 50 % reduction in toxic releases within five years of Bill 167 coming into force;
- a 20 % reduction in toxic use within five years; and
- a 40 % reduction in toxic use within ten years.

(RNAO)

Bill 167 should set a target of reducing the use, creation and release of toxics by 50% within 5 years.

(CELA, OPHA)

Reduction targets should be renewable.

(OPHA, PP)

Role of the Public

To enhance the role of the public, Bill 167 should be amended to provide for the following:

- the public's right to know information, beyond what is contained in the summaries of toxics reduction plans (as prescribed by s. 4 of the Bill), compiled under the authority of existing environmental laws;
- the public's right to apply to the Minister for review of toxics use reduction (and safe alternative) plans, or in the alternative, amendment of the *Environmental Bill of Rights*, 1993 (*EBR*) to ensure that such plans are included in the definition of "instruments," and therefore, subject to review under the *EBR*; and
- a public right of action to enforce key provisions of Bill 167.

(CELA)

Sector-Specific Approach

Include specific provision for a sector-based approach that includes risk-based prioritization of substances.

(CAC/SMC)

Substitution of Safer Alternatives

Substitution should be mandated in situations where a safer alternative exists or where the use is non-essential, as is now required under the European Union's *REACH* regulation.

(CCS, CELA, ED/USW, OPHA, PP, RNAO)

The Minister should be able to create a schedule listing specific substances to be reduced, as well as suggested alternative substances acceptable to the Minister. The Minister should certify that suggested alternatives are approved for use by Health and Environment Canada, do not have greater risks than the substance they are replacing, and perform as ingredients as portrayed.

(CCTFA, ED/USW)

The Minister should be able to create a schedule listing specific substances to be reduced, as well as suggested alternative substances acceptable to the Minister. There must be a substitution assessment framework and methods for evaluating alternatives.

(Diamond)

There should be mandatory phase-outs or substitution of high-hazard substances. The provisions for regulating such substances should ensure that the process for doing so is open, be subject to regular review so the list of substances reflects scientific developments and have specific dates for companies to achieve elimination or substitution of high-hazard substances.

(TPH)

Technical Assistance Programs

Bill 167 should establish a regime to provide technical assistance to employees who require re-employment assistance, vocational retraining or other assistance as a result of the implementation of the Bill.

(CELA)

Bill 167 should establish a regime to provide technical and financial assistance to small businesses, regardless of whether they are subject to Bill 167, that require assistance in reducing their use of toxic substances.

(CELA)

Capacity building under Bill 167 should support small- and medium-sized facilities that use and release priority substances under National Pollutant Release Inventory thresholds.

Toxic Use Reduction Institute

The government should establish an independent university-based research Toxic Use Reduction Institute (TURI) to advance the province's capacity for toxic use reduction activities, safe substitution, green chemistry, education and information outreach and training on toxics reduction planning. TURI should provide mandatory training on toxics reduction plans and offer training and certification for toxics reduction planners.

(CCS, CELA, Diamond, ED/USW, OPHA, PP, RNAO, TPH)

A third-party institute should be established to work with industry and government to research and develop toxic reduction strategies and outreach. It should coordinate information sharing and encourage research and development in partnership with industry. It should foster the sharing of solutions and case studies in a Web 2.0 environment through an efficient and logical data collection process, and assist manufacturers with Environmental Value Stream Mapping.

(APMA)

(TPH)

Toxic Use Reduction Planners

The role of toxic use reduction planners should be expanded and specifically linked with water and energy conservation.

(ED/USW)

WITNESS LIST

Abbreviation	Organization/Individual	Date of Appearance
APMA	Automotive Parts Manufacturers' Association	25 May 2009
CAC/SMC	Cement Association of Canada/St. Marys Cement	25 May 2009
CCS	Canadian Cancer Society, Ontario Division	13 May 2009
CCPA	Canadian Chemical Producers' Association	25 May 2009
CCSPA	Canadian Consumer Specialty Products Association	25 May 2009
CCTFA	Canadian Cosmetic, Toiletry and Fragrance Association	13 May 2009
CELA	Canadian Environmental Law Association	13 May 2009
CME	Canadian Manufacturers and Exporters	25 May 2009
CPCA	Canadian Paint and Coating Association	25 May 2009
CPIA	Canadian Plastics Industry Association	25 May 2009
CPPI	Canadian Petroleum Products Institute	13 May 2009
CVMA	Canadian Vehicle Manufacturers' Association	25 May 2009
Diamond	Miriam Diamond	25 May 2009
ED/USW	Environmental Defence/United Steelworkers	13 May 2009
EP	Environmental Plastics	25 May 2009
OBAC	Ontario BioAuto Council	25 May 2009
OCETA	Ontario Centre for Environmental Technology Advancement	25 May 2009
OMA	Ontario Mining Association	25 May 2009
OPHA	Ontario Public Health Association	13 May 2009
PP	Pollution Probe	25 May 2009
RNAO	Registered Nurses Association of Ontario	13 May 2009
SLEA	Sarnia Lambton Environmental Association	13 May 2009
ТРН	Toronto Public Health	25 May 2009

Legislative Assembly of Ontario



Assemblée législative de l'Ontario

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> 090094 May 28, 2009

MEMORANDUM TO:	Standing Committee on General Government
FROM:	Marta Kennedy Research Officer

SUBJECT:

Background Information on Bill 167, the *Toxics Reduction Act, 2009*

The Standing Committee on General Government is currently holding public hearings on Bill 167, the *Toxics Reduction Act*, 2009.

After a presentation by the Ontario Mining Association, the Committee requested that the Legislative Research Service provide a copy of a review done in Massachusetts that lead to the exclusion of metals and alloys from the Massachusetts *Toxics Use Reduction Act* (Massachusetts TURA). A copy of this document was provided by the Ontario Mining Association to the Legislative Research Service and is attached.

The Committee also asked for a memorandum that specified the reasons for removing metals and alloys from the Massachusetts TURA.

This memo includes information about alloys that have been removed from the application of the Massachusetts TURA. It appears, though, that certain metals may still be subject to that Act. We are awaiting confirmation on this and will provide a follow-up memo once we receive further information.

MASSACHUSETTS TOXICS USE REDUCTION ACT

The Massachusetts *Toxics Use Reduction Act* was signed into law by the Governor of Massachusetts on July 24, 1989.¹ The legislation requires Massachusetts companies that use large quantities of specific toxic chemicals to evaluate pollution prevention opportunities, implement them if practical, and measure and report their results every year. Companies must also evaluate their efforts and update their toxics use reduction plans every other year. The chemicals that must be reported are those on the Toxic or Hazardous Substances List established pursuant to the Massachusetts TURA.²

Delisting of certain alloys from the Massachusetts Toxic or Hazardous Substances List

The Toxic or Hazardous Substances List (Substances List) was originally created from two federal lists: the Toxics Release Inventory (TRI) list created by the *Emergency Planning and Community Right to Know Act* and the *Comprehensive Environmental Response, Compensation, and Liability Act* (CERCLA) list.³ Substances on the CERCLA list originate from four other U.S. federal regulatory lists: the *Clean Air Act* list of hazardous air pollutants (HAPs), the *Clean Water Act* list of hazardous substances and priority pollutants, the *Solid Waste Disposal Act* list of hazardous wastes, and the *Toxic Substances Control Act* list of substances that pose an imminent hazard.⁴

A person may petition the Science Advisory Board (SAB) for the delisting of a chemical from the Substances List. The SAB was established under the Massachusetts TURA to make recommendations with respect to the addition and deletion of substances from the Substances List. The SAB considers the petition and makes recommendations to the Administrative Council on whether to add or remove a substance from the list. The Administrative Council is the governing body of the Massachusetts TURA program and has the power to officially list or delist substances.⁵

¹ Massachusetts, Department of Environmental Protection, *Fact Sheet: About the Toxics Use Reduction Act (TURA) Program*, Internet site at <u>http://www.mass.gov/dep/toxics/turafact.pdf</u>, accessed 27 May 2009.

² Massachusetts Department of Environmental Protection, "Toxics Use Reduction Act (TURA) Program Overview," Internet site at <u>http://www.mass.gov/dep/toxics/tura/turaover.htm</u>, accessed 26 May 2009.

³ Massachusetts Toxics Use Reduction Act, Mass. Gen. Laws ch. 21I, § 9(A)-(B) (2009), Commonwealth of Massachusetts, General Court, Internet site at

http://www.mass.gov/legis/laws/mgl/21i-9.htm, accessed 26 May 2009.

⁴ Massachusetts Toxics Use Reduction Institute, *Policy Analysis: Recommendations on CERCLA chemicals that have never been reported by TURA filers* (Lowell, MA: Massachusetts TURI, 2008), 2. TURI Internet site at

http://www.turi.org/policy/ma_tura_program/tura_administrative_council/recent_council_consider ations/policy_analyses_for_cercla_chemicals/policy_analysis_for_cercla_chemicals_never_report ed_under_tura/non_reported_cercla_chemicals, accessed 27 May 2009.

⁵ Administrative Council on Toxics Use Reduction, *Toxics Use Reduction in Massachusetts* (Massachusetts: Administrative Council on Toxics Use Reduction, November 2008). TURI Internet site at <u>http://www.turi.org/policy/ma_tura_program/tura_report_to_governor_patrick</u>, accessed 26 May 2009.

Effective 1995, the following alloys were delisted:

copper, nickel, chromium, cobalt or manganese in a solid or molten metal alloy, but not including aerosols, where aerosols are defined as particles less than 50 um (microns) in diameter.⁶

An alloy is

a substance possessing metallic properties and composed of two or more elements of which at least one must be a metal. The term refers to those cases where there is an intentional addition to a metal for the purpose of improving certain properties.⁷

The Guidance Document: Reporting and Planning for Certain Metal Alloys, produced by the Massachusetts Department of Environmental Protection, says

the delisting covers metal alloys in solid or molten form, such as copper in bronze and brass, and nickel and chromium in stainless steel.

This delisting does not extend to dissolved metals, metal alloys in aerosol form, metal alloy particles less then 50 microns in diameter, or *any of the five metals in non-alloy form;* i.e., a metal containing a low level of impurities is NOT an alloy. For example, wire comprised of 99+% copper and less than 1% impurities is considered to be essentially pure copper, and not an alloy.

The use of metals in pure form can be subject to reduced TURA reporting requirements and exempted from TURA planning and fees provided specific requirements are met. Detailed information on this expanded article exemption can be found in the "BWP 94-014 Policy for TURA Reporting and Planning for Certain Metalworking Operations" and the

⁶ Toxic or Hazardous Substance List, 301 Mass. Code Regs. 41.03(1)(a) (2009), Executive Office of Energy and Environmental Affairs Internet site at

http://www.mass.gov/Eoeea/docs/eea/ota/tur_prog/301cmr41_toxic_haz_sub_list.pdf, accessed 27 May 2009.

⁷ Massachusetts Toxics Use Reduction Institute, *Glossary*, s.v. "alloy," TURI Internet site at <u>http://www.turi.org/turadata/glossary/alloy</u>, accessed 27 May 2009.

"Recycling Activity Report Form."⁸ [emphasis added]

The Policy for TURA Reporting and Planning for Certain Metalworking Operations is attached as Appendix B. The Guidance Document: Reporting and Planning for Certain Metal Alloys is attached as Appendix A.

The delisting of these alloys is mentioned in the review referred to by the Ontario Mining Association, *Categorization of the Toxics Use Reduction List of Toxic and Hazardous Substances.*⁹ Note, though, that this document deals only peripherally with the delisting of these alloys.

Sterling silver alloy and silver in alloy form have also been delisted.¹⁰

The following table summarizes the delisting decisions made to date with respect to alloys. All but two delisting were the result of industry petitions.

Summary of Delisting Decisions ¹¹					
Chemical	SAB Recommendation	Supplemental Information	Status or Outcome		
Nickel in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per TURI/SAB recommendation.		
Chromium in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per TURI/SAB recommendation.		
Copper in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per TURI/SAB recommendation.		
Manganese in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per TURI/SAB recommendation.		

⁸ Massachusetts Department of Environmental Protection, *Guidance Document: Reporting & Planning for Certain Metal Alloys* (Massachusetts: Mass DEP, 1995), Internet site at <u>http://www.mass.gov/dep/toxics/laws/metals.htm</u>, accessed 27 May 2009.

⁹ Massachusetts Toxics Use Reduction Institute, *Categorization of the Toxics Use Reduction List of Toxic and Hazardous Substances. Methods and Policy Report No. 18* (Lowell, MA: Massachusetts TURI, 1999), 3.

¹⁰ Toxic or Hazardous Substance List, 301 Mass. Code Regs. 41.03(1) (2009).

¹¹ Email, Policy Analyst, Massachusetts Toxics Use Reduction Institute, 27 May 2009.

Summary of Delisting Decisions ¹¹					
Chemical	SAB Recommendation	Supplemental Information	Status or Outcome		
Cobalt in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per TURI/SAB recommendation.		
Sterling silver alloy	delist copper-silver alloys except for aerosols (less than 50 um)	Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council with qualifications as per TURI/SAB recommendation.		
Copper in alloy form	delist except for aerosols (less than 50 um)	Reconsideration. Delisting originated in SAB to be consistent with previous decisions. Unanimous vote. Aerosols should be reported under TURA because planning for efficient use is beneficial.	TURI/SAB recommendation accepted by Admin Council		
Silver in alloy form	delist except for aerosols (less than 50 um)	Delisting originated in SAB to be consistent with previous decisions. Unanimous vote Aerosols should be reported under TURA because planning for efficient use is beneficial.	TURI/SAB recommendation accepted by Admin Council		

Process for delisting substances

Currently, the SAB considers the following information before making a recommendation to the Administrative Council for the delisting of a substance:

- International Agency for Research on Cancer rating,
- data from the Environmental Protection Agency PBT Profiler (persistence in water, soil, sediment, and air; bioconcentration factor; and chronic toxicity in fish),
- neurotoxicity,
- developmental/reproductive toxicity,
- mutagenicity,
- lethal dose or concentration information (LD50 and LC50),
- exposure limits required or recommended by U.S. federal agencies,
- flash point.¹²

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¹² Massachusetts Toxics Use Reduction Institute, *Policy Analysis: Recommendation to take no action on certain CERCLA chemicals that have been reported by TURA filers* (Lowell, MA:

However, only limited information is available about the process the delisted alloys actually went through before being delisted. Whether these factors were considered before SAB made its recommendation to the Administrative Council is unknown at present. What is clear is that the SAB of the time unanimously voted to recommend delisting of these alloys, and the Administrative Council accepted their recommendations.

Massachusetts TURI, 2008), 1. TURI Internet site at

http://www.turi.org/policy/ma_tura_program/tura_administrative_council/recent_council_consider ations/policy_analyses_for_cercla_chemicals/policy_analysis_for_cercla_chemicals_reported_und er_tura/reported_cercla_chemicals_recommended_for_no_action, accessed 27 May 2009.

APPENDIX A

GUIDANCE DOCUMENT: TURA REPORTING AND PLANNING FOR CERTAIN METAL ALLOYS¹³

The following substances have been delisted from the Toxics and Hazardous Substance List (301 CMR 41.00) effective reporting year 1995:

Copper, nickel, chromium, cobalt or manganese in a solid or molten metal alloy, but not including aerosols, where aerosols are defined as particles less than 50 um (microns) in diameter.

Questions have been raised regarding implementation of this delisting. This guidance has been produced to assist the regulated community in complying with TURA. If you have any remaining questions after reading this document, please contact the Department of Environmental Protection (MassDEP) for compliance assistance or the Office of Technical Assistance for further information.

General Information

For the purposes of determining what a metal alloy is, a useful definition of a metal alloy is given by "An Encyclopedia of Metallurgy and Materials" (MacDonald and Evans, 1984):

"An alloy is a substance possessing metallic properties and composed of two or more elements of which at least one must be a metal. The term refers to those cases where there is an intentional addition to a metal for the purpose of improving certain properties."

The delisting covers metal alloys in solid or molten form, such as copper in bronze and brass, and nickel and chromium in stainless steel.

This delisting does not extend to dissolved metals, metal alloys in aerosol form, metal alloy particles less then 50 microns in diameter, or any of the five metals in non-alloy form; i.e., a metal containing a low level of impurities is NOT an alloy. For example, wire comprised of 99+% copper and less than 1% impurities is considered to be essentially pure copper, and not an alloy.

The use of metals in pure form can be subject to reduced TURA reporting requirements and exempted from TURA planning and fees provided specific requirements are met. Detailed information on this expanded article exemption can be found in the "BWP 94-014 Policy for TURA Reporting and Planning for Certain Metalworking Operations" and the "Recycling Activity Report Form". Both of these documents can be obtained from MassDEP by calling the TURA Program at 617-292-5982.

The 50 micron size which describes metal alloy aerosols and particles refers to the particle size of the metal alloy, not the particle size of the individual listed metals.

Specific Process Questions

The following sections provide guidance for specific processes in which metal alloys are used:

• Etching processes. When a metal is removed from the metal alloy stock through chemical processing, the resultant metal salt is reportable if it is produced in threshold quantities (in this case, the metal salt is "coincidentally manufactured"). The 25,000 pound threshold would apply if the metal salt is the only reportable chemical; the threshold drops to 10,000 pounds if a facility is using other reportable chemicals. It should be noted that the entire weight of the metal salt is considered when determining

¹³ Massachusetts Department of Environmental Protection, *Guidance Document: TURA Reporting and Planning For Certain Metal Alloys* (Massachusetts: Massachusetts DEP, n.d.), Massachusetts DEP Internet site at <u>http://www.mass.gov/dep/toxics/laws/metals.htm</u>, accessed 27 May 2009.

whether a threshold has been met, and not just the amount of the metal in the salt. The stock itself, if it is a metal alloy as defined above, is not reportable.

- Electropolishing operations. The metal salts in solution are reportable as "coincidentally manufactured" if they are produced in threshold quantities. Again, the stock itself, if it is a metal alloy as defined above, is not reportable.
- Grinding Operations. The listed metals are not exempt from reporting when contained in aerosols of a metal alloy, i.e., particles less than 50 microns in diameter (in this case, the metal is reported as "processed"). These metals are reportable if they are present in a metal alloy aerosol in threshold quantities. When determining whether the reporting threshold has been exceeded, only the weight of the listed metal is considered, and not the total weight of metal alloy aerosol. The stock itself, if it is a metal alloy as defined above, is not reportable.
- Multiple processes. In a production unit where the metal alloy stock is being used in several different processes, none of the stock is reportable if it is a metal alloy as defined above. However, the individual production processes may produce metal alloy in aerosol form or metal salts in solution. Those aerosols and metal salts in solution would be reportable if produced in threshold quantities, as described above.
- Alloys in powdered form. If a metal alloy is used in particle form, the facility will need to perform a particle size analysis to determine the amount of material less than 50 microns in size. The listed metals (e.g., Cr, Cu) are reportable if they are present in metal alloy particles less than 50 microns in size in threshold quantities. When determining whether the reporting threshold has been exceeded, only the weight of the listed metal is considered, and not the total weight of metal alloy particles.
APPENDIX B

POLICY FOR TURA REPORTING AND PLANNING FOR CERTAIN METALWORKING OPERATIONS¹⁴

Bureau of Waste Prevention Policy BWP-94-014 Signed by Steve DeGabriele for Patricia Deese Stanton, Assistant Commissioner on June 8, 1994

1. POLICY STATEMENT

This document sets forth the DEP's policy [Endnote 1] regarding reduced reporting requirements and exemption from TURA planning and fees for specified metals and metal alloys processed in certain metalworking operations, including but not limited to processes involving bending, cutting, stamping, and extruding.

The reduced reporting provisions and the exemption from planning and fees may be claimed only for those metals specified in Part A of this policy and only when the metalworking operation meets the conditions, requirements, and procedures set forth, respectively, in Parts B, C and D of this policy.

A. Applicability

- 1. This policy applies only to copper and its alloys, and steels including stainless steel.
- 2. This policy applies only where a metal is processed; the policy does not apply to metals that are manufactured or otherwise used. [Endnote 2]
- This policy does not apply to (a) metallic compounds, (b) metals or metal alloys not specified in paragraph 1(a) above or (c) lead even if the lead is a constituent of an alloy listed in paragraph 1(a).

B. Conditional Exemption and Reduced Reporting for Certain Metals

DEP will allow reduced reporting and an exemption from TURA planning and fees for a metal covered by this policy *when the following conditions are met for each metal claimed under this policy.*

- 1. The metal must be an "input" to a process.
- 2. While the physical form or shape of the metal may change during processing, the metal must remain solid and the other properties of the metal, including its chemical properties, must not change.
- 3. Except for transfers of the metal as a commodity or item with value to scrap metal brokers or other recycling or reuse operations, the handling, storage, and processing activities at the facility must result, on a calendar year basis, in no releases Endnote 3] of the metal to the environment as reportable in the following sections of the Form R: Sections 5 (RELEASES OF THE TOXIC CHEMICAL TO THE ENVIRONMENT ON-

¹⁴ Massachusetts Department of Environmental Protection, *Policy For TURA Reporting and Planning For Certain Metalworking Operations* (Massachusetts: Massachusetts DEP, 1994), Massachusetts DEP Internet site at <u>http://www.mass.gov/dep/toxics/laws/metals.htm</u>, accessed 27 May 2009.

SITE) and Section 6 (TRANSFERS OF THE TOXIC CHEMICAL IN WASTES TO OFF-SITE LOCATIONS).

- 4. The scrap metal generated by the metalworking operation must be collected under a "comprehensive and aggressive" recycling program as described by EPA. [Endnote 4] The scrap metal must also be sent without treatment or further processing -- except mechanical separation processing -- to a scrap metal broker or other recycling or reuse operation.
- 5. The toxics user must comply with the reporting requirements described below.

C. Reporting Requirements

- 1. For each metal that the toxic user seeks to have covered under this policy, the toxic user shall complete and submit to the Department: a Form R, a Form S Cover Sheet, and Section 1 of the Form S. (Completion of sections 2 and 3 of the Form S will not be required if all the conditions of this policy are met.) In Section 1, the user shall report as separate items the amount of metal shipped in or as product and the amount shipped as scrap. The certification statement in the Form S Cover Sheet shall apply not only to the Form S and Form R information but also to the information described in paragraph 2 below.
- 2. Recycling Activity Report Form: The toxics user shall provide, on a form specified by DEP, such additional information necessary to demonstrate that the metalworking operation satisfies the conditions of this policy. Such information shall include, but not be limited to:
 - i. a description of the steps taken, if any, to minimize the production of scrap metal, and metal disposed of, treated, or otherwise released to the environment;
 - ii. a statement that the scrap metal was sent without treatment or further processing -- except mechanical separation processing -- to a scrap metal broker or other recycling or reuse operation.
 - iii. a description of the toxic user's recycling or reuse strategy sufficient to demonstrate that this strategy is "comprehensive and aggressive." This description shall include, without limitation, the following information:a) methods employed for collecting the metal for recycling or reuse and ensuring that all discernable metal particles are so collected; b) the methods of storing and handling of the scrap metal prior to recycling or reuse; c) the name(s) and address(es) of the scrap metal broker or other recycling or reuse operation as well the type of operation(s) to which the scrap metal was transferred (unless this information was reported in the Form R); d) information to substantiate the commodity-like nature of the scrap metal, e.g., information indicating that the scrap metal has value.

D. Procedure for Modified Reporting, and Planning and Fee Exemption, for Metalworking Operations

- The toxic user shall submit to the Department the Form R and Form S documents described above on or before July 1st for each prior calendar year for which the user seeks to proceed under this policy. The first page of the Form S Cover Sheet and Form S shall be clearly marked "submitted under the provisions of BWP Policy-94-014."
- The toxic user shall submit to the Department the Recycling Activity Report Form described above on or before July 1st for the first prior calendar year for which the user seeks to proceed under this policy. This form need not be submitted annually; provided,

however, that if a change occurs in the toxic user's recycling operations that would make it less aggressive and comprehensive, the toxic user shall so notify the Department. [Endnote 5]

- 3. Any toxic user seeking to claim confidentiality of information in the documents submitted under this policy shall follow the procedures set forth in 310 CMR 3.30.
- 4. If the Department determines, based on information in the reports submitted under this policy or other relevant information, that the metal does not qualify under this policy, then the Department shall so notify the toxic user in writing, and the toxics user shall submit to DEP a toxics use report, develop a plan and pay applicable fees by the deadline specified in the written notice.
- 5. This policy is based on information available to the Department as of the date of the policy. Should additional information indicate a need to repeal or modify the policy, the Department may take such action but any repeal or modification of this policy shall be effective for the calendar year reporting period following the calendar year in which the repeal or modification occurs.

2. BACKGROUND STATEMENT

A. Problem Definition

In part, this policy relies on the rationale behind the "article exemption" [Endnote 6] to establish the appropriate level of regulation for certain metals. Under the article exemption, toxic chemicals contained in articles are not regulated under TURA. Because, in general, an article is a discrete item when it enters the facility, and remains so during its normal use at the facility, the environmental and public health risks of a toxic substance contained in an article are generally less than the use of same toxic substance when it is not contained in an article.

Many metals that would otherwise qualify under the article exemption fail to do so because the processing of the item involves a change in shape. The DEP believes, if the conditions in this policy are met, that certain uses of the metals specified in the applicability section of this policy pose a risk similar to the reduced risk posed by metals in an article.

This is not to say, however, that there are no risks associated with metal use. For example, mining, manufacture, smelting, and general handling of metals may pose issues of concern. The DEP believes that these risks must continue to be assessed, and that, consistent with TURA and EPCRA, the public has a right to information about metal use. The level of reporting required under this policy will allow the continued assessment of risks associated with metals, and will provide the public with information about metal use.

In addition, this policy seeks to encourage metal reuse and recycling by requiring an "aggressive and comprehensive" strategy for metal reuse and recycling in order to qualify for the planning and fee exemption. While reuse/recycling, under these circumstances is NOT toxics use reduction as defined in TURA, it is beneficial as a resource conservation measure. Furthermore, although the remelting of scrap metal may entail certain risks, it does not involve the handling of ores that can contain such toxics as arsenic (in copper ores).

B. **Implementation and Evaluation** The information in these modified reports will provide a clearer picture of metal use in Massachusetts, and will allow the agencies responsible for implementing TURA to continue to assess risks associated with metal use throughout its lifecycle. The modified reports will be publicly available, as are toxics use

reports. The DEP will collect and manage data from the reports in a manner consistent and compatible with data from toxics use reports.

ENDNOTES:

- 1. While this policy is based on certain concepts and conditions found in the article exemption, this policy does not change the definition of "article" under TURA, or the interpretation or application of the article exemption.
- 2. See 310 CMR 50.00 for the definitions of the terms "manufacture," "process," and "otherwise use."
- 3. Consistent with EPA, if any amount (calculated separately for each metal) is less than or equal to 0.5 pounds, the amount may be rounded to zero. Also consistent with EPA policy, "(i)f the owner/operator has instituted a comprehensive and aggressive program for the recycle/recovery of all released material then small amounts of toxic chemical that are not, in fact, recycled/recovered due to the imperfect efficiency of virtually any recycling/recovery system should not count towards the 0.5 pound cut-off value." See EPA memorandum entitled Clarification of Article Exemption, dated June 14, 1991, from Sam Sasnett to Robert W. Hicklin.
- 4. See EPA memorandum entitled Clarification of Article Exemption, dated June 14, 1991, from Sam Sasnett to Robert W. Hicklin. In effect, this memorandum defines a "comprehensive and aggressive" recycling program as one in which the facility takes all practicable steps to collect and recycle scrap. For example, a small amount of metal that may cling to worker's clothing would not negate an otherwise comprehensive and aggressive recycling program. Sweeping up metal dust and discarding it in a dumpster could negate the comprehensiveness and aggressiveness of the program.
- DEP recognizes that metalworking operations typically choose their scrap metal dealers based on the best available price. Changes in scrap metal dealers based on pricing decisions would not be considered a change that would normally require notice to the Department.
- 6. The TURA regulations (310 CMR 50.10) define "article" as follows:
 - "a manufactured item, other than an item which is manufactured at the facility: (1) which is formed to a specific shape or design during manufacture; (2) which has end use functions dependent in whole or in part upon its shape or design during end use; and (3) which does not release a toxic substance under normal conditions of processing or use of that item at the facility or establishments."



THE MASSACHUSETTS TOXICS USE REDUCTION INSTITUTE

Categorization of the Toxics Use Reduction List of Toxic and Hazardous Substances

Methods and Policy Report No. 18

 $\hat{r}_{\mathcal{B}}$

... ,* 1999

University of Massachusetts Lowell

Categorization of the Toxics Use Reduction List of Toxic and Hazardous Substances

Report on the work of the Toxics Use Reduction Science Advisory Board

The Toxics Use Reduction Institute University of Massachusetts Lowell

March 1999



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The Toxics Use Reduction Institute is a multi-disciplinary research, education, and policy center established by the Massachusetts Toxics Use Reduction Act of 1989. The Institute sponsors and conducts research, organizes education and training programs, and provides technical support to promote the reduction in the use of toxic chemicals or the generation of toxic chemical byproducts in industry and commerce. Further information can be obtained by writing the Toxics Use Reduction Institute, University of Massachusetts Lowell, One University Avenue, Lowell, Massachusetts 01854.

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34 Coddington Street, Quincy, Massachusetts 02169 617-984-1600

Dr. Michael Ellenbecker Acting Director Toxics Use Reduction Institute One University Avenue Lowell, MA 01854-2866

Dear Dr. Ellenbecker,

Attached is a report by the Toxics Use Reduction Science Advisory Board on its "Categorization of the Toxics Use Reduction List of Toxic and Hazardous Substances" project. The Board has been working on this project for the past 18 months and is very pleased to submit this work product. It represents a concerted effort on the part of the Board to categorize 258 chemicals into three categories, high hazard, low hazard and uncategorized chemicals.

Many Board members contributed their time and expertise to this project. Those Board members are: James J. Ahearn Jr., Ph.D. from Polaroid Corporation, Andrew F. Beliveau of the U. S. Environmental Protection Agency, Richard Clapp, Sc.D. of B.U. School of Public Health, George M. Gray, Ph.D. of the Harvard School of Public Health, Center for Risk Analysis, Thomas Trayers from the Division of Occupational Safety's Occupational Hygiene Program, and Lawrence H. Boise from the Gloucester Co., Inc.; and two former Board members, Halina Brown, Ph.D. from Clark University and Christine Oliver, M. D. from Mass. General Hospital.

The Board realizes that this work will never be complete as new data on existing chemicals are generated and as new chemicals are reported in the Commonwealth. Adjustments to the list will be made based on new information in these areas. We trust that the list will aid the decisions made concerning TURA Program priorities. In addition, the Board respectfully requests to be informed of any policy decisions resulting from the use of these lists. Thank you.

Sincerely,

and J. Williams

David T. Williams Executive Director Quincy College Center for Technology & Health and Chair, Toxic Use Reduction SAB

Summary

For the past eighteen months, the Toxics Use Reduction Science Advisory Board has been working on a project to categorize the 258 chemicals which have ever been reported under the Massachusetts Toxics Use Reduction Act (TURA). The resulting lists of more hazardous¹, less hazardous and uncategorized chemicals will be used by the Toxics Use Reduction Program to aid in setting priorities and will serve as guidance for companies making chemical substitution decisions. The lists of more hazardous (Category 1) and less hazardous (Category 2) substances follow. The specific chemicals in the categories may change based on new data becoming available or new chemicals being used above TUR reportable threshold quantities in the Commonwealth.

	Table 1: Category 1 Chemic	cals ²
Acrylamide	Diethylsulfate	Nickel compounds
Acrylonitrile	Dimethylformamide	Nitrobenzene
Arsenic compounds	Dioxane	Phosgene
Arsenic	Epichlorohydrin	Propyleneimine
Cadmium compounds	Ethylene oxide	Propyleneoxide
Cadmium	Formaldehyde	Selenium and selenium
Carbon tetrachloride	Hydrazine	compounds
Chlorine	Hydrogen cyanide	Silver chromate
Chloroform	Hydrogen fluoride	Sulfuric acid
Chromic acid	Lead	Sulfuric acid (fuming)
Chromium compounds (+6)	Lead compounds	Tetrachloroethylene
Cyanide compounds	Methylene bisphenyl	Toluenediisocyanate
Dibromochloropropane	isocyanate	Trichloroethylene
Dichloroethane		

² Chemical names with CAS numbers can be found in Table 4 of this report.

¹ For this work, "hazard" includes inherent toxicity, potential for exposure through dispersal in the workplace (based on the physico-chemical properties of the chemicals, e.g., vapor pressure) and indicators of safety of use (e.g., flammability). Potential for exposure and indicators of safety do not include site-specific conditions.

Table 2: Category 2 Chemicals ³				
Acetic acid	Ferrous chloride			
Acetone	Ferrous sulfate			
Ammonium bicarbonate	Isobutyl alcohol			
n-Butyl alcohol	Methylethylketone			
sec-Butyl alcohol	Methanol			
Chromium compounds (+3)	Silver in alloy form			
Ethyl acetate	Sodium phosphate, dibasic			
Ethylene glycol	Sodium phosphate, tribasic			
Ferric chloride	Zinc in alloy form			
Ferric sulfate	Zinc borate			
Ferrous ammonium sulfate	Zinc sulfate			

Introduction

As required under the Toxics Use Reduction Act (M.G.L. c.21I) the Toxic Use Reduction Science Advisory Board serves in an advisory capacity to the Toxics Use Reduction Institute (the Institute) in the following three areas, 1) adding chemicals to or deleting chemicals from the reporting list, 2) establishing priority user segments, and 3) general advice to the Institute on other related matters. In December 1994, the Institute organized the first meeting of the TUR Science Advisory Board. The full Board is composed of eleven members with expertise in the areas of toxicology, epidemiology, medicine, worker issues, industry issues, environmental chemistry and risk assessment. A list of the members who worked on the Chemical Categorization Project is included in Appendix A.

For the past three years, the Toxics Use Reduction Science Advisory Board has assisted the Institute in preparing recommendations for the Administrative Council for delisting chemicals from the TURA Toxic and Hazardous Substance List⁴. Fourteen industry petitions requesting delisting were submitted. The Board recommended delisting in ten cases. Appendix B provides a summary of the recommendations. Throughout this petitioning process the Board has struggled with decisions which seemed to require, at least implicitly, a ranking of the relative hazards of chemicals. Delisting (or refusal to delist) particular chemicals was seen as having the potential to effect the use of one material in preference to another. Therefore, the Board has spent the last eighteen months discussing the categorization of chemicals on the list into one of the following three groups:

³Chemical names with CAS numbers can be found in Table 5 of this report.

⁴Massachusetts Department of Environmental Protection, Bureau of Waste Prevention, Toxics Use Reduction 1995 Reporting Package.

- Category 1 more hazardous chemicals
- Category 2 less hazardous chemicals
- Category 3 uncategorized chemicals which includes chemicals not reported under TURA since 1990 and chemicals reported under TURA but not categorized as more or less hazardous due to insufficient information or because the chemical was deemed to be of medium hazard.

The resulting categorized list is intended to provide guidance to companies and technical assistance providers making chemical substitution decisions, to aid in targeting technical assistance and research efforts and, ultimately, to aid in reducing overall risk to workers and the environment. It is also intended to provide information to the Institute and the other entities created under TURA, for their use in guiding the implementation of the TURA program. Categorization will not address the issue of varying risk associated with the same chemical used in different processes; this issue has been discussed frequently by the Board during the petition review processes.

Approach

To begin the Categorization project, the Board reviewed many existing models for chemical prioritization which are briefly described in Appendix C. All models, with the exception of the Swedish National Chemicals Inspectorate's system, rely entirely on a scoring system based on health and environmental data. Only one model used by the Indiana Clean Manufacturing Technology and Safe Materials Institute considers occupational safety issues which are of particular concern to the Board. In the initial stages the Board assumed that they, like the other groups, would create a model based on an algorithm using environmental, health and safety data. The Board was concerned, however, that the necessary data might not be available to accurately assess chemical hazard.

Using the criteria from existing models as a starting point, the Board chose their own set of criteria. In choosing criteria for categorizing the list, the following three items were discussed:

- the data should be generally available
- the data should be reliable
- the scheme should be defensible and understandable

Data points were discussed in the following four major areas:

- human health
- environmental
- ♦ safety
- persistence/bioaccumulation

After lengthy discussions, the Board choose the following eight criteria:

- Carcinogenicity (IARC Classification)
- Oral LD_{50}
- Reference dose (RfD)
- Threshold limit value (TLV) / time weighted average (TWA)
- Aquatic LC_{50}
- ♦ Flash point (FP)
- ◆ pH (used pKa and pKb)
- Bioconcentration factor (BCF)

These criteria are defined in Appendix D. The Board requested that the data for each chemical be provided to them for further discussion. In order to make the task less daunting, the Board decided to categorize only the 258 chemicals that had ever been reported under the Toxics Use Reduction Act⁵.

It was necessary to choose surrogate compounds for chemicals listed as groups (e.g., lead compounds). Using the surrogate choices from other chemical ranking schemes as a guide, the criteria used for this selection were as follows: most toxic member of a group, most data available, most widely used. A list of the chemical group name, the surrogate used and an explanation of the choice can be found in Appendix E. In addition, the listings for individual metals and metal compounds were defined based on similar toxicity. These definitions can be found in Appendix F.

The Institute contracted with the Tellus Institute to collect the available data. The data for each chemical, which was provided to each Board member on computer disk, is in Appendix G along with data sources and General Comments authored by the Tellus Institute concerning the collection of data. Table 3 shows the availability of data for the 258 chemicals.

Other models for chemical prioritization reviewed and considered by the Board, rely completely on algorithms which assign either a value of zero to a missing datum point, or use quantitative or qualitative structure-activity relationships⁶ to derive an estimate of the value. The Board rejected these ideas due to the lack of available data for many compounds and the crude assumptions used in algorithms to complete data sets. Instead of developing an algorithm that might be difficult to understand or could ignore known risks, the Board chose to use an expert judgment method

⁵The list of substances reportable under TURA contains the approximately 1500 substances reportable under the federal laws, EPCRA and CERCLA, with the exception of a few federally listed substances that have been delisted by the TURA Administrative Council. Only those chemicals which a TURA filer uses or processes in quantities of 25,000 pounds or more per year, or otherwise uses in the amount of 10,000 pounds or more per year at any one facility are reportable.

⁶Quantitative Structure Activity Relationship software is available (e.g., Ecosar and MicroQSAR can be obtained from the U.S. EPA).

(based on the principles of the Delphi Method). This approach was used by Polaroid in developing their chemical ranking system⁷, and it allows for incorporation of the Board members' professional experiences which is especially important for chemicals that have little or no data available. The Board members supported the expert judgment method and found it to be more satisfactory than the algorithm method

Table 3: Percent Data Available for each Criterion				
Criterion	Percent data available			
IARC Classification	40%			
LD ₅₀	55%			
RfD	38%			
TLV (TWA)	70%			
TLV (STEL)	19%			
LC ₅₀	41%			
FP	41%			
рКа	8%			
рКЬ	6%			
BCF	54%			

The Delphi Method and the Expert Judgement Method

The term Delphi Method came from a study concerning the use of expert opinion called Project Delphi performed by the Rand Corporation in the 1950s for the U. S. Air Force. This study aimed to "obtain the most reliable consensus of opinion of a group of experts."⁸ The Delphi method is appropriate when "accurate information is unavailable or expensive to obtain or evaluation models require subjective inputs to the point where they become the dominating parameters."⁹ The rationale behind the method is that "if the opinion of one expert on an uncertain point is useful, the opinion of many experts - when boiled down to a single group opinion - should be even better."¹⁰

⁹Ibid.

¹⁰Gautschi, T.F., "Delphi Method Predicts the Future," Design News, Feb. 1990, p. 414.

⁷Ahearn, J., Fatkin, H., and Schwalm, W., "Polaroid Corporation's Systematic Approach to Waste Minimization," Pollution Prevention Review, Summer 1991, pp. 257-271.

⁸Linstone, H.A., and Turoff, M., "The Delphi Method: Techniques and Applications," Addison-Wesley, Reading, Mass., 1975, pp. 3-12.

The original method uses a series of questionnaires to solicit the opinions of the experts. The results of the questionnaires are summarized by an investigator who provides feedback to the experts. A modified questionnaire is then used to obtain a second round of opinions and the process continues until consensus is reached.

The Science Advisory Board's Expert Judgement Method began with each expert choosing fifty "more hazardous chemicals" and fifty "less hazardous chemicals", subsequently named Category 1 and Category 2 respectively. Each member used their own ranking scheme based on the data, their area of expertise and personal experiences. The votes from each expert were tabulated and the chemicals were ranked by the number of expert votes received for the category.

When asked to describe the criteria used to categorize the chemicals, the following statements were made by Board members for the more hazardous list of chemicals: "data revealed at least two criteria of concern and toxicity was rated higher than flammability", "aquatic toxicity ranked highest", "focused on carcinogenicity", "looked at potential for exposure to workers", "ratio of TLV/LD₅₀". These comments illustrate the diversity of expert opinion that contributed to the creation of these categories. Commonly, the criteria used for the less hazardous list were simply opposite of the more hazardous list or "didn't raise any concerns".

Refinement

Following the initial vote, two lists were prepared of chemicals that received a number of votes for each Category. The number cutoff was determined in order to produce lists of 25-30 chemicals each for further discussion. This ended up being 4 votes for Category 2 and 5 votes for Category 1 as there was considerable consensus for many of the chemicals on the Category 1 list. Each list was then discussed chemical by chemical. In some cases, additional data were requested. In some cases, chemicals receiving one vote less than the cuttoff were discussed. For the more hazardous chemical list, the Board decided to discuss every chemical that had an IARC classification of 1, 2a or 2b. As each chemical was discussed, consensus decisions were made to put the chemical in Category 1, 2 or 3.

For chemicals that received more than one vote for each list, the Board reviewed the data that were available and discussed the chemicals at length. In all cases the discrepancy was due to either conflicting data (e.g., low TLV and high LC_{50}), the lack of data or, in the case of metals, different definitions being used by members. This exercise resulted in a complete review of all metals and metal categories to be certain that all Board members were making the same assumptions. (See Appendix F, mentioned previously.)

The Board also compared its list to other lists of hazardous chemicals such as EPA's list of Extremely Hazardous Substances, the list created by EPA's Waste Minimization Prioritization Tool, and the Swedish National Chemicals Inspectorate list of Chemical Substances Which Require Particular Attention. In each case, the Board discussed similarities and discrepancies, and concluded that their process and resulting categories were more appropriate for the purposes of the project.

The Resulting Categories

Following are the Category 1 and 2 chemicals with CAS numbers as defined by the TUR Science Advisory Board. For chemicals in Category 1, along with the chemical name are a few phrases summarizing the discussion that resulted in the chemical's placement in this category. The list of Category 3 chemicals can be found in Appendix H. It is important to note that these three categories represent only the 258 chemicals that have been reported under TURA at the time of this project. For a complete list of the approximately 1200 chemicals on the TURA List of Toxic and Hazardous Substances, please refer to the Massachusetts Department of Environmental Protection, Bureau of Waste Prevention, Toxics Use Reduction Reporting Package.

	Table 4: Category 1 Chemica	Is with CAS Numbers
-CAS Number(s)	Chemical Name	Summary
79-06-1	Acrylamide	IARC 2a, potential worker exposure problem
107-13-1	Acrylonitrile	IARC 2a, evidence of human carcinogenicity
01-00-1, 7440-38-2	Arsenic and arsenic compounds	IARC 1, acutely toxic
01-00-4, 7440-43-9	Cadmium and cadmium compounds	IARC 1
56-23-5	Carbon tetrachloride	Montreal Protocol chemical, liver toxin, suspect human carcinogen, IARC 2b
7782-50-5	Chlorine	low RfD, gas, toxic, corrosive to skin, heavier than air, stable in air, used in large quantities, toxic to aquatic organisms
67-66-3	Chloroform	low RfD, acute effects to the liver, medium bioaccumulation factor, IARC 2b
7738-94-5, 11115-74-5	Chromic acid	hexavalent chromium
01-01-2, 7440-47-3	Chromium compounds (+6 valence)	IARC 1, confirmed carcinogens
01-01-6, 143-33-9	Cyanide compounds and sodium cyanide	acutely toxic
96-12-8	Dibromodichloropropane (DBCP)	banned as a fumigant in 1977, PEL 1 ppb, IARC 2b
107-06-2	1,2-Dichloroethane (Ethylene dichloride)	acute toxicity, IARC 2b
64-67-5	Diethyl sulfate	IARC 2a, incompatible with water, highly irritating, reactive alkylating agent
68-12-2	Dimethylformamide	limited evidence of testicular cancer in humans, inadequate in animals, very soluble in water, PEL 10 ppm, easily absorbed into skin, exposures are likely to be high because of its large use, occupational hazard, highly mobile in soil
123-91-1	1,4-Dioxane	acute effects, strong skin absorber, IARC 2b

	Table 4: Category 1 Chemica	als with CAS Numbers
-CAS Number(s)	Chemical Name	Summary
106-89-8	Epichlorohydrin	IARC 2a, reportable quantity 10 pounds
75-21-8	Ethylene oxide	IARC 2a, mutagenic, reactive, eye and skin irritant, carcinogenic, and highly flammable
50-00-0	Formaldehyde	reactive, irritating, IARC 2a, acutely toxic
302-01-2	Hydrazine	eye and skin irritant, flammable, IARC 2b, TLV 10 ppb
01-02-6, 7439-92-1, 10099-74-8	Lead and lead compounds	neuro-toxic and impairs reproduction, IARC 2b
101-68-8	Methylenebis(phenylisocyanate)	
01-02-9	Nickel compounds	IARC 1 classification for nickel and nickel compounds (1990)
98-95-3	Nitrobenzene	carcinogen, causes liver damage, eye and skin irritant, smells foul, very low RFD and TLV=1 ppm
75-44-5	Phosgene	Leukocyte, severe eye, skin, mucous membrane irritant, TLV=0.1 ppm
75-55-8	Propyleneimine	very reactive, PEL=2ppm, skin absorbing, sufficient carcinogenic evidence in humans, IARC 2b
75-56-9	Propyleneoxide	TLV 20 ppm, evidence of mutagenicity, carcinogenicity, acute hazard
01-03-6, 7782-49-2	Selenium and selenium compounds	
7664-93-9, 8014-95-7	Sulfuric acid and fuming sulfuric acid	IARC 1, fuming is the actual toxic factor (due to its vapor pressure), corrosive, carcinogenic, reactive, causes lung damage
127-18-4	Tetrachloroethylene	IARC 2b, suspected carcinogen
584-84-9, 91-08-7, 26471-62-5	Toluenediisocyanate (2,4 and 2,6 and mixed isomers)	irritating to eyes, nose, skin and TLV 5 ppb, IARC 2b
79-01-6	Trichloroethylene	causes eye, skin, liver and central nervous system damage and low TLV

Table 5: Category 2 Cher	nicals with CAS Numbers
64-19-7	Acetic Acid
67-64-1	Acetone
1066-33-7	Ammonium bicarbonate
71-36-3	n-Butyl Alcohol
78-92-2	sec-Butyl Alcohol
	Chromium ³⁺ compounds
141-78-6	Ethyl Acetate
107-21-1	Ethylene Glycol
7705-08-0	Ferric chloride
10028-22-5	Ferric Sulfate
10045-89-3	Ferrous Ammonium Sulfate
7758-94-3	Ferrous Chloride
7720-78-7, 7782-63-0	Ferrous Sulfate
78-83-1	Isobutyl Alcohol
78-93-3	Methylethylketone
67-56-1	Methanol
7558-79-4, 10039-32-4, 10140-65-5	Sodium Phosphate, dibasic
7601-54-9, 7758-29-4, 7785-84-4, 10101-89-0, 10124-56-8, 10361-89-4	Sodium Phosphate, tribasic
1332-07-6	Zinc Borate
7733-02-0	Zinc Sulfate

Maintenance and Further Work

The Board realizes that the chemicals in the specific categories may change based on new data becoming available or new chemicals being used above TUR reportable threshold quantities in the Commonwealth. The Board will establish a review process whereby the Category 1 and Category 2 lists will be reviewed annually and new chemicals reported in Massachusetts will be evaluated. This review process will begin at the Board meeting following the release of Toxics Use Réduction data by the Department of Environmental Protection.

Appendix A: List of Members

David T. Williams Director, Center for Technology and Health Quincy College

James J. Ahearn Jr., Ph.D. Polaroid Corporation

Andrew F. Beliveau Environmental Protection Agency

Richard Clapp, Sc.D. Boston University School of Public Health

George M. Gray, Ph.D. Harvard School of Public Health Center for Risk Analysis

Thomas Trayers Division of Occupational Safety Occupational Hygiene Program

Lawrence H. Boise Gloucester Co., Inc.

Halina Brown, Ph. D. (ex officio) Clark University

Christine Oliver, M. D. (ex officio) Public Health Resource Group Mass. General Hospital

Appendix B: Summary of Toxics Use Reduction Science Advisory Board Recommendations

Chemical Name	Recommendation	Supplemental Information	Status or Outcome
Nickel in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote to accept recommendation. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per SAB recommendation.
Chromium in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote to accept recommendation. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per SAB recommendation.
Copper in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote to accept recommendation. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per SAB recommendation.
Manganese in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote to accept recommendation. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per SAB recommendation.
Cobalt in alloy form	delist except for aerosols (less than 50 um)	Unanimous vote to accept recommendation. Aerosols should be reported under TURA because planning for efficient use is beneficial.	Delisting petition request accepted by Admin Council per SAB recommendation.
Chromium (III) oxide	delist	Unanimous vote to accept recommendation. Chromium (III) oxide is not known to cause significant human health effects, is not known to cause significant adverse effects on the env., does not bioaccumulate and the oxidation of chromium (III) to chromium (VI) is not likely to occur.	Delisting petition request accepted by Admin Council per SAB recommendation.
Sodium hydroxide	not delist	Majority decision to accept recommendation. Decision based primarily on its potential for acute toxicity to workers. For specific applications, there may be uses of sodium hydroxide for which there is scientific justification to determine that sodium hydroxide is the least hazardous material and presents the least risk; this should be considered by the Administrative Council	Delisting petition request denied by Admin Council per SAB recommendation.
Hydroquinone	delist, except for manufacture	Unanimous vote to accept recommendation. Material has moderate to low toxicity. Recommendation to delist was made because material did not satisfy the criteria of "significant health effects"	Delisting petition request accepted by Admin Council per SAB recommendation.

Prepared 12/3/96, Updated 10/29/97

Chemical Name	Recommendation	Supplemental Information	Status or Outcome
Butyl benzyl phthalate	delist	Unanimous vote to accept recommendation. The Board recommended delisting in the absence of science to prove that butyl benzyl phthalate is estrogenic despite emerging science that suggests that this potential exists.	From a policy perspective, the Institute questioned whether the absence of knowledge is a sufficient basis to support a delisting at this time. The Admin Council denied the delisting petition.
Ethyl Acetate	not delist	Unanimous vote to accept recommendation. Recommendation based primarily on its potential for acute toxicity to workers.	Delisting petition request denied by Admin Council per SAB recommendation.
Acetic Acid	delist at conc. below 12%	Unanimous vote to accept recommendation.	Delisting petition request accepted by Admin Council per SAB recommendation.
Sodium Hypochlorite	not delist	Majority decision to accept recommendation.	Delisting petition request denied by Admin Council per SAB recommendation.
Acetone	no recommendation	Board vote was split.	Delisting request denied. Acetone will be reviewed again in one year and categorization of the list of chemicals will be evaluated.
Zinc oxide	delist	Unanimous vote to accept recommendation.	Delisting petition request accepted by Admin Council per SAB recommendation.

Prepared 12/3/96, Updated 10/29/97

Appendix C: Bibliography of Categorization/Prioritization Schemes

Davis, Gary et al., Center for Clean Products and Clean Technologies, University of Tennessee, "Chemical Hazard Evaluation for Management Strategies: A Method for Ranking and Scoring Chemicals by Potential Human Health and Environmental Impacts", EPA Document EPA/600/R-94/177, June 1994. This model uses risk-based chemical ranking and scoring combining the toxic effects of chemicals and the potential for exposure to those chemicals. The report ranks 140 TRI chemicals based on 99% of total releases. The method does not include secondary global impacts such as ozone depleting and global warming, nor does it include worker saféty. Potential uses of the methodology are: priority setting for regulatory action, for business decisions and to set priorities for pollution prevention.

Davis, Gary et al., Center for Clean Products and Clean Technologies, University of Tennessee, "Comparative Evaluation of Chemical Ranking and Scoring Methodologies", April 7, 1994.

Gray, George and Jennifer Hartwell, Harvard Center for Risk Analysis, Harvard School of Public Health, "<u>The Role of Risk in Chemical Substitution Decisions</u>," prepared for the Massachusetts Toxics Use Reduction Institute, July 1994. *Outlines a risk-based substitution decision-making* framework, the chemical substitution tree (CST). Suggests looking at both the application exposure and the disposal exposure for potential effects on the environment, workers and the public. Gives some ideas of chemical characteristics to consider and where to find relevant information. The model seeks to identify areas of potentially high risk so that companies can make informed decisions on how to reduce the risk.

Grimsted, Bradley, et al., <u>"A Multimedia Assessment Scheme to Evaluate Chemical Effects on</u> the Environment and Human Health" Pollution Prevention Review, Summer 1994, pp. 259-268. This article presents a model for calculating a common unit of measure - the Pollution Unit that allows comparisons of potential relative effects of chemicals on different environmental media. The scheme incorporates environmental and human health factors (using ambient standards and regulatory criteria) but can be adjusted to stress one over the other or may be developed to incorporate occupational standards if worker health is of primary concern. Authors boast "easy to use", "technically defensible" and "versatile" as words to describe the model.

Indiana Clean Manufacturing Technology and Sage Materials Institute, <u>Pollution Prevention</u> <u>Progress Measurement Method (3P2M)</u>, Purdue University, February 1998. This work builds on the Center for Clean Products and Clean Technologies algorithm to include factors for worker exposure and atmospheric hazards. This model does not include releases to the environment as did the Clean Products work. It has an option for inputting number of pounds of a chemical used in the workplace. The worker exposure component has three parts: health effects (chronic and acute), routes of exposure (vapor pressure, oral, skin, dust/mist) and safety (flammability, reactivity, corrosivity). For carcinogenicity, the most protective rating of EPA, ACGIH and IARC was used. The acute hazard value is based on the short term exposure limit (STEL); if an STEL does not exist, the score is 0. For oral exposure, the only compounds with scores other than zero are lead compounds. Massachusetts Toxics Use Reduction Institute, "<u>Blanket Wash Technology Study: An Evaluation</u> of Commercially Available Blanket Washes," Technical Report No. 16, 1994. This study gives comparative information on the performance, environmental, health and safety characteristics of blanket washes commonly used in sheetfed offset lithography. Each attribute was given a good, fair or poor score. The non-performance attributes scored included VOC content, flash point, health hazard and potential regulatory impact. For determining a score for the health hazard, mixtures were given the highest score of any ingredient and data were obtained from REPROTEXT. For determining the potential regulatory impact, chemicals were given scores based on how many times they appeared on nineteen regulated chemical lists.

Swedish National Chemicals Inspectorate, "<u>Observation List of Chemical Substances Which</u> <u>Require Particular Attention</u>," Printgraf AB, Stockholm, Sweden, January 997. This work was done to guide users of chemicals to pay particular attention to the use of chemicals on the Observation List which contains 199 chemicals. A substance was placed on the list if it met any of an established list of criteria indicating environmental or health hazards. These criteria include bioaccumulation, aquatic toxicity, ozone depletion, acute toxicity, sensitizer, chronic toxicity, neurotoxicity, reproductive toxicity and carcinogenicity.

Swedish National Chemicals Inspectorate, "Selecting Multiproblem Chemicals for Risk Reduction." This work began with 7000 chemicals which was a combination of 70 national and international lists of chemicals hazardous to human health or the environment. In Step I, the list was narrowed to 500 chemicals that appeared on several lists. In Step II, the list was narrowed to 100 chemicals using 18 criteria of equal weight in the categories of environmental properties, health properties, and exposure potential. If no data was available, the criteria was not used. From this list of 100 chemicals, 45 were chosen based on additional data and use patterns in Sweden. Finally, 27 chemicals were chosen as candidates for risk reduction by a panel of experts who used the available data and member's experience and knowledge.

Tiley, Jaimie, <u>"Solvent Substitution Methodology using Multiattribute Utility Theory and the</u> <u>Analytical Hierarchical Process</u>", Department of the Air Force, Air Force Institute of Technology, Wright-Patterson Air Force Base, OH. *This thesis presents a multicriteria decision making methodology for ranking alternatives to solvent cleaning. It compares Multiattribute Utility Theory and the Analytical Hierarchical Process. The cleaning situation studied is general cleaning of aircraft engine components. There were problems associated with both decision models including independence constraints and scaling issues. The author used group decision making scoring (1-7) in four areas: environmental impact, health/safety, process compatibility, cleaning effectiveness. Important attributes within each category were chosen by survey. Interesting to note which attributes were chosen in the environmental impact and health/safety categories (p 46.)*

US Environmental Protection Agency, "<u>Waste Minimization Prioritization Tool</u>," EPA 530-R-97-019, June 1997. This work began with the adoption and modification of earlier work on the Use Cluster Scoring System. The Tool uses persistence, bioaccumulation and toxicity (human cancer, human non-cancer and ecological) characteristics for chemical risk screening. The mass of a chemical can be input into the tool's software. 900 chemicals were scored due to the

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availability of data. Partial chemical data is available for an additional 3800 chemicals which were not scored. If data did not exist, the chemical is not scored. The tool ignores acute effects, including those to workers.

Wolf, Katy, "The Generic Classification System: A Simplified Approach to Selecting

<u>Alternatives to Chlorinated Solvents</u>" Pollution Prevention Review, Winter 1993-94, p 15-29. The author sets up a generic classification system for choosing alternative to a chlorinated solvents. The properties/classifications of PEL, VOC, HAP, flash point, evaporation rate, solvent strength, ozone depleting potential, global warming potential and toxicity are covered. Good reference for data on the available solvent alternatives. Methodology is practical but very specific to solvents alternatives.

Working Group of Accelerated Reduction/Elimination of Toxics (ARET), "<u>Environmental</u> <u>Leaders - Voluntary Commitments to Action on Toxics through ARET</u>," Ontario, March 1995. This work began with 2000 substances from the Chemical Evaluation Search and Retrieval System. Approximately 500 of these substance had sufficient information to screen them for the ARET list. The criteria were chosen in the areas of toxicity, persistence and bioaccumulation. The toxicity criteria were in the following seven groups: acute lethality, chronic toxicity nonmammals, chronic toxicity plants, chronic toxicity mammals, teratogenicity, carcinogenicity, genotoxicity.

Appendix D: Criterion Definitions

<u>Oral LD</u>₅₀: A single calculated dose of a substance administered through food or gavage (tube feeding) in mg per kg of body weight, which kills 50% of a group of test animals within 14 days. A lower LD₅₀ indicates a more toxic substance.

<u>Reference Dose (RfD)</u>: An estimate of the daily exposure level for the human population that is likely to be without an appreciable risk of adverse effects over a lifetime. RfDs are often estimated from the highest dose at which no adverse effects are observed in animals, the No Observed Adverse Effects Level (NOAEL). The Environmental Protection Agency has defined RfD's for a number of chemicals.

<u>Carcinogen</u>: International Agency for Research on Cancer (IARC) uses the term "carcinogen to denote an agent that is capable of increasing the incidence of malignant meoplasms; the induction of benign neoplasms may in some circumstances contribute to the judgement that an agent is carcinogenic..."

IARC Classification. Carcinogens are rated in 1 of 5 groups: (1) Group 1 - the agent is carcinogenic to humans; (2) Group 2A - the agent is probably carcinogenic to humans; (3) Group 2B - the agent is possible carcinogenic to humans; (4) Group 3 - the agent is not classifiable as to its carcinogenicity to humans (when agents cannot be placed in any other group); and (5) Group 4 - the agent is probably not carcinogenic to humans.

<u>TLV (Threshold Limit Value)</u>: Published by the American Conference of Governmental Industrial Hygienists (ACGIH), defined as airborne concentrations under which it is believed that nearly all workers may be repeatedly exposed day after day without adverse effects. TLV's are generally established on a consensus basis; as such, some workers may be affected at or below these limits due to unusual susceptibility or pre-existing conditions. A lower TLV indicates a more toxic substance.

<u>Aquatic LC₅₀</u>: The concentration of a chemical, in water, that causes death in 50% of the fish tested. Aquatic LC₅₀ can be calculated for both freshwater and saltwater fish (and sometimes for other aquatic organisms).

<u>Bioconcentration</u>: Describes the tendency for a chemical to accumulate in biological systems, and more specifically the ability of a substance to accumulate in the tissues of organisms. Bioconcentration is a function of the physicochemical properties of a chemical, especially the chemical's lipid solubility (solubility in fat). Two parameters most frequently used to express bioconcentration are the octanol-water partition coefficient (K_{ow}) and the Bioconcentration factor (BCF).

[•] Bioconcentration factor is the ratio of the concentration of a chemical in an organism to its concentration in the test medium or environment, typically water, at steady-state conditions. This factor is a measure of the chemical's ability to bioaccumulate K_{ow} is defined as the ratio of a chemical's concentration in the octanol phase to its concentration in the aqueous phase of a two-phase 1-octanol/water system at equilibrium. In other words, it represents the distribution tendency of organic chemicals between organic and aqueous phases. As lipid soluble chemicals are generally also soluble in solvents such as octanol and are relatively insoluble in water, K_{ow} can be used to predict the bioconcentration factor. A low log K_{ow} value is considered hydrophilic and has a low fat solubility and high water solubility. K_{ow} is generally expressed in log units.

<u>Flash point</u>: The temperature at which material gives off sufficient vapor to form an ignitable mixture with the air near the surface of the material. The lower the flash point, the more probability an explosion could occur under normal working conditions.

<u>pH</u>: A logarithmic index for the hydrogen ion concentration in an aqueous solution. A pH below 7 indicates acidity, and one above 7 alkalinity (at 25C). The pH scale ranges from 0-14, with extreme values representing a more corrosive aqueous solution. Values closest to 7 represent the lowest hazard.

Appendix E: Surrogate Chemicals

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Mass. #	CAS #	Chemical Group Name	Suggested Surrogate	Notes
01-00-0	1309-64-4	Antimony compounds	Diantimony trioxide	Tennessee surrogate
01-00-1	1303-28-2	Arsenic compounds	Arsenic pentoxide	Tennessee surrogate
01-00-2	10361-37-2	Barium compounds	Barium chloride	Tennessee surrogate
. 01-00-4	10108-64-2	Cadmium compounds	Cadmium chloride	Tennessee surrogate
01-01-2	1333-82-0	Chromium Compounds	Chromium oxide	Tennessee surrogate
01-01-3	7646-79-9	Cobalt compounds	Cobalt chloride	Tennessee surrogate
01-01-5	7758-98-7	Copper compounds	Copper sulfate	Tennessee surrogate
01-01-6	143-33-9	Cyanide compounds	Sodium cyanide	Most widely used, most toxic, most data available
01-02-2	110-80-5	Glycol ethers	Glycolmonoethylether	Most common, most data available
01-02-6	7758-95-4	Lead compounds	Lead chloride	Tennessee surrogate
01-02-7	1344-43-0	Manganese compounds	Manganese oxide	Tennessee surrogate
01-02-9	37211-05-5 373-02-4 6018-89-9	Nickel and compounds	Nickel chloride Nickel acetate Nickel acetate tetrahydrate	Tennessee surrogate for all except mammalian oral toxicity For mammalian oral toxicity due to availability of data Choose specific Nickel acetate with the most data
01-03-3	117-81-7	Phthalate esters	Diethylhexylphthalate	Most common, most data available
01-03-6	7446-08-4 7783-00-8 7488-56-4 7783-79-1 12033-59-9 14832-90-7 57-12-5	Selenium and compounds	Selenium IV dioxide Selenium IV disulfide Selenium hexallouride Selenium nitride Selenium oxide Selenium	Tellus choose chemical in this group that has the most data

Mass. #	CAS#	Chemical Group Name	Suggested Surrogate	Notes
01-03-7	7783-90-6 7761-88-8 7783-91-7 7784-01-2 506-64-9	Silver and compounds	Silver chloride Silver nitrate Silver chlorite Silver chromate Silver cyanide	Tellus choose chemical in this group that has the most data
01-03-9	1314-13-2 7733-02-0	Zinc and compounds	Zinc oxide Zinc sulfate	Tennessee surrogate for all except fish toxicity For fish toxicity b/c Zinc oxide not soluble in water
01-09-0	7761-88-8	Nitrate compounds	Zinc nitrate	Note: water dissociable nitrate compounds reportable only in aqueous solutions on SARA 313. Silver nitrate most soluble.

Notes: For Nickel acetate, Selenium compounds and Silver compounds, choose the specific chemical based on the availability for the most data. Glycomonoethyl and Diethylhexylphthalate are listed separately also.

Appendix F: Metals

At the November, 1997 meeting of the Science Advisory Board, members proposed groupings for metals and metal compounds to represent similarities in metal toxicities. If the base metal is a category by itself (e.g., copper), that particular category represents the toxicity of the metal itself (e.g., metallic copper). If the base metal is in a category with other metal compounds (e.g., lead and lead compounds), it was the opinion of the group that the metal toxicities of the base metal and the metal compounds were basically similar. For categories that are comprised of more than one compound, the CAS number for the group is given.

Base Metal		Metal Categories	CAS#
Aluminum	1	Aluminum, Aluminum oxide	7429-90-5
	2	Aluminum sulfate	
Antimony	1	Antimony, Antimony cmpds, Antimony trioxide	1309-64-4
Arsenic	1	Arsenic, Arsenic compounds	1303-28-2
Barium	1	Barium	
	2	Barium compounds	
Cadmium	1	Cadmium, Cadmium compounds	10108-64-
Character	1	Chromium, Chromium and cmpds [valence 0 and 3]	7440-47-3
Chromium	2	Hexavalent chromium [valence 6]	1333-82-0
Cobalt	1	Cobalt	
	2	Cobalt compounds	
Copper	1	Copper	
	2	Copper compounds	
Iron	1	Ferric chloride, Ferric sulfate, Ferrous ammonium sulfate, Ferrous chloride, Ferrous sulfate	7705-08-0

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Base Metal		Metal Categories	CAS#
Lead ·	1	Lead, Lead chromate, Lead compounds, Lead nitrate	7758-97-6
Manganese	1	Manganese, Manganese cmpds, Potassium permanganate	7439-96-5
Nickel	1	Nickel	
	2	Nickel acetate, Nickel acetate tetrahydrate, Nickel and compounds	37211-05-
Potassium	1	Potassium hydroxide	
Selenium	. 1	Selenium, Selenium & cmpds, Selenium cmpd, Selenium dioxide, Selenium hexaflouride, Selenium nitride, Selenium oxide	7488-56-4
	1	Silver	
	2	Silver and compounds	
Silver*	3	Silver chlorite	
	4	Silver chromate	
	5	Silver cyanide	
	6	Silver nitrate	
*For the Silver presence of the	comp ose con	ounds, it was noted that the Silver chloride is not very soluble and that the Chromate and Cyanide compounds are more toxic due moments, not the silver.	to the
	1	Sodium	
	2	Sodium bichromate	
Sodium	3	Sodium bisulfite	
	4	Sodium cyanide	
	5	Sodium dimethyldithiocarbamate	
	6	Sodium dodecylbenzenesulfonate	

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Base Metal		Metal Categories	CAS#
-	7	Sodium fluoride	
	8	Sodium hydrosulfide	
	9	Sodium hydroxide	
	10	Sodium hypochlorite	
	11	Sodium methylate	
	12	Sodium nitrite	
-	13	Sodium phosphate, dibasic	
i	14	Sodium phosphate, tribasic	
	15	Sodium phosphate, tribasic dodecahydrate	
-	16	Sodium phosphate, tribasic anhydrous	-
	1	Zinc	
	2	Zinc and compounds	
Zinc	3	Zinc oxide fume	
	4	Zinc sulfate	
	5	Zinc ammonium chloride	-
	6	Zinc borate	
	7	Zinc sulfate	

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Appendix G: Data, Collection and Sources

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MASTER DATA FILE

CLAS Number Control Name (Hall C) RE0 mg/ms/s (PC) (PC) (PC) (PC) Aquatic LCS0 notes (Do0 poles) TU Notes 0000777-06.9 (Lar) CLOROT FACUE (Lar)]]				TLV(TWA in				рКа	рК	LD50,	c LC50			
000177.006 (1, HUCHCORO 1 FLUGROFTWARE 0.0736 0.0736 0.0736 000078.006 (HEPML2, PERRODNE) 0.0736 0.0736 0.0736 000078.006 (HEPML2, PERRODNE) 0.0736 1.07 Rei, Territor 000078.006 (HEPML2, PERRODNE) 0.0736 1.07 Rei, Territor 000078.006 (HEPML2, PERRODNE) 0.0736 1.07 Rei, Territor 000078.006 (HEPML2, PERRODNE) 0.07 1.06 1.06 1.07 000078.006 (HEPML2, PERRODNE) 0.07 1.06 1.06 1.07 1.06 1.07 1.07 1.06 1.07 <th>CAS Number</th> <th>Chemical Name</th> <th>IARC</th> <th>RFD</th> <th>mg/m^3)</th> <th>FP (C)</th> <th>BCF</th> <th>рКа</th> <th>(ca)</th> <th>notes</th> <th>рТ</th> <th>, pT</th> <th>Aquatic LC50 notes</th> <th>LD50 notes</th> <th>TLV Notes</th>	CAS Number	Chemical Name	IARC	RFD	mg/m^3)	FP (C)	BCF	рКа	(ca)	notes	рТ	, pT	Aquatic LC50 notes	LD50 notes	TLV Notes
0000276-00-142 PRINCIPUANE 0.0738 0.0738 0.0738 0000276-00-122 PRINCIPUANE 7 86.7 0.0716 Ref. TenToc 0000276-00-122 PRINCIPUANE 28 no TWA 46.0 0.12 Bohr, FIAL, State, TenToc 0000276-00-122 PROFINATE 28 no TWA 46.0 0.12 Bohr, FIAL, State, TenToc 0000276-00-122 PROFINATE 28 no TWA 46.0 0.12 Bohr, FIAL, State, TenToc 0000276-00-122 PROFINATE 0.01 118 17 0.090 0.02 FIAL, TenToc 00000276-00-122 PROFINATELE 0.01 118 17 0.090 0.02 1.01 Bohr, FIAL, State, TenToc Mouse 00000276-00-122 PROFINATELE 0.01 47 80.080 0.02 1.01 Bohr, FIAL, State, TenToc Mouse Bohr, FIAL,	0001717-00-6	1,1-DICHLORO - 1 FLUOROETHANE								1	<u> </u>]		<u></u>	
000064-8-1122.4-TRIMETHY DEPTRATE 7 94.7 94.7 PRI Territor PRI Territor 0000671-85-33 DODO2-POPWINT 28 POTWA 40.228 POTWA PRI Territor PRI Territor 0000071-85-32 DODO2-POPWINT 28 POTWA 40.228 POTWA PRI Territor Mouse 0000071-85-32 DODO2-POPWINT 28 POTWA 40.228 POTWA PRI Territor Mouse 0000071-85-32 DODO2-POPWINT 28 POTWA POTWA POTWA PRI Territor Mouse 0000074-10-72 DOTWA POTWA POTWA POTWA POTWA Mouse Mouse POTWA Mouse POTWA Mouse POTWA Mouse POTWA POTWA POTWA POTWA Mouse POTWA	0000872-50-4	1-METHYL-2-PYRROLIDONE					0.0738					1		{	
0000258-93-33D(CHLORDCRE/ZDINE	0000540-84-1	2,2,4-TRIMETHYLPENTANE				-7	848.7			1	1	1		1	
0655405-35,93-0002-24PROPEYV, Image: Signal Si	0000612-83-9	3,3'DICHLOROBENZIDINE									-1.07			Rat, TerraTox	
00000763-07-04/CETALGENYDE 28 40 0.402 1.44 0.75 980, FFM, Sale, TerraTox Rat, TerraTox 0000064-97-04/CETC ACID 22 46 0.448 1.42 0.94 1.23 987, FFM, Sale, TerraTox Mouse, 0000163-84-74/CETC AKIVGRUE 0.01 781 170 0.029 FAIL TerraTox Mouse, 0000163-84-74/CETONER 0.01 781 60 0.068 -1.77 1.51 987, FFM, Sale, TerraTox Mouse, 0000076-84-74/CETONER MUNDLE 0.01 462 4600 -0.78 -0.02 1.61 987, FFM, Sale, TerraTox Mouse, skin 0000076-84-74/CETONERNUE 24 0 0.03 0.031 -0.13 -0.19 Mouse, skin 0000076-74-74/CETONERNUE 24 6 4.60 1.623 6.47 989, FFM, Sale, TerraTox Mouse, skin 0000076-74-74/CETONERNUE 24 6 6.671 0.03 6.47 989, FFM, Sale, TerraTox Mouse, skin 0000076-74-74/CANUARIDE 24	0055406-53-6	3-IODO-2-PROPYNYL									Ì	1	· · · · · · · · · · · · · · · · · · ·		
000006419-7/ACETIC-ARD 12 54 0.11 4.78 -0.98 -0.12 Sept. FFM. Static, TerraTox Mouse, 0000067-64-1/ACETICANIPORDE 0.11 78 -17 0.099 -1.71 2.15 Sept. FFM. Static, TerraTox Mouse, 0000075-64-1/ACETICANIPORDE 0.01 67 5 0.0395 -0.78 -0.15 Sept. FFM. Static, TerraTox Mouse, sept. FFM. Sta	0000075-07-0	ACETALDEHYDE	2B		no TWA	-40	0.4202			1	-1.64	0.15	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000028-47,ACETICA.NHYCRIDE 0.1 17.4 1.2.4 Ret. TerraTox 0000078-0-6,ACETORE 0.1 1188 1.7 2.5 Sehr. KM Static, TerraTox Mouse, 0000078-0-6,ACETORE 0.01 67 5 0.0303 -0.7 2.5 Sehr. KM Static, TerraTox Mouse, 0000078-0-1,ACETORE 0.01 68 6.7 0.33 0.031 0.011 Control Mouse, etin 0000078-0-1,ACETORE 0.05 6.5 5 0.033 0.031 0.011 Control Mouse, etin 0000078-0-1,ACETORE 0.5 5 0.033 0.031 0.031 Control Mouse, etin 0000078-0-1,ACETORE 0.5 5 0.033 0.031 Control Mouse, etin<	0000064-19-7	ACETIC ACID			25	40	0.1146	4.75		[-0.94	-0.12	96hr, FHM, Static, TerraTox	Mouse,	
0000076-4-1 ACETONIE 0.1 1188 .17 0.098 .1.7 1.21 Gen, FMA, State, TerraTox Mouse 0000076-4-1/ACETONITRUE 0.1 467 8 0.038 .0.22 1.61 667, FMA, State, TerraTox Mouse 0000074-0-7/ACETONITRUE 0.1 467 0.13 666, FMA, State, TerraTox Mouse skin 0000074-0-7/ACETONITRUE 0.5 5.5 54 0.33 FRA, TerraTox Mouse skin 0000074-0-7/ACETONITRUE 2.4 0.276 0.23 647, FMA, State, TerraTox Mouse skin 000017-3-1/ACETONITRUE 2.4 0.276 0.276 0.5 667, FHA, State, TerraTox Mouse skin 000170-53 1.41 0.276	0000108-24-7	ACETIC ANHYDRIDE			21	54	0.0486			ĺ	-1.24			Rat, TerraTox	·······
0000078-65-8/ACCTOMENDE 0.01 67 5 0.0003 0.62 1.61 96hr, FHM, State, TerraTox Mouse, 0000078-65-MACRTOREMONE 0.03 0.031 0.0319 -0.13 64hr, FHM, State, TerraTox Mouse, skin 0000078-65-MARVICACID 0.53 54 -0.33 Ret, TerraTox Mouse, skin 0000173-61-MARVICACID 0.5 54 -0.33 Ret, TerraTox Mouse, skin 0000174-61-MARVICACID 0.5 59 54 -0.28 0.42 Skin, FHM, State, TerraTox Mouse, skin 0000174-61-MARVICARID 1 3 24 677 -0.28 64r, FHM, State, TerraTox Mouse, skin 000174-61-MARVICARID 1 -	0000067-64-1	ACETONE		0.1	1188	-17	0.099				-1.71	-2.15	96hr, FHM, Static, TerraTox	Mouse,	······
0000098-06-2 ACTOPFIENONE 0.1 469 82/ 49000 0.79 0.11 DBR, FHM, State, TerraTox Mouse. 0000079-0-7)-07/ACRYLLACDID 0.5 5.9 54 0.33 FRI, FHM, State, TerraTox Mouse. ekin 0000171-31-ACRYLLACDID 0.5 5.9 54 0.33 FRI, TerraTox Mouse. ekin 0000171-31-ACRYLLACDID 0 4.5 0 0.276 0.74 0.38 GRY, FHM, State, TerraTox Mouse. ekin 0000170-51-ALUNNILMONDE 1 2.8 4.677 0.37 0.74 0.5 GRY, FHM, State, TerraTox Mouse. ekin 000134-33-ALUNNILMA 5 2 4.72 0.74 0.6 GRY, FHM, State, TerraTox Mouse. multiple TLVs, wabling turnes and purp powders 000134-33-ALUNNILMANULFATE 10 - - 1.26 0.69 Serv, FHM, State, TerraTox Mouse. multiple TLVs, wabling turnes and purp powders 0.01214-53-4 AMOUNUMUNILLORDID 0.001 0.01215-01-4 A.726 0.68 0.69 Ser	0000075-05-8	ACETONITRILE	1	.0.01	67	5	0.0803				-0.82	-1.61	96hr, FHM, Static, TerraTox	Mouse,	
0000079-06-4 ACRYLANDE C 24 0 0.03 0.0319 -0.18 0.15 96hr, FHM, Static, TerraTox Mouse, Mouse ekin 0000071-3-14,ACRYLOACID 0.5 5.9 54 0.33 FRATERTOX Sin 000017-3-14,ACRYLOACID 1 169 1536 442 0.33 FRATERTOX Mouse, sin 000017-05-14,ALVLCH-LORIDE 3 228 4.677 0.74 0.5 96hr, FHM, Static, TerraToX Mouse, sin 00001740-49,ALUMINUM 5 10 0.74 0.74 0.5 96hr, FHM, Static, TerraToX Mouse, sin 0001740-49,ALUMINUMSCHEF 10 -1.26 Mouse, Juminum soluble saits 0.00174-0-44,MANNA 0.017 0.74 0.95 96hr, FHM, Static, TerraToX Mouse, Juminum soluble saits 0001740-43-AMANNA 17 0.266 4.75 CRC 0.98 96hr, FHM, Static, TerraToX Mouse, Juminum soluble saits 0001740-43-AMANNUMMELORIDE 10 -1.26 -1.26 Mouse, Juminum soluble saits 0.001112-5.5.5.5.5 0.67 96h	0000098-86-2	ACETOPHENONE		0.1	49	82	4.9805				-0.79	-0.13	96hr, FHM, Static, TerraTox	Mouse.	
0000079-10-/IACRYLCACID 0.5 5.9 64 0.33 Eat, TerraTox Behr 0000173-0-4/ACRYLONTRILE 2A 0 4.7 500r, FHM, Static, TerraTox Mouse, skin 0000173-0-4/ADPIC ACID 3 196 0.133 4.42 0.47 Sonr, FHM, Static, TerraTox Mouse, skin 000173-0-4/AUVICHLCRIDE 3 24 4.77 0.74 0.5 Sonr, FHM, Static, TerraTox Mouse, umdified TLVs, welding Amees and pyro powders 000134-3-4/AUMINUMSULFATE 10 - - 1.28 multiple TLVs, welding Amees and pyro powders 000134-3-3/AUMINUMSICARCONATE 17 0.265 4.76 CRC 0.98 Sonr, FHM, Static, TerraTox Mouse, Juminum soluble safis 000134-3-3/AUMONUMSICARCONATE 10 - - 0.99 Sonr, FHM, Static, TerraTox TerraTox Mouse, Juminum soluble safis 000134-3-4/AUMONUMCHLORIDE 10 - - - 0.99 Sonr, FHM, Static, TerraTox Mouse, Juminum soluble congost 0001332-43/AUMONUMCHLORIDE 10 -	0000079-06-1	ACRYLAMIDE	2A	0	0.03		0.0319				-0.18	-0.19	96hr, FHM, Static, TerraTox	Mouse.	skin
0000173-13_LACPYLONITRILE 2A 0 274 90Hr, FHM, Static, TerraTox Mouse, skin 0000174-05_LALTYLCHORIDE 0 3 228 4.6771 -0.74 0.5 96Hr, FHM, Static, TerraTox Mouse, skin 000174-05_LALLYLCHORIDE 0 3 228 4.6771 -0.74 0.5 96Hr, FHM, Static, TerraTox Mouse, skin 000174-05_LALLMINUMOXUDE 10 - -1.28 multiple TUX, weighing mess and pyro powders 0001764-05_LALMINUMOXUDE 12 -1.28 - Mouse, skin 0001764-13_LAMINONUMELLORIDE 17 0.265 4.75 CRC 0.59 96Hr, FHM, Static, TerraTox Mouse, Juninom soluble salts 0001736-12_AMMONIMAHUORIDE 10 - - - - - - - Mouse, Juninom soluble salts 0001736-12_AMMONIMAHUORIDE 10 - - - - - - - - - - - - - - - - - <td>0000079-10-7</td> <td>ACRYLICACID</td> <td></td> <td>0.5</td> <td>5.9</td> <td>54</td> <td></td> <td></td> <td></td> <td></td> <td>0.33</td> <td></td> <td></td> <td>Rat. TerraTox</td> <td>skin</td>	0000079-10-7	ACRYLICACID		0.5	5.9	54					0.33			Rat. TerraTox	skin
000129-0-49_ADIPIC ACID 5 166 0.1566 4.42 0.74 0.5 Self, TerraTox Mouse; 000129-0-49_ALLMINUM 5 4.6771 0.74 0.5 56/rr, FHM, Static, TerraTox Mouse; 000134-0-33_LLMINUMSULFATE 2 - -1.28 multiple TLVs, welding formes and pyro powders 0001664-01_ALMINUMSULFATE 2 - -1.28 Mouse; Mouse; 0001664-01_AMMONIABICAREONATE 17 0.265 4.75 CRC 0.98 96/rr, FHM, Static, TerraTox Mouse; 000134-05_AMMONIABICAREONATE 17 0.265 4.75 CRC 0.98 96/rr, FHM, Static, TerraTox Mouse; 001314-07_AMMONIABICHORDE -	0000107-13-1	ACRYLONITRILE	2A	0	4.3	0	0.2764				0.29	0.47	96hr, FHM, Static, TerraTox	Mouse.	skin
00001742-51_ALLYLCH CRIDE 3 28 4.6771 -0.74 0.6 Opint, FHM, Static, TerraTox Mouse, Implicing Turnes and providers of the state of the s	0000124-04-9	ADIPIC ACID			5	196	0.1936	4.42						i	
000728-90-5/ALUMINUM 5 multiple TLVs, welding furnes and pyro powders 0001344-28.1/LUMINUMSULEATE 1 0 -1.28 Mouse, luminum soluble saits 0007684-17, MAMONIA 17 0.265 4.75 CRC 0.98 98hr, FHM, Static, TerraTox Mouse, luminum soluble saits 0001342-82, MAMONIUMEICARBONATE 10 - <	0000107-05-1	ALLYLCHLORIDE			3	-28	4.6771				-0.74	0.5	96hr, FHM, Static, TerraTox	Mouse.	
0001344-28-1, ALUMINUMONICE 10 128 00010043-013, ALUMINUMSULFATE 128 1.28 Mouse, Juminum soluble salts 00010643-17, AMMONIUMBICAREONATE 17 0.265 4.75 CRC 0.89 96hr, FHM, Static, TerraTox Mouse, Juminum soluble salts 0001664-37, AMMONIUMBICAREONATE 10 4.75 CRC 0.99 96hr, FHM, Static, TerraTox fume 00012125-01-8, AMMONIUMELIDORIDE 10 4.75 CRC 0.99 96hr, FHM, Static, TerraTox fume 0001236-21-6, AMMONIUMELIDORIDE 10 4.72 4.66 0.87 0.99 96hr, FHM, Static, TerraTox monologues 0001305-21-6, AMMONIUMELIDORIDE 10 4.72 4.66 0.87 0.99 96hr, FHM, Static, TerraTox Rat, TerraTox 0000126-23-3, AMILINE 7.6 70 2.5704 4.66 0.87 99hr, FHM, Static, TerraTox Rat, TerraTox Vinci. Anonologues 00013026-24-4, ANTIMONY COMPOUNDS 28 0 0.5 4.66 0.87 9hr, FHM, Static, TerraTox Antimony TLY 0001302-24-24, RASENIC <td>0007429-90-5</td> <td>ALUMINUM</td> <td></td> <td></td> <td>5</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>multiple TL</td> <td>Vs, welding fun</td> <td>es and pyro powders</td>	0007429-90-5	ALUMINUM			5								multiple TL	Vs, welding fun	es and pyro powders
001003-01-3] ALUMINUMSULFATE 2 1.26 Mouse_luminum soluble saits 0001066-33-7_AMMONIUMBICAREDNATE 17 0.265 4.75 CRC 0.89 96hr, FHM, Static, TerraTox 0001066-33-7_AMMONIUMBICAREDNATE 10 2 <td>0001344-28-1</td> <td>ALUMINUMOXIDE</td> <td></td> <td>Ī</td> <td>10</td> <td></td> <td></td> <td>Ì</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>_</td> <td></td>	0001344-28-1	ALUMINUMOXIDE		Ī	10			Ì						_	
0007644-17, AMMONIA 17 0.265 4.76 CRC 0.88 96hr, FHM, Static, TerraTox 0001064-37, AMMONIUMBICARDONATE 10 10 10 10 10 00012125-01-8 AMMONIUMBICARDONATE 10 10 10 10 10 0012125-01-8 AMMONIUMPLORIDE 10 10 10 10 10 10 10 0012125-01-8 AMMONIUMPLORIDE 10 1.124 0.99 96hr, FHM, Static, TerraTox 10 0001305-24-4 AMMONIUMPLORAUDE 1.124 Rat, TerraTox 10 1.124 Rat, TerraTox 10 0001305-24-4 AMMONIUMPLORAUDE 28 0 0.5 10 Amfmony TLV 0001309-44-4 ANTIMONY COMPOUNDS 28 0 0.5 10 Amfmony TLV 0001309-44-4 ANTIMONY FOROUDE 28 0 0.61 10 Amfmony TLV 0001309-44-4 ANTIMONY FOROUDE 28 0 0.61 10 Amfmony TLV 0001309-44-4	0010043-01-3	ALUMINUMSULFATE	1		2			1			-1.26			Mouse.	luminum soluble salts
0001063-37-7AMMONIUMBICARBONATE Image: Control of the second	0007664-41-7	AMMONIA			17		0.265		4.75	CRC		0.98	96hr, FHM, Static, TerraTox		
0001314-49-7/AMMONUMBIFLUORIDE 10 10 6012125-01-8 AMMONUMEUCRIDE 10 6012125-01-8 AMMONUMFLUORIDE 10	0001066-33-7	AMMONIUMBICARBONATE	1	1											
0012125-02-9 AMMONUMCHLORIDE 10 0.99 96hr, FHM, Static, TerraTox fume 0001338-21-6 AMMONUMHYDROXIDE -0.99 96hr, FHM, Static, TerraTox -	0001341-49-7	AMMONIUMBIFLUORIDE	1												
0012122-01-9[AMMONIUMFLUCRDE -0.99 96hr, FHM, Static, TerraTox 0001332-16 AMMONIUMFUCRDXDE 10 -1.24 Rat, TerraTox 0007773-06-0[AMMONIUMFUDROXDE 76 70 2.5704 4.69 -0.99 96hr, FHM, Static, TerraTox 0007403-56-0[ANTIMONY 0 0.5 -1.24 Rat, TerraTox Rat orago 0001306-64-4[ANTIMONY COMPOUNDS 2B 0 0.5 - - Antimony TLV 0001308-64-4[ANTIMONY RICKIDE 2B 0 0.5 - - - Antimony TLV 0001308-64-4[ANTIMONY RICKIDE 2B 0 0.5 - - - Antimony TLV 0001308-64-4[ANTIMONY RICKIDE 0 0.01 - - - Antimony TLV 0001308-64-4[ANTIMONY RICKIDE 0 0.01 - - - Antimony TLV 0001308-64-4[ANTIMONY RICKIDE 0 0.01 - - - - - Arteminory TLV Antimony TLV - - - - - - </td <td>0012125-02-9</td> <td>AMMONIUMCHLORIDE</td> <td>1</td> <td></td> <td>10</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>fume</td>	0012125-02-9	AMMONIUMCHLORIDE	1		10										fume
0001338:21-6] AMMONIUMHYDROXIDE 10 -1.24 Rat, TerraTox, Rat, TerraTox, LV incl. homologues, and compounds 0000062:53-3] ANILINE 7.6 70 2.5704 4.69 -0.67 96hr, FHM, Static, TerraTox, Rat, TerraTox, LV incl. homologues, and compounds 0001308-64:4] ANTIMONY COMPOUNDS 28 0 0.5 -	0012125-01-8	AMMONIUMFLUORIDE										-0.99	96hr, FHM Static TerraTox		
0007773-06-0/AMONUMSULFAMATE 10 -1.24 Rat, TerraTox 0000062-53-3 (ANLINE) 7.6 70 2.5704 4.69 0.67 96hr, FHM, Static, TerraTox Rat, TerraTox 0001302-64-4 (ANTIMONY 0 0.5 - </td <td>0001336-21-6</td> <td>AMMONIUMHYDROXIDE</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>1</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	0001336-21-6	AMMONIUMHYDROXIDE						1							
0000062-33-3[ANILINE 7.6 70 2.5704 4.69 -0.67 0 96hr, FHM, Static, TerraTox Rat, TerraTox LV incl. homologues 0007440-36-0[ANTIMONY 0 0.5 Antimony TCV 0001309-64-4[ANTIMONY COMPOUNDS 2B 0 0.5 Antimony TLV 0001309-64-4[ANTIMONY TRIOXIDE 2B 0 0.5 Antimony TLV 0001309-64-4[ANTIMONY TRIOXIDE 2B 0 0.5 Antimony TLV 0001309-64-4[ANTIMONY TRIOXIDE 2B 0 0.01 Antimony TLV 0001309-28-2[ARSENIC COMPOUNDS 0.01 Antimony TLV 0001309-28-2[BARIUM COMPOUNDS 0.07 0.5 And soluble empds 0000092-32-4 BIPHENYL 0.6 205 43.6.52 -1.37 Mouse 0000092-32-4 BIPHENYL 0.6 205 54.36 0000012-38-9 BROMOMETHANE 0.6 205 54.36	0007773-06-0	AMMONIUMSULFAMATE			10			1			-1.24			Rat TerraTox	
0007440-38-0[ANTIMONY 0 0.5 Antimony TLV 0001309-64-4[ANTIMONYTRIOXIDE 28 0 0.5 Antimony TLV 0001303-64-4[ANTIMONYTRIOXIDE 28 0 0.01 Antimony TLV 0001303-28-2[ARSENIC 0 0.01 Antimony TLV Antimony TLV 000740-39-3[BARIUM 0.07 0.5 Image: Compounds and inorg. cmpds, not Arsine Arsinie 0000361-37-2[BARIUM 0.07 0.5 Image: Compounds and inorg. cmpds, not Arsine Barium TLV 0000361-37-2[BARIUM 0.07 0.5 Image: Compounds and inorg. cmpds, not Arsine Barium TLV 0000032-23-1[BISETHYLHEXYL 0.6 205 546.36 Image: Compounds and inorg. cmpds, not Arsine 0000013-23-56 Image: Compounds and inorg. cmpds, not Arsine Image: Compounds and inorg. cmpds, not Arsine 0000012-23-1[BISETHYLH	0000062-53-3	ANILINE	1	1	7.6	70	2.5704		4.69		-0.67	0	96hr FHM Static TerraTox	Rat TerraTox	1 V incl. homologues
0001309-64-4 Antimony COMPOUNDS 2B 0 0.5 Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 2B 0 0.5 Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 2B 0 0.65 Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 0 0.01 Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 0 0.01 Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 0 0.01 Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 0.01 Antimony TLV Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 0.01 Antimony TLV Antimony TLV 0001309-64-4 ANTIMONYTRICXIDE 0.01 Antimony TLV Antimony TLV 0001309-64-4 Antimony TLV 0.05 - - Antimony TLV 00001309-64-4 Antimony TLV 0.61 0.05 - - - 00001239-18-5 BROMOCHLORODIFLUCROMETHANE 0.66 none - - - - <td< td=""><td>0007440-36-0</td><td>ANTIMONY</td><td></td><td>0</td><td>0.5</td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td>and compounds</td></td<>	0007440-36-0	ANTIMONY		0	0.5										and compounds
0001309-64-4 ANTIMONYTRIOXIDE 28 0 0.5 Antimony TLV 0007440-38-2 ARSENIC 0 0.01 and inorg. cmpds, not Arsine Antimony TLV 0001303-22-2 ARSENIC COMPOUNDS 0.01 Antimony TLV Arsenic TLV 0001303-22-2 BARIUM 0.07 0.5 Arsenic TLV Arsenic TLV 000094-36-30 BARIUM COMPOUNDS 0.5 Interview Arsenic TLV 000092-52-4 BIPHENYL 0.6 205 545.36 Interview 000012-33-1 BISETHYLHEXYL 0.6 205 546.36 Interview Mouse, 0000074-83-9 BROMOCHLORODIFLUOROMETHANE 0 3.9 1.981 -0.35 Rat, TerraTox Mouse, 000012-38-64 BUTYL ACETATE-T 200 16 7.7332 -0.45 96hr, FHM, Static, TerraTox Mouse, 000001-3-0-9 BUTYL ACETATE-T 200 16 7.7332 -0.45 96hr, FHM, Static, TerraTox Mouse, 0000071-36-82 BUTYLACETATE 713 22<	0001309-64-4	ANTIMONY COMPOUNDS	2B	0	0.5			Í							Antimony TLV
0007440-38-2 ARSENIC 0 0.01 and inorg. cmpds, not Arsine 0001303-28-2 ARSENIC COMPOUNDS 0.01 Arsenic TLV 000740-39-3 BARIUM 0.07 0.5 and soluble cmpds 0001361-37-2 BARIUM COMPOUNDS 0.5 and soluble cmpds 000094-38-0 BERZOYLPEROXIDE 5 -1.37 Mouse, 0000094-38-0 BISETHYLHEXYL 0.6 205 546.36 -1.09 1.9 96hr, FHM, Static, TerraTox Mouse, 0000728-95-8 BROMINE 0.66 none -0.35 Rat, TerraTox Mouse, 000010-19-0 BUTYL ACETATE-I 0.13 21 6.8197 -0.45 96hr, FHM, Static, TerraTox skin 000011-19-0 BUTYL ACETATE-I 0.1 7.7322 -0.45 96hr, FHM, Static, TerraTox Mouse, 0000071-3-8-3 BRUMOMETHANE -0.35 Rat, TerraTox Mouse, 000011-19-0 BUTYL ACETATE-I 0.1 7.7322 -0.45 96hr, FHM, Static, TerraTox 0000012-3-28-3 <	0001309-64-4	ANTIMONYTRIOXIDE	2B	0	0.5	i				· · · · · · · · · · · · · · · · · · ·	······································				Antimony TLV
0001303-28-2 ARSENIC COMPOUNDS 0.01 Arsenic TLV 0007440-39-3 BARIUM 0.07 0.5 and soluble cmpds 001303-28-2 BARIUM COMPOUNDS 0.5 and soluble cmpds 000094-38-0 BENZOYLPEROXIDE 5 -1.37 Mouse 000092-52-4 BIPHENYL 0.6 205 546.36 -1.09 1.9 96hr, FHM, Static, TerraTox Mouse 000013-23-1 BISETHYLHEXYL 0.6 205 546.36 -	0007440-38-2	ARSENIC		0	0.01									and ino	ra, cmpds, not Arsine
0007440-39-3 BARIUM 0.07 0.5 and soluble cmpds 0010361-37-2 BARIUM COMPOUNDS 0.5 -1.37 Mouse, Barium TLV 000092-52-4 BIPHENYL 1.3 436.52 -1.09 1.9 96hr, FHM, Static, TerraTox Mouse, 000013-23-1 BISETHYLHEXYL 0.6 205 546.36 - - - - 0000726-95-6 BROMINE 0.66 none -	0001303-28-2	ARSENIC COMPOUNDS			0.01									ĺ	Arsenic TLV
0010361-37-2 BARIUM COMPOUNDS 0.5 Barium TLV 0000094-36-0 BENZOYLPEROXIDE 5 -1.37 Mouse, 0000092-52-4 BIPHENYL 1.3 436.52 -1.09 1.9 96hr, FHM, Static, TerraTox Mouse, 000012-23-1 BISETHYLHEXYL 0.6 205 546.36 - <	0007440-39-3	BARIUM		0.07	0.5								· · · ·		and soluble cmods
0000094-36-0 BENZOYLPEROXIDE 5 -1.37 Mouse, 0000092-52-4 BIPHENYL 1.3 436.52 -1.09 1.9 96hr, FHM, Static, TerraTox Mouse, 000103-23-1 BISETHYLHEXYL 0.6 205 546.36 -	0010361-37-2	BARIUM COMPOUNDS			0.5										Barium TI V
0000092-52-4 BIPHENYL 1.3 436.52 -1.09 1.9 96hr, FHM, Static, TerraTox Mouse, 0000103-23-1 BISETHYLHEXYL 0.6 205 546.36 -	0000094-36-0	BENZOYLPEROXIDE		1	5				****	······	-1.37		**********	Mouse	
0000103-23-1 BISETHYLHEXYL 0.6 205 546.36 1 <th1< th=""> 1 1 <</th1<>	0000092-52-4	BIPHENYL		1	1.3		436.52				-1.09	1.9	96hr, FHM, Static, TerraTox	Mouse	· · · · · · · · · · · · · · · · · · ·
0007726-95-6 BROMINE 0.66 none	0000103-23-1	BISETHYLHEXYL	1	0.6		205	546.36							modeo,	
0000353-59-3 BROMOCHLORODIFLUOROMETHANE	0007726-95-6	BROMINE	1	1	0.66	none		í	····						
0000074-83-9 BROMOMETHANE 0 3.9 1.981 -0.35 Rat, TerraTox skin 0000110-19-0 BUTYL ACETATE-I 713 21 6.8197 - <td>0000353-59-3</td> <td>BROMOCHLORODIFLUOROMETHANE</td> <td>1</td> <td></td> <td></td> <td>i</td> <td>******</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td>	0000353-59-3	BROMOCHLORODIFLUOROMETHANE	1			i	******								
0000110-19-0 BUTYL ACETATE-I 713 21 6.8197 713 21 6.8197 713	0000074-83-9	BROMOMETHANE		0	3.9		1.981				-0.35			Rat TerraTox	skin
0000540-88-5 BUTYL ACETATE-T 200 16 7.7332 -0.45 96hr, FHM, Static, TerraTox 0000123-86-4 BUTYLACETATE 713 22 7.4159 -1.76 0.81 96hr, FHM, Static, TerraTox Mouse, 0000141-32-2 BUTYLACRYLATE 52 39 22.986 -0.85 Rat, TerraTox Mouse, 0000071-36-3 BUTYLALCOHOLA 0.1 no TWA 35 1.0347 -1.03 -1.37 96hr, FHM, Static, TerraTox Rat, TerraTox 0000078-92-2 BUTYLALCOHOLB 303 26 -0.87 -1.94 96hr, FHM, Static, TerraTox Rat, TerraTox 0000075-65-0 BUTYLALCOHOLC 303 4 0.3408 -1.67 -1.94 96hr, FHM, Static, TerraTox Rat, TerraTox 0000085-68-7 BUTYLBENZYLPHTHALA 0.2 213 776.25 -0.87 2.13 96hr, FHM, Static, TerraTox Rat, TerraTox 0000084-74-2 BUTYLBHTHALATE 0.1 171 3013.9 -1.28 2.4 96hr, FHM, Static, TerraTox Mouse,	0000110-19-0	BUTYL ACETATE-I			713	21	6.8197					······································			
0000123-86-4 BUTYLACETATE 713 22 7.4159 -1.78 0.81 96hr, FHM, Static, TerraTox Mouse, 0000141-32-2 BUTYLACRYLATE 52 39 22.986 -0.85 Rat, TerraTox Mouse, 0000071-36-3 BUTYLALCOHOLA 0.1 no TWA 35 1.0347 -1.03 -1.37 96hr, FHM, Static, TerraTox Rat, TerraTox 0000078-92-2 BUTYLALCOHOLB 303 26 -	0000540-88-5	BUTYL ACETATE-T	11		200	16	7.7332	· · · ·			······	-0.45	96br EHM Static TerraTox		
0000141-32-2 BUTYLACRYLATE 52 33 22.986 -0.85 Rat, TerraTox 0000071-36-3 BUTYLALCOHOLA 0.1 no TWA 35 1.0347 -1.03 -1.37 96hr, FHM, Static, TerraTox Rat, TerraTox 0000078-92-2 BUTYLALCOHOLB 303 26 -	0000123-86-4	BUTYLACETATE			713	22	7.4159				-1.78	0.81	96br FHM Static TerraTox	Mouse	
0000071-36-3 BUTYLALCOHOLA 0.1 no TWA 35 1.0347 -1.03 -1.37 96hr, FHM, Static, TerraTox Rat, TerraTox 0000078-92-2 BUTYLALCOHOLB 303 26 -1.67 -1.94 96hr, FHM, Static, TerraTox Rat, TerraTox 0000075-65-0 BUTYLALCOHOLC 303 4 0.3408 -1.67 -1.94 96hr, FHM, Static, TerraTox Rat, TerraTox 0000085-68-7 BUTYLBENZYLPHTHALA 0.2 213 776.25 -0.87 2.13 96hr, FHM, Static, TerraTox Rat, TerraTox 0000084-74-2 BUTYLPHTHALATE 0.1 171 3013.9 -1.28 2.4 96hr, FHM, Static, TerraTox Mouse, 0000123-72-8 BUTYRALDEHYDE -11 1.0347 -1.54 0.65 96hr, FHM, Static, TerraTox Rat, TerraTox	0000141-32-2	BUTYLACRYLATE	1		52	39	22.986				-0.85			Rat TerraTox	
0000078-92-2 BUTYLALCOHOLB 303 26 100	0000071-36-3	BUTYLALCOHOLA	1	0.1	no TWA	35	1.0347				-1.03	-1.37	96hr EHM Static TerraTox	Rat TerraTox	
0000075-65-0 BUTYLALCOHOLC 303 4 0.3408 -1.67 -1.94 96hr, FHM, Static, TerraTox Rat, TerraTox 0000085-68-7 BUTYLBENZYLPHTHALA 0.2 213 776.25 -0.87 2.13 96hr, FHM, Static, TerraTox Rat, TerraTox 0000084-74-2 BUTYLPHTHALATE 0.1 171 3013.9 -1.28 2.4 96hr, FHM, Static, TerraTox Mouse, 0000123-72-8 BUTYRALDEHYDE -11 1.0347 -1.54 0.65 96hr, FHM, Static, TerraTox Rat, TerraTox	0000078-92-2 6	BUTYLALCOHOLB	1		303	26		1							
0000085-68-7 BUTYLBENZYLPHTHALA 0.2 213 776.25 -0.87 2.13 96hr, FHM, Static, TerraTox Rat, TerraTox 0000084-74-2 BUTYLPHTHALATE 0.1 171 3013.9 -1.28 2.4 96hr, FHM, Static, TerraTox Rat, TerraTox 0000123-72-8 BUTYRALDEHYDE -11 1.0347 -1.54 0.65 96hr, FHM, Static, TerraTox Rat, TerraTox	0000075-65-0	BUTYLALCOHOLC	1		303	4	0.3408				-1 67	-1 94	96hr FHM Static TerraTov	Rat TerraTov	
0000084-74-2 BUTYLPHTHALATE 0.1 171 3013.9 -1.28 2.4 96hr, FHM, Static, TerraTox Mouse, 0000123-72-8 BUTYRALDEHYDE -11 1.0347 -1.54 0.65 96hr, FHM, Static, TerraTox Rat TerraTox	0000085-68-7	BUTYLBENZYLPHTHALA	<u>†</u> †	0.2		213	776.25				-0.87	2 13	96hr EHM Static TerraTox	Rat TerraTov	
0000123-72-8 BUTYRALDEHYDE -11 1.0347 -1.54 0.65 96hr FHM Static TerraTox Rat TerraTox	0000084-74-2	BUTYLPHTHALATE	1	0.11		171	3013.9				-1 28	24	96br EHM Static TerraTox	Mouse	
	0000123-72-8	BUTYRALDEHYDE				-11	1.0347				-1.54	0.65	96hr EHM Static TerraTox	Rat TerraTor	

CACNUmber	Chamient Manua	1		ILV(IWA in		-	14	рКа	pK	LD50,	CLC50			
CAS Number		IARC	RFD	mg/m^3)	FP (C)	BCF	рКа	(ca)	notes	pT	, pT	Aquatic LC50 notes	LD50 notes	TLV Notes
000/440-43-9	CADMIUM	1	0	0.01							ļ	·		and compounds
0010108-64-2	CADMIUM COMPOUNDS	1	ļ	0.01					ļ	0.49	2.08	96hr, FHM, Static, TerraTox	Mouse,	Cadmium TLV
0000075-20-7			ļ					ļ	L					
0007778-54-3	CALCIUM HYPOCHLORITE	1							ļ		[· · · · · · · · · · · · · · · · · · ·		
0000105-60-2	CAPROLACTAM DUST AND VAPOR	Į	0.5	1		0.6525							Dust	LV=1, Vapor TLV=23
0000056-23-5	CARBONTETRACHLORIDE	<u>2B</u>	0	31	none	17.378			ļ	-1.06	0.55	96hr, FHM, Static, TerraTox	Rat, TerraTox	skin
0007782-50-5	CHLORINE		0.1	1.5							ļ			
0000108-90-7	CHLOROBENZENE	ļ	0.02	46	24	446.68			ļ	-1.31	0.7	96hr, FHM, Static, TerraTox	Mouse,	
0000075-45-6	CHLORODIFLUOROMETHANE	ļ		3540		1.5732		<u> </u>						
0000067-66-3	CHLOROFORM	<u>2B</u>	0.01	49	none	6.0256				0.52	0.23	96hr, FHM, Static, TerraTox	Mouse,	
0000074-87-3	CHLOROMETHANE	<u></u>		103	-40	1.1018				-1.55	-1.04	96hr, bgill, static, TerraTox	Rat, TerraTox	skin
0000095-57-8	CHLOROPHENOL	<u> </u>	0.01		64	213.8	8.56			-0.43	1.02	96hr, FHM, Static, TerraTox	Mouse,	
0007790-94-5	CHLOROSULFONIC ACID				none				ļ					
0001897-45-6	CHLOROTHALONIL	<u></u>	0.02			41.319				-1.14	2.83	6hr, bgill, static, TerraTox	Mouse,	
0007738-94-5	CHROMIC ACID	<u></u>	<u> </u>	0.01	ļ		0.74		CRC		ļ			Insol Cr (VI) cmpds
0010101-53-8	CHROMIC SULFATE			0.5							ļ		TLV for Cr	metal & Cr(III) cmpds
0007440-47-3	CHROMIUM	1	0.01	0.5								•	Cr	metal & Cr(III) cmpds
0001333-82-0	CHROMIUM AND COMPOUNDS	1	ĮĮ	0.01			0.74		CRC					Insol Cr (VI) cmpds
0028407-37-6	CI DIRECT BLUE 218	ļ	<u> </u>											
0002832-40-8	CIDISPERSEYELLOW	<u></u>	ļ											·····
0000081-88-9	CIFOODRED15									-0.27			Mouse,	
0000097-56-3	CISOLVENTYELLOWA					1278.2					[
0007440-48-4	COBALT	2B		0.02						:			-	and inorganic cmpds
0007646-79-9	COBALT COMPOUNDS	2B		. 0.02										TLV for Cobalt
0007440-50-8	COPPER	1		0.05								& inorg cmpds;	LV for fume ar	nd respirable particles
0007758-98-7	COPPER COMPOUNDS			0.05						-0.27			Rat, TerraTox	TLV for Copper
0008001-58-9	CREOSOTE	2A												
0000108-39-4	CRESOLA		0.05	22	86	10.369	10.1			-0.35	0.29	96hr, FHM, Static, TerraTox	Rat, TerraTox	Cresol mixed isomer
0000095-48-7	CRESOLB		0.05	22	81	10.715	10.3			-0.05	0.77	96hr, FHM, Static, TerraTox	Rat, TerraTox	Cresol mixed isomer
0001319-77-3	CRESOLMIXEDISOMER	<u> </u>	0	22										skin; all isomers
0000098-82-8	CUMENE	ļ	0.04	246	46	348.29				-1.07	1.28	96hr, FHM, Static, TerraTox	Rat, TerraTox	, skin
0003251-23-8	CUPRIC NITRATE			0.05										TLV for Copper
0007758-98-7	CUPRIC SULFATE		•	0.05						-0.27			Rat, TerraTox	TLV for Copper
0000143-33-9	CYANIDE COMPOUNDS		0.04	no TWA		0.0047				0.88	2.46	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000110-82-7	CYCLOHEXANE			1030	-18	220.19				-0.98	0.38	96hr, bgill, static, TerraTox	Mouse,	
0000108-94-1	CYCLOHEXANONE .	1	5	100	47	0.8935		L		-1.22	-0.81	96hr, FHM, Static, TerraTox	Mouse,	skin
0000096-12-8	DBCP	2B				80.736				0.14			Rat, TerraTox	
0001163-19-5	DECABROMODIPHENYLOX		0.01			0.0315							****************	
0000095-50-1	DICHLOROBENZENEA		0.09	150	66	89.125				-0.53	1.4	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000106-46-7	DICHLOROBENZENEC	2B		60	66	60.256				-0.53	1.62	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0025321-22-6	DICHLOROBENZENEMIX			60	64								TL	V for p-dichlorobenze
0000075-27-4	DICHLOROBROMOMETHANE	2B	0.02		none	10.813				-0.44			Mouse,	
0000075-71-8	DICHLORODIFLUOROMETHANE		0.2	4950		15.119								
0000107-06-2	DICHLOROETHANE	2B		40	16	1.9953				-0.69	-0.14	96hr, FHM, Static, TerraTox	Mouse,	
0000156-60-5	DICHLOROETHYLENE	1	0.02		6	13.057			1	-1.1			Rat, TerraTox	
0000075-09-2	DICHLOROMETHANE	2B	0.06	174	none	2.2464				-1.39	-0.56	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000076-14-2	DICHLOROTETRAFLUOROETHANE	1	Ì	6990		60.23				-				
0000111-42-2	DIETHANOLAMINE			2	138	0.0082		8.88		-0.83	-1.65	96hr, FHM, Static, TerraTox	Rat, TerraTox	skin
0000109-89-7	DIETHYLAMINE			15	-28	0.5518		11.09		-0.83	-1.07	96hr, FHM, Static, TerraTox	Mouse,	skin
0000117-81-7	DIETHYLHEXYLPHT	2B	0.02	5	207	851.14				-1.89	1.8	96hr, goldf, static, TerraTox	Rat, TerraTox	

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		1		TLV(TWA in				рКа	рК	LD50,	c LC50			1
CAS Number	Chemical Name	IARC	RFD	mg/m^3)	FP (C)	BCF	рKa	(ca)	notes	pT	, pT	Aquatic LC50 notes	LD50 notes	TLV Notes
0000084-66-2	DIETHYLPHTHALATE		0.8	5	162	117.49		1	1	-1.44	0.84	96hr, FHM, Static, TerraTox	Mouse	
0000064-67-5	DIETHYLSULFATE	2A	4		78	1.784		1		-0.62		· · · · · · · · · · · · · · · · · · ·	Mouse	
0000822-06-0	DIISOCYANATES	1		0.034	140	133.38		1	1	-0.32			Mouse	
0000124-40-3	DIMETHYLAMINE	}		9.2	15	0.0738		10.78	3	-0.85	-0.67	96hr, guppy, static, TerraTox	Mouse	
0000121-69-7	DIMETHYLANILINE		0	25	63	20.701		[· 1	-1.07	0.19	96hr, FHM, Static, TerraTox	Rat, TerraTo	
0000068-12-2	DIMETHYLFORMAMIDE	2B	0.1	30	58	0.0197		1	1	-1.58	-2.16	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000131-11-3	DIMETHYLPHTALATE		10	5	156	57.544				-1.54	0.21	96hr, FHM, Static, TerraTox	Rat. TerraTox	
0000117-84-0	DIOCTYLPHTHALATE	}	0.02		219	7455.7			T ·	-2.08	1		Rat, TerraTox	
0000123-91-1	DIOXANE	2B		90	12	0.0679				-1.68	-2.05	96hr, FHM, Static, TerraTox	Rat, TerraTox	skin
0027176-87-0	DODECYLBENZENESULFONIC ACID									-0.3	1		Rat, TerraTox	
0000106-89-8	EPICHLOROHYDRIN	2A	. 0	1.9	33	0.4202		1		0.01	0.86	96hr, FHM, Static, TerraTox	Rat, TerraTox	skin
0000110-80-5	ETHOXYETHANOL	1	0.4	18	44	0.0528				-1.43	-2.05	96hr, bgill, static, TerraTox	Mouse	skin
0000141-78-6	ETHYLACETATE		0.9	1440	-3	0.7556			1	-1.67	-0.42	96hr, FHM, Static, TerraTox	Mouse	
0000140-88-5	ETHYLACRYLATE	2B		20	16	2.6013				-1.25	1.6	96hr, FHM, Static, TerraTox	Mouse.	
0000100-41-4	ETHYLBENZENE		0.1	434	22	120.14]	1	-1.52	0.4	96hr, FHM, Static, TerraTox	Rat. TerraTox	
0000074-85-1	ETHYLENE			listed, no TLV		1.747		Ì	1					
0000142-59-6	ETHYLENE BIS DITHIOCARBAMATE					0.0017		1	1	-0.19	1.65	96hr, guppy, static, TerraTox	Rat. TerraTox	
0000107-15-3	ETHYLENEDIAMINE	1	0.02	25	34	0.0112		9.93	1	-0.91	-0.28	96br, FHM, Static, TerraTox	Rat TerraTox	skin
0000060-00-4	ETHYLENEDIAMINE-TETRAACETIC]			33	5E-05				-0.01	0.69	96hr, FHM, Static, TerraTox	Mouse	
0000107-21-1	ETHYLENEGLYCOL		2	no TWA		0.0029				-2.08	-2.93	96hr, FHM, Static, TerraTox	Mouse	aerosol
0000075-21-8	ETHYLENEOXIDE	2A		1.8	>110	0.0873				-0.88	-0.28	96hr FHM Static TerraTox	Rat TerraTox	4010001
0000096-45-7	ETHYLENETHIOUREA	2B	0			0.0411				-1.25	-2.87	96hr. guppy, static, TerraTox	Rat TerraTox	
0000060-29-7	ETHYLETHER	1	. 0.2	1210	-40	1.0347				-1.22	-1.54	96hr, FHM, Static, TerraTox	Rat TerraTox	
0007705-08-0	FERRICCHLORIDE	1		1		ĺ								soluble iron salts
0010028-22-5	FERRICSULFATE			1			Í							soluble iron salts
0010045-89-3	FERROUSAMMONIUM SULFATE			1										soluble iron salts
0007758-94-3	FERROUSCHLORIDE			1								· · · ·		soluble iron salts
0007720-78-7	FERROUSSULFATE			1									······	soluble iron salts
0007782-63-0	FERROUSSULFATE	1	1	1									·	soluble iron salts
0000133-07-3	FOLPET	1	0.1			327.2			· · · ·	-0.72	3	96hr, FHM, Static, TerraTox	Mouse	
0000050-00-0	FORMALDEHYDE	2A	0.2	no TWA	56	0.3408	13.29			-0.15	0.1	96hr FHM Static TerraTox	Mouse	
0000064-18-6	FORMIC ACID		2	9.4	69	0.0528	3.75			-1.18			Mouse	
0000076-13-1	FREON113		30			122.68						······		
0000110-17-8	FUMARIC ACID				·		3.03		CRC	-1.9		·	Rat TerraTox	
0000109-99-9	FURAN, TETRAHYDRO-			590	-17	0.4291		-2 1		-1.5	-1 48	96hr FHM Static TerraTox	Mouse	
0000098-01-1	FURFURAL	j	0	7.9	73	538.75				0.17	0.67	96br EHM Static TerraTox	Rat TerraTox	skin
0000110-80-5	GLYCOL ETHERS	1	0.4	18	44	0.0528				-1 43	-2.05	96hr baill static TerraTox	Mouse	skin
0000110-54-3	HEXANE (N-HEXANE)		0.06	176	-23	705				-2 52	1 54	96br EHM Static TerraTox	Pat TerraToy	31(11)
0000302-01-2	HYDRAZINE	2B		0.013	52	0.0021		*****	······	-0.26	1 51	96hr baill static TerraTox		skin
0007647-01-0	HYDROCHLORICACID			no TWA		0.5075				0.20		Join, Jgm, State, Tenarox	wouse,	
0007664-39-3	HYDROGENFLUORIDE			no TWA	поле	0.0010								
0000078-83-1	ISOBUTYL ALCOHOL	<u> </u>	0.3	152	37	0.8046				-1 52	_1 3	96br EHM Static TerraTox	Pat TerraTev	
0000078-59-1	ISOPHORONE		0.2	no TWA	84	7 0795				-1 23	-0.22	96br EHM Static TerraTox	Pat TerraTex	
0004098-71-9	ISOPHORONE DIISOCYANATE			0.045	84						U.22		nac, renarox	
0000067-63-01	ISOPROPYLALCOHOL	11		983	22	0.1818				-1 78	-2 21	96hr EHM Static TerraToy	Mourea	
0000080-05-71	ISOPROPYLIDENED	<u>i </u>	0.05							1 02			Mouse,	
0007439-92-11	LEAD	2B		0.05						-1.02		******	wouse,	inorganic compounda
0007758-97-6	LEAD CHROMATE	1		0.012									and	as Cr. 05 po Lood
0007758-95-4		'		0.05										TIV for Lood
001009-74-8		<u>├</u>		0.05	ł									
0010000-14-01		1		0.001			1					· · ·		ILV IOF Lead

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				TLV(TWA in				рКа	рК	LD50,	c LC50			
CAS Number	Chemical Name	IARC	RFD	mg/m^3)	FP (C)	BCF	рКа	(ca)	notes	рТ	., pT	Aquatic LC50 notes	LD50 notes	TLV Notes
0014307-35-8	LITHIUM CHROMATE	1		0.05							Ì		wa	ater sol Cr (VI) cmpds
0000110-16-7	MALEICACID					•	1.83		CRC					
0000108-31-6	MALEICANHYDRIDE		0.1	. 1	103	4.8773				-0.61			Rat. TerraTox	······
0007439-96-5	MANGANESE		0.14	0.2								· · · · · · · · · · · · · · · · · · ·	and	inorganic compounds
0001344-43-0	MANGANESE COMPOUNDS	}		0.2								· · · · · · · · · · · · · · · · · · ·		TLV for Manganese
0000126-98-7	METHACRYLONITRILE	}	0	2.7	12	0.6805				0.6			Mouse.	skin
0000067-56-1	METHANOL	1	0.5	262	11	0.0326	15.5	······		-2.25	-2.95	96hr, FHM, Static, TerraTox	Rat. TerraTox	skin
0000109-86-4	METHOXYETHANOL	·		16	46	0.0326				-1.49		****	Rat. TerraTox	skin
0000096-33-3	METHYLACRYLATE	1	0.03	7	6	0.348		·		-0.51			Rat. TerraTox	skin
0000079-22-1	METHYLCHLOROFORMATE	1			17								,	
0005124-30-1	METHYLENE BIS(4-			0.054							<u></u>			
0000101-14-41	METHYLENEBISCHLORO	2A	0	0.11		313.85				-0.38			Mouse	skin
0000101-68-81	METHYLENEBISPHENYL	1		0.051						-0.94			Mouse	
0000078-93-31	METHYLETHYLKETONE	1	C.6	590	-3	0.3005				-1 67	-1.65	96hr FHM Static TerraTox	Rat TerraToy	
0000108-10-1	METHYLISOBUTYLKETO	1	0.08	205	13	2.5473		·····		-1.32	-0.71	96hr FHM Static TerraTox	Rat TerraTox	
0000080-62-6	METHYLMETHACRYLATE	1	0.08	410	10	2 9498				-1.56	-0.41	96br EHM Static TerraTox	Mouse	
0001634-04-4	METHYLTBUTYLET	+		144	-10	1 5136				-1.66	-0.91	96br EHM Static TerraTox	Pat TerraTov	
0000090-94-8	MICHLERSKETONE	<u> </u>				538 75				-1.00	-0.00	sonn, i rinn, Static, renarox	Nat, Tenatox	
0000075-04-7	MONOETHYLAMINE			92	-16	0 1247		10.81		0.05			Bat TorraTay	akin
0000924-42-51					-10	0.1241		10.01		-0.35			Rat, Tenarox	SKII
0000021 42 01			0.04	52	70	426 58				0.62	1 22	Offer FUM Statia TorraTor	Mayaa	· · · · · · · · · · · · · · · · · · ·
0007440-02-01	NICKEI	28	0.07	0.5		420.00				-0.02	1.52	Solit, Frim, Static, Tenatox	Mouse,	
0000373-02-4			0.02	0.05								· · · · · · · · · · · · · · · · · · ·		coluble Ni emeda
0006018-89-91				0.00										Soluble NI Cimpus
0037211-05-5				0.05			· · · ·						·	soluble NI cmpds
0007761 99 91	NITRATE COMPOLINIDS			0.03							1.00			soluble NI cmpas
0007607 27 21				5.2						0.53	4.28	96hr, FHM, Static, TerraTox	Mouse,	soluble Ag cmpds
0007097-37-21				0.2			>0							
0010102-43-91														
0000098-95-3		ļ	0	5	88	15.136				-0.6	0.01	96hr, FHM, Static, TerraTox	Rat, TerraTox	skin
0010102-44-0	NITROGEN DIOXIDE			5.6								· · · · · · · · · · · · · · · · · · ·		
0000088-75-5	NITROPHENOLA					6.9641	7.23			-0.38	0	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000099-55-8	NITROTOLUIDINE					1.784				-0.58	0.35	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0020325-40-0	D-DIANISIDINE DIHYDROCHLORIDE												Į	
0030525-89-4 F	PARAFORMALDEHYDE				71									
0000594-42-3 F	PERCHLOROMETHYLMERCAPTAN			0.76	none		·							
0000108-95-2 F	PHENOL		0.6	19	79	17.378	9.99			-0.46	0.51	96hr, FHM, Static, TerraTox	Mouse,	skin
0000106-50-3 F	PHENYLENEDIAMINE		0.19	0.1		0.0411				0.13			Rat, TerraTox	
0000075-44-5 F	PHOSGENE			0.4		0.037			-					
0007664-38-2 F	PHOSPHORICACID			1			2.12		CRC			-		· · · · · · · · · · · · · · · · · · ·
0010025-87-3 F	PHOSPHORUS OXYCHLORIDE			0.63			1			-0.39	, in the second s		Rat, TerraTox	
0000117-81-7 F	PHTHALATE ESTERS	2B	0.02	5	207	851.14				-1.89	1.8	96hr, goldf, static, TerraTox	Rat, TerraTox	
0000085-44-9 F	PHTHALICANHYDRIDE		2	6.1	151	0.0446		·····		-1.43			Rat, TerraTox	
0000109-06-8 F	PICOLINE				26	1.6753	i			-0.86	-0.98	96hr, FHM, Static. TerraTox	Mouse	
0001336-36-3 F	POLYCHLORINATEDBIPH	2A	0	0.5					· · · · · · · · · · · · · · · · · · ·			skin; TL	V for chlorodipl	nenyl (0011097-69-1)
0009016-87-9 F	POLYMERIC DIPHENYLMETHANE	11					1					1	1	
0001310-58-3 F	POTASSIUMHYDROXIDE	j{		no TWA				>0			· · ·			
0007722-64-7 F	POTASSIUMPERMANGANATE	11		0.2			i					TLV for M	anganese and i	norganic compounds
0000079-09-4	PROPIONICACID	İ		30	51	0.3268	4 87			-1.55			Rat TerraTov	
0000107-12-0	PROPIONITRILE	i			6	0.2289				0.19	-1 44	96hr EHM Static TerraTox	Morise	
0000075-55-8		28		47	15					0.19	1.74		Rat TerraTov	ekin
0000070-00-00	· · · · · · · · · · · · · · · · · · ·	<u>)</u>	1		-10;		1	1		0.40			nut, renaroz	57011

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CAS Number (Chemical Name (NRC) RPC (I) RCP PT (I) Accurate LCS0 notes LDS0 notes TV Notes 000007356-05 PRODING 2 0.533 - 0.52 0.33 0.591, bijl, solas, torral Tox Franta ocurate torral tox Internation 00000715-05 PRODING 0.00 0.22 0.0		[·			TLV(TWA in				pKa	рК	LD50,	c LC50			
00007-80-9PROPUEDENDEE 28 40 37 0.533 537 0.533 537 0.533 537 0.533 537 0.533 537 0.533 537 0.533 0.537 <th0.537< th=""> 0.537 0.53</th0.537<>	CAS Number	Chemical Name	IARC	RFD	mg/m^3)	FP (C)	BCF	рКа	(ca)	notes	рТ	, pT	Aquatic LC50 notes	LD50 notes	TLV Notes
0000108-85-IP/FRIDNE 0.0 16 200 6.33 5.23 1.13 0.00 69hr, FMA, Stato, Terrai Too Mouse 0000108-65, ISEENUAL 0.02 0.27 0.4 0.28 0.09 69hr, FMA, Stato, Terrai Too Mouse 0000108-65, ISEENUAL AND/DO 0.01 0.2 0.27 0.4 0.28 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.02 0.00 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.00 0.01	0000075-56-9	PROPYLENEOXIDE	2B	8	48	-37	0.1936]	1	-0.82	-0.39	96hr, bgill, static, TerraTox	Rat. TerraTox	
0000104-93 (RESORCINOL - 46 0.273 64 - 2.83 Str. FHJL, Static, TerraTor Mouse. 00000771-89-03 (ELNUMA COMPOLINDS 0.001 0.22 - <td>0000110-86-1</td> <td>PYRIDINE</td> <td>1</td> <td>0</td> <td>16</td> <td>20</td> <td>0.639</td> <td></td> <td>5.23</td> <td></td> <td>-1.05</td> <td>-0.1</td> <td>96hr, FHM, Static, TerraTox</td> <td>Rat. TerraTox</td> <td></td>	0000110-86-1	PYRIDINE	1	0	16	20	0.639		5.23		-1.05	-0.1	96hr, FHM, Static, TerraTox	Rat. TerraTox	
0000071-2-358LENLM 0.02 0.22 0.22 0.22 0.24 <td>0000108-46-3</td> <td>RESORCINOL</td> <td>1</td> <td></td> <td>45</td> <td></td> <td>0.875</td> <td>9.4</td> <td>1</td> <td></td> <td>-0.26</td> <td>0.04</td> <td>96hr, FHM, Static, TerraTox</td> <td>Mouse.</td> <td>· · · · · · · · · · · · · · · · · · ·</td>	0000108-46-3	RESORCINOL	1		45		0.875	9.4	1		-0.26	0.04	96hr, FHM, Static, TerraTox	Mouse.	· · · · · · · · · · · · · · · · · · ·
0007742-0-3/SELENUM ALCO COMPOLINDS 0.01 0.2 0.01 0.01 0.01 0.01 0.01 0.01 0.02	0000057-12-5	SELENIUM	[0.02	0.2				[1		1			and compounds
0007748-00-4]SELENUM AND COMPOUNDS 0	0007782-49-2	SELENIUM	1	0.01	0.2		Ì]	1	<u> </u>	<u>j</u>			and compounds
00077830-04 SELENUM DIOLOGE 0.01 0.2 0.02 Rat Terra Too Rat To	0007446-08-4	SELENIUM AND COMPOUNDS	İ		0.2				1						and compounds
00070985-04_SELENUM DICKUE Image: constraint of the second	0007783-00-8	SELENIUM CMPD?	1	0.01	0.2				<u> </u>		1				and compounds
0007738-9-ISELENUM NEXAFLUCRIDE 0.06 0.02 0.07 0.02 0.02 0.02 0.02 0.02 0.02 0.01 0.02 0.01 0.02 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.03 0.01 0.01 0.01 0.01 0.01 0.01 0.01	0007488-56-4	SELENIUM DIOXIDE	1		0.2		Ì		1	-	0.02			Rat. TerraTox	and compounds
0010033-09/SELENUM NITRIDE 0.02 0.03 0.01 0.02 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.03 4.28 0.01 0.01 0.01 0.01 0.03 4.28 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 <t< td=""><td>0007783-79-1</td><td>SELENIUM HEXAFLUORIDE</td><td>1</td><td></td><td>0.16</td><td></td><td></td><td></td><td></td><td>1</td><td>1</td><td><u> </u></td><td></td><td></td><td>as Selenium</td></t<>	0007783-79-1	SELENIUM HEXAFLUORIDE	1		0.16					1	1	<u> </u>			as Selenium
0014832-9-07/SELENIUM OXIDE 0.01 0.02 0.01 <	0012033-59-9	SELENIUM NITRIDE	1		· 0.2				1				· ·		and compounds
000742-24_SLVER 0.01 0.03 4.28 0017, F443, S10/EC NUMARTE 0.01	0014832-90-7	SELENIUM OXIDE	1	1	0.2		1		ĺ		1	1	***************************************	·····	and compounds
0007783-0-0.5]ULVER AND COMPOUNDS 0.01 0.07 0.03 4.28 98hr, FMM, Stalls, TernTox Mouse, sol ompds, as Silver 00070143-04_SUCENTITATE 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.00 <	0007440-22-4	SILVER		0.01	0.1		1		1			1		····.	
0007733-91-7,SILVER CHLORITE 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.01 0.007 0.008 0.007 0.007 0.008 0.0007 0.004 0.007 0.004 0.007 0.004 0.004 0.0007 0.004 0.004 0.0007 0.004 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.004 0.0007 0.0001 0.0007 0.00000000000000	0007783-90-6	SILVER AND COMPOUNDS]		. 0.01				1						sol cmpds, as Silver
000724-01-2]SILVER CHROMATE 0.01 0.01 0.01 0.03 2.0 e0 ompts, as Silver 0000506-64-39 SILVER YANDDE 0.01 0.01 0.03 4.28 9hr, FHM, Static, TerraTox Mouse, sol ompts, as Silver 000761-88-8 SILVER NTRATE 0.01 0.01 0.03 4.28 9hr, FHM, Static, TerraTox Mouse, sol ompts, as Silver 0007631-90-5 SODUM BIGUERTTE 0.01 0.01 0.03 4.28 9hr, FHM, Static, TerraTox Mouse, sol ompts, as Silver 0007631-90-5 SODUM BIGUERTE 1 0.01 0.04 no TWA 0.044 0.88 2.44 9hr, FHM, Static, TerraTox Rat, TerraTox Mouse, sol ompts, as Silver 00007631-90-5 SODUM DIGUENTE 0.04 no TWA 0.047 0.88 2.44 9hr, FHM, Static, TerraTox Rat, TerraTox Mouse, sol ompts, as Silver 00007631-94-5 SODUM VORDE 0.04 1 0.047 0.88 1.74 9hr, gupp, static, TerraTox Rat, TerraTox Mouse, sol ompts, as Silver 0007631-94-5 SODUM VPROSULPIDE 0.011 0.04 1 0.04 1 0.013 1 1 1	0007783-91-7	SILVER CHLORITE	1	1	0.01		[sol cmpds, as Silver
000506-64-63 LIVER VYANDE 0.1 0.01 0.03 0	0007784-01-2	SILVER CHROMATE			0.01		1		1				********		sol cmpds, as Silver
00077613-88-BILVER NITRATE 0.01 0.53 2.28 96hr, FHM, Static, TerraTox Mouse, sol compts, as Silver 00077613-88-BILVERNITRATE 0.01 0.53 4.28 96hr, FHM, Static, TerraTox Mouse, sol compts, as Silver 0007613-00-SIDOUM BICURROMATE 1 0.01 0.53 4.28 96hr, FHM, Static, TerraTox Mouse, sol compts, as Silver 0007613-00-SIDOUM DICHROMATE 1 0.01 -1.28 0.64 96hr, FHM, Static, TerraTox Rat, TerraTox 000712-04-11 SODUM DICHROMATE -1.28 0.64 96hr, FHM, Static, TerraTox Rat, TerraTox 000763-04-3 SODUM PURPHONENCIPE -0.84 1.74 96hr, FHM, Static, TerraTox Rat, TerraTox 000763-43-4 SODUM PURPHONENCIPE -0.84 1.74 96hr, FHM, Static, TerraTox -0.76 000763-43-59 SODUM HYDORSULPE -0.04<	0000506-64-9	SILVER CYANIDE]	0.1	0.01		1			1					sol cmods, as Silver
000716-86-8[SLVERNITRATE 0.01 0.53 4.28 96hr, FHM, Static, TerraTox Mouse, sol ompts, as Silve 001588-01-9 SODUM BICHROMATE 1 0.01 1.28 0.64 96hr, FHM, Static, TerraTox Mouse, sol ompts, as Silve 0007631-93-5 SODUM BICHROMATE 5 -1.28 0.64 96hr, FHM, Static, TerraTox Rat, TerraTox Mouse, sol ompts, as Silve 000143-33-5 SODUM CNNDE (Ma(CN)) 0.04 no TWA 0.0047 0.88 2.49 96hr, FHM, Static, TerraTox Rat, TerraTox 000143-33-5 SODUM DIMETHYLDITHICCARBAMATE 0.0047 0.84 1.74 96hr, guppy, static, TerraTox Rat, TerraTox 000163-14-9 SODUM HYDORDE 1 0.0047 0.84 1.74 96hr, FHM, Static, TerraTox Rat, TerraTox 000163-14-9 SODUM HYDORDKUEE 1 0.004 1.48 96hr, FHM, Static, TerraTox 0.0014 000162-14-15 1 0.0011 1 2.0 0.0014 0.0014 0.0014 0.0014 0.0014 0.0014 0.0014 0.0014 0.0014 0.00017	0007761-88-8	SILVER NITRATE	1		0.01		1		1		0.53	4.28	96hr, FHM, Static, TerraTox	Mouse	sol cmods, as Silver
0007440 23-5 SODIUM Insol Cr (VI) cmpds Insol Cr (VI) cmpds 0007831-90-5 SODIUM BISULFITE 5 -1.28 0.64 96hr, mosqf, static, TerraTox, Rat, TerraTox 0007831-90-5 SODIUM BISULFITE 0.04 no TWA 0.0047 0.88 2.46 96hr, mosqf, static, TerraTox, Rat, TerraTox 0001728-41-15 SODIUM BISULFITE 0.04 no TWA 0.0047 0.88 2.46 96hr, mosqf, static, TerraTox, Rat, TerraTox 0001728-41-15 SODIUM MISOLICARBAMATE 0.04 -0.34 1.74 96hr, FHM, Static, TerraTox Rat, TerraTox 0007611-94-9 SODIUM HYDORXIDE no TWA 0.0011 -0.44 -0.44 -0.44 0001728-11-25 SODIUM HYDORXIDE no TWA >0 -0.44 -0.44 -0.44 0000124-114 SODIUM INTPOCHLORITE 0.0011 1.48 96hr, FHM, Static, TerraTox -0.44 0000124-114 SODIUM HYDORHATE, TRIBASIC -0.0011 1.48 96hr, FHM, Static, TerraTox -0.44 0000124-124 SOLIUM HYDORHATE, TRIBASIC -0.0011 1.47 96hr, FHM, Static, TerraTox -0.44 0000124-54 SOLIUM HOSPHATE, TRIBASIC -0.48 <	0007761-88-8	SILVERNITRATE	1		0.01					[0.53	4.28	96hr, FHM, Static, TerraTox	Mouse	sol cmpds, as Silver
0010588-01-93 SODUM BICHROMATE 1 0.01 insol Cr (VI) cmpds 00007631-09-250DIM BICHROMATE 5 -1.28 0.64 96hr, FHM, Static, TerraTox insol Cr (VI) cmpds 0000123-39-950DIM BICHROMATE 0.04 no TWA 0.0047 0.68 2.44 96hr, FHM, Static, TerraTox Rat, TerraTox 0000123-94-15 SODIUM FLUORDE - <	0007440-23-5	SODIUM]			3								1100000,	
00073190-5 SODIUM PISULFITE 1 5 -128 0.64 96hr, mosqf, static, TerraTox Notion (v), Mipolity 0001143-39-5 SODIUM 0.04 no TWA 0.0047 0.88 2.46 96hr, mosqf, static, TerraTox Rat, TerraTox 0001128-04-1 SODIUM CMNDE (Na(CN)) 0.04 no TWA 0.0047 0.88 2.46 96hr, mosqf, static, TerraTox Rat, TerraTox 0007681-34-94 SODIUM CMNDRX - <td>0010588-01-9</td> <td>SODIUM BICHROMATE -</td> <td>1</td> <td></td> <td>0.01</td> <td></td> <td></td> <td></td> <td>İ</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>insol Cr (VI) cmpds</td>	0010588-01-9	SODIUM BICHROMATE -	1		0.01				İ						insol Cr (VI) cmpds
000143-33-9 SODIUM CYANIDE (Na(CNI)) 0.04 no TWA 0.0047 0.86 2.46 96h; FHM, Static, TerraTox 000128-04-1 SODIUM DIMETHYLDITHIOCARBAMATE -0.84 1.74 96h; gupp, state. TerraTox Rat, TerraTox 0007581-09-SODIUM FLUORIDE -0.84 1.74 96h; gupp, state. TerraTox Rat, TerraTox 0007681-09-SODIUM FLUORIDE -0.84 1.74 96h; gupp, state. TerraTox Rat, TerraTox 0007681-09-SODIUM HYDROSULFIDE -0.84 -0.84 -0.84 -0.84 -0.84 0007681-09-SODIUM HYDOCHLORITE -0.84 -0.84 -0.84 -0.84 -0.84 0007622-00-SODIUM HYDOCHLORITE -0.0011 1.48 96h; FHM, State, TerraTox -0.001 0007632-00-SODIUM HYDOSHATE, TRIBASIC -0.0011 1.48 96h; FHM, State, TerraTox -0.001 0007769-29-4 SODIUM PHOSPHATE, TRIBASIC -0.076 96hr, mosqf, state, TerraTox -0.001 0000769-20-4 SODIUM PHOSPHATE, TRIBASIC -0.076 96hr, FHM, State, TerraTox -0.001 0000769-24-9 SODIUM PHOSPHATE, TRIBASIC -0.06	0007631-90-5	SODIUM BISULFITE	· · ·		5						-1 28	0.64	96br mosof static TerraToxi	Rat TerraTox	
0000728-04-1 SODIUM DIMETHYLDITHIOCARBAMATE 0.8.4 1.74 96hr, guppy, static, TerraTox Rat, TerraTox 0007681-49-4 SODIUM FLUORIDE 0.8.4 1.74 96hr, guppy, static, TerraTox Rat, TerraTox 0007681-49-4 SODIUM HYDROSULFIDE 0.0 0.0 0.0 0.0 0007681-49-4 SODIUM HYDROXIDE 0.0 0.0 0.0 0.0 0007681-52-9 SODIUM HYDCHLORITE 0.0011 0.0011 0.0011 0.0011 00007254-73-4 SODIUM HYDCHLORITE 0.00011 0.0011 0.0011 0.0011 00007555-73-4 SODIUM PHOSPHATE, TRIBASIC 0.00011 1.48 96hr, FHM, Static, TerraTox 0.00011 00007563-73-4 SODIUM PHOSPHATE, TRIBASIC 0.0011 1.48 96hr, FHM, Static, TerraTox 0.0011 0000763-73-4 SODIUM PHOSPHATE, TRIBASIC 0.0011 1.48 96hr, FHM, Static, TerraTox Mouse, skin 0000104-25-5 SYTERLMONOMER 28 0.2 85 197.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin	0000143-33-9	SODIUM CYANIDE (Na(CN))		0.04	no TWA		0.0047				0.88	2.46	96hr FHM Static TerraTox	Rat TerraTox	
0026156-30-0 SODIUM Image: Solid Structure Solid Stru	0000128-04-1	SODIUM DIMETHYLDITHIOCARBAMATE									-0.84	1 74	96hr guppy static TerraTox	Rat TerraTox	
0007881-494 SODIUM FLUORIDE no TWA >0 00180721-80-S SODIUM HYDROSULFIDE no TWA >0 0001810-72-25 SODIUM HYDROCHLORITE 0 0 0000124-41-4 SODIUM HYDROCHLORITE 0 0 0000124-41-4 SODIUM HYDROCHLORITE 0 0 0000124-41-4 SODIUM HYDROCHLORITE 0 0 0000124-41-4 SODIUM HYDROSHATE, DIBASIC 0 0 000758-29-4 SODIUM HYDROSHATE, TRIBASIC 0 0 0000758-29-4 SODIUM HOSPHATE, TRIBASIC 0 0.001 14.8 0000768-29-4 SODIUM HOSPHATE, TRIBASIC 0 0 0.001 0000768-29-4 SODIUM HOSPHATE, TRIBASIC 0 0 0.001 0000764-29-5 SUFURICACID 1 1 >0 0 0000764-29-5 SUFURICACID 1 1 >0 0 000010-42-5 STRENEMONOMER 28 0.01 170 none 39.05 -1.26 1.09 96hr, FHM, Static, TerraTox	0025155-30-0	SODIUM			· · · ·								oom, gappy, static, renarios		
0016721-80-\$ SODIUM HYDROXIDE no TWA >0 0001310-73-2 SODIUM HYDROXIDE no TWA >0	0007681-49-4	SODIUM FLUORIDE			· · · · ·										
0001310-73-2 (SODIUM HYPOCHLORITE no TWA >0 >0 0007681-52-9 (SODIUM HYPOCHLORITE	0016721-80-5	SODIUM HYDROSULFIDE											······		
0007681-52-9 SODIUM HYPOCHLORITE Image: Constraint of the second	0001310-73-2	SODIUM HYDROXIDE			no TWA				>0						
0010022-70-S SODIUM HYPOCHLORITE	0007681-52-9	SODIUM HYPOCHLORITE		├ ─── ┤										· · · · ·	
0000124-41-4 SODUM METHYLATE 0.0011 1.48 96hr, FHM, Static, TerraTox 0007632-00-3 SODUM NITRITE 0.0011 12.67 CRC 12.67 000756-74-9 SODUM PHOSPHATE, DIBASIC 12.67 CRC 12.67 CRC 000756-74-9 SODUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 12.67 000010-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 000010-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 000010-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0000127-18-4 TETRACHLORCETHYLENE 28 0.01 170 none 38.905 -1.26 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 0000012-718-4 TETRACHL	0010022-70-5	SODIUM HYPOCHLORITE													
0007632-00-0 SODIUM NITRITE 0.0011 1.48 96hr, FHM, Static, TerraTox 000758-79-4 SODIUM PHOSPHATE, DIBASIC 12.67 CRC 96hr, mosqf, static, TerraTox 000758-29-4 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 000101-89-0 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 0000100-42-5 STRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox 0000100-42-5 STURENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0000104-25 STUEFURICACID 1 1 >0 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0000127-18-4 TETRACHLOROCTHYLENE 28 0.01 none 38.905 -1.26 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 00000108-85-3 TIOLPHENOL 0 513.2137 6.52 0.38 Rat, TerraTox skin	0000124-41-4	SODIUM METHYLATE			· · · · · ·										
0007568-79-4 SODIUM PHOSPHATE, DIBASIC 12.67 CRC 0007601-54-9 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 0000758-29-4 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 0000101-89-0 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 0000100-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 000010-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0000124-35-7 SULFURICACID 1 1 >0 -	0007632-00-0	SODIUM NITRITE					0.0011					1.48	96br EHM Static TerraTox		
0007601-54-9 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 0007758-29-4 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 000101-89-0 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 000101-89-0 SODIUM PHOSPHATE, TRIBASIC 0.76 96hr, mosqf, static, TerraTox 0000100-42-5 STYRENEMONOMER 2B 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0000164-95-7 SULFURICACID 1 >0 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 000012-54-5 STYRENEMONOMER 2B 0.01 10 >0 -0.2 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 0000108-98-5 THIOPHENOL 0 170 none 38.905 -1.26 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 00000108-98-5 THIOPHENOL 0 132.137 6.52 0.38 Rat, TerraTox Rat, TerraTox 0000010	0007558-79-4	SODIUM PHOSPHATE, DIBASIC						·····	12.67	CRC					
0007758-29-4 SODIUM PHOSPHATE, TRIBASIC 0000101-39-0 SODIUM PHOSPHATE, TRIBASIC 0000100-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 000010-42-5 STYRENEMONOMER 28 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0000104-95-7 SULFURICACID 1 1 >0 H2SO4 in strong inorg acid mists 00001127-18-4 TETRACHLOROETHYLENE 28 0.01 170 none 38.905 -1.26 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 0000108-98-5 THIOPHENOL 0 51 32.137 6.52 0.38 Rat, TerraTox 00000108-88-3 TOLUENE 0.2 188 45579 -1.74 0.43 96hr, FHM, Static, TerraTox skin 00000108-88-3 TOLUENE 0.2 188 4411.34 -1.52 0.02 96hr, FHM, Static, TerraTox skin 0000075	0007601-54-9	SODIUM PHOSPHATE, TRIBASIC			·····							0.76	96hr, mosof, static, TerraTox		
0010101-89-0 SODIUM PHOSPHATE, TRIBASIC	0007758-29-4	SODIUM PHOSPHATE, TRIBASIC													
0000100-42-S STYRENEMONOMER 2B 0.2 85 31 97.464 -0.48 1.41 96hr, FHM, Static, TerraTox Mouse, skin 0007664-93-9 SULFURICACID 1 1 >0 H2SO4 in strong inorg acid mists 0000142-55-7 SULFURICACID (FUMING) -0 -0 H2SO4 in strong inorg acid mists 0000127-18-4 TETRACHLOROETHYLENE 2B 0.01 170 none 38.905 -1.26 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 0000102-56-6 THIOPHENOL 0 51 32.137 6.52 0.38 Rat, TerraTox 0000108-88-3 TOLUENE 2B 0.0206 -0.22 Rat, TerraTox skin 00000108-88-3 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox 0000054-84-9 TOLUENEDIISOCYANATEA 2B -1.74 0.43 96hr, FHM, Static, TerraTox skin 0000271-55-6 TRICHLOROETHANEA 2B -1.05 Mous	0010101-89-0	SODIUM PHOSPHATE, TRIBASIC		1					· ·						
0007664-93-9 SULFURICACID 1 1 0 Hobel, 1	0000100-42-5	STYRENEMONOMER	2B	0.2	85	31	97,464				-0.48	1 41	96br EHM Static TerraTox	Mouse	skin
0008014-95-7 SULFURICACID (FUMING) Intervention of the state of t	0007664-93-9	SULFURICACID	1		1			>0			0.10			H2SO4 in s	trong inorg acid mists
0000127-18-4 TETRACHLOROETHYLENE 2B 0.01 170 none 38.905 -1.26 1.09 96hr, FHM, Static, TerraTox Rat, TerraTox 000018-98-5 THIOPHENOL 0 51 32.137 6.52 0.38 Rat, TerraTox 0000062-56-6 THIOUREA 2B 0.0206 -0.22 Rat, TerraTox 0000108-88-3 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox 0000058-88-3 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox 00000584-84-9 TOLUENEDIISOCYANATEB 2B 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox 0026471-62-5 TOLUENEDIISOCYANATEC 2B -1.05 Mouse, -1.05 Mouse, 0000071-65-6 TRICHLOROETHANEA 1910 none 8.9125 -1.92 0.45 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROET	0008014-95-7	SULFURICACID (FUMING)						>0							
0000108-98-5 THIOPHENOL 0 51 32.137 6.52 0.38 Rat, TerraTox 0000062-56-6 THIOUREA 2B 0.0206 -0.22 Rat, TerraTox 0000108-88-3 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox 0000091-08-7 TOLUENEDIISOCYANATEA 2B 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox 0000071-65-7 TOLUENEDIISOCYANATEB 2B 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox 0026471-62-5 TOLUENEDIISOCYANATEC 2B 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox 0026471-62-5 TOLUENEDIISOCYANATEC 2B -1.05 Mouse, Mouse, 0000071-55-6 TRICHLOROETHANEA 1910 none 8.9125 -1.92 0.45 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROETHYLENE 2A 269 none 19.851 -1.2	0000127-18-4	TETRACHLOROETHYLENE	2B	0.01	170	none	38 905				-1 26	1 09	96br EHM Static TerraTox	Rat TerraToy	
0000062-56-6 THIOUREA 2B 0.0206 -0.22 Rat, TerraTox 0000108-88-3 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox skin 0000091-08-7 TOLUENEDIISOCYANATEA 2B 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox skin 0000584-84-9 TOLUENEDIISOCYANATEB 2B 0.036 121 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox 0026471-62-5 TOLUENEDIISOCYANATEC 2B -1.05 Mouse, Mouse, 0000071-55-6 TRICHLOROETHANEA 1910 none 8.9125 -1.92 0.45 96hr, FHM, Static, TerraTox Rat, TerraTox 0000075-69-4 TRICHLOROETHYLENE 2A 269 none 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no TWA none 32.817 -1.77 Rat, TerraTox Skin	0000108-98-5	THIOPHENOL		0		51	32,137	6.52			0.38			Rat TerraTox	······································
O000108-88-3 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox skin 0000091-08-7 TOLUENE 0.2 188 4 45.879 -1.74 0.43 96hr, FHM, Static, TerraTox Rat, TerraTox skin 0000091-08-7 TOLUENEDIISOCYANATEA 2B 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox skin 0000584-84-9 TOLUENEDIISOCYANATEB 2B 0.036 121 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox skin 0026471-62-5 TOLUENEDIISOCYANATEC 2B -1.05 Mouse, Mouse, 0000071-55-6 TRICHLOROETHANEA 1910 none 8.9125 -1.92 0.45 96hr, FHM, Static, TerraTox Rat, TerraTox 0000075-69-4 TRICHLOROETHANE 2A 269 none 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 <td< td=""><td>0000062-56-6</td><td>THIOUREA</td><td>2B</td><td></td><td></td><td></td><td>0.0206</td><td></td><td></td><td></td><td>-0.22</td><td></td><td></td><td>Rat TerraTox</td><td></td></td<>	0000062-56-6	THIOUREA	2B				0.0206				-0.22			Rat TerraTox	
0000091-08-7 TOLUENEDIISOCYANATEA 2B 411.34	0000108-88-3	TOLUENE		0.2	188	4	45 879				-1 74	0.43	96br EHM Static TerraTox	Rat TerraTox	skin
0000584-84-9 TOLUENEDIISOCYANATEB 2B 0.036 121 411.34 -1.52 0.02 96hr, FHM, Static, TerraTox Rat, TerraTox 0026471-62-5 TOLUENEDIISOCYANATEC 2B -1.05 Mouse, 0000071-55-6 TRICHLOROETHANEA 1910 none 8.9125 -1.92 0.45 96hr, FHM, Static, TerraTox Rat, TerraTox 0000079-01-6 TRICHLOROETHANEA 1910 none 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROETHYLENE 2A 269 none 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no 19.821 -1.77 Rat, TerraTox skin 0000121-44-8 TRIETHYLAMINE 4.1 -6 3.3449 10.72 -0.66 Rat, TerraTox skin	0000091-08-7	TOLUENEDIISOCYANATEA	2B]	т 	411 34				1./*	0.40			JUIL
O026471-62-5 TOLUENEDIISOCYANATEC 2B -1.02 0.02 Son, FHW, State, Tenaro, Rat, T	0000584-84-9	TOLUENEDIISOCYANATEB	2B	<u> </u>	0.036	121	411 34				-1 52	0.02	96hr FHM Static TerraTovi	Rat TerraTor	
0000071-55-6 TRICHLOROETHANEA 1910 none 8.9125 -1.92 0.45 96hr, FHM, Static, TerraTox Rat, TerraTox 0000079-01-6 TRICHLOROETHYLENE 2A 269 none 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Rat, TerraTox 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no 19.851 -1.77 Rat, TerraTox 00000121-44-8 TRIETHYLAMINE 4.1 -6 3.3449 10.72 -0.66 Rat, TerraTox skin 0000075-63-8 TRIEL LIOROBBROMOMETHANE 6090 8.0641 -1.26 0.66 Rat, TerraTox skin	0026471-62-5	TOLUENEDUSOCYANATEC	2B								_1.02			Mouso	474-1 m. 1000 m. (m. 6. 7 m.) 1 m. 10.
0000079-01-6 TRICHLOROETHYLENE 2A 269 none 19.851 -1.26 0.47 96hr, FHM, Static, TerraTox Mouse, 0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no TWA none 32.817 -1.77 Rat, TerraTox Mouse, 0000121-44-8 TRIETHYLAMINE 4.1 -6 3.3449 10.72 -0.66 Rat, TerraTox skin 0000075-63-8 TRIETHYLAMINE 6090 8.0641 -1.77 Rat, TerraTox skin	0000071-55-6	TRICHLOROETHANEA			1910	none	8,9125				-1.00	0.45	96hr EHM Static TerraTov	Rat TerraTov	····
0000075-69-4 TRICHLOROMONOFLUOROMETHANE 0.3 no TWA none 32.817 -1.77 Rat, TerraTox 0000121-44-8 TRIETHYLAMINE 4.1 -6 3.3449 10.72 -0.66 Rat, TerraTox skin 0000075-63-8 TRIETHYLAMINE 6090 8.0641 -1.77 Rat, TerraTox skin	0000079-01-6	TRICHLOROETHYLENF	24		269	none	19 851				_1 26	0.47	96hr EHM Static TerraTev	Moveo	
0000121-44-8 TRIETHYLAMINE 4.1 -6 3.3449 10.72 -0.66 Rat, TerraTox skin	0000075-69-4	TRICHLOROMONOFLUOROMETHANE		0.3	no TWA	none	32 817		······		-1.20	<u> </u>		Rat TerraTov	
	0000121-44-8	TRIETHYLAMINE			4.1	-6	3,3449		10.72		-0.66		· · · · · · · · · · · · · · · · · · ·	Rat TerraTox	skin
	0000075-63-8	TRIFLUOROBROMOMETHANE			6090		8.0641				0.00				
Chemical Data

	A ·		·	TLV(TWA in				рКа	рК	LD50,	c LC50			· · · · · · · · · · · · · · · · · · ·
CAS Number	Chemical Name	IARC	RFD	mg/m^3)	FP (C)	BCF	рКа	(ca)	notes	.pT	, pT	Aquatic LC50 notes	LD50 notes	TLV Notes
0000075-50-3	TRIMETHYLAMINE	1		12	-6	0.2289		9.79						
0000095-63-6	TRIMETHYLBENZ				48	327.2		1			1.21	96hr, FHM, Static, TerraTox		· ·
0000108-05-4	VINYLACETATE	2B	1	35	-6	0.5754		1		-1.27	0.57	96hr, FHM, Static, TerraTox	Mouse,	
0000095-47-6	XYLENEB		2	434	32	112.83		1		-1.67	0.81	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0000106-42-3	XYLENEC		2	434	27	120.14				-1.67	1.21	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0001330-20-7	XYLENEMIXEDISOMER		2	434	29	120.14			1	-1.61	0.87	96hr, FHM, Static, TerraTox	Rat, TerraTox	
0007440-66-6	ZINC	}	0.3				· · ·							
0001314-13-2	ZINC AND COMPOUNDS		·	5						-1.99			Mouse,	TLV for Zinc dust=10
0001314-13-2	ZINC OXIDE FUME			5						-1.99			Mouse,	
0007733-02-0	ZINC SULFATE							}	·					
0014639-98-6	ZINCAMMONIUM CHLORIDE							1		[
0001332-07-6	ZINCBORATE		·]						
0007733-02-0	ZINCSULFATE							1						
		47	ຍ5	190	114	136	21	14	[137	102	102	137	94

1/1/-00-6	1,1-DICHLORO - 1 FLUOROETHANE
507-55-1	1,3-DICHLORO-1,1,2,2,3-PENTAFLUOROPHOPANE
8/2-50-4	1-METHYL-2-PYRROLIDONE
540-84-1	2,2,4-TRIMETHYLPENTANE
612-83-9	3,3'DICHLOROBENZIDINE DIHYDROCHLORIDE
55406-53-6	3-IODO-2-PROPYNYL BUTYLCARBAMATE
75-07-0	ACETALDEHYDE
108-24-7	ACETIC ANHYDRIDE
75-05-8	ACETONITRILE
98-86-2	ACETOPHENONE
79-10-7	ACRYLICACID
124-04-9	ADIPIC ACID
107-05-1	ALLYLCHLORIDE
7429-90-5	ALUMINUM
1344-28-1	ALUMINUMOXIDE
10043-01-3	ALUMINUMSULFATE
7664-41-7	AMMONIA
1341-49-7	AMMONIUMBIFLUORIDE
12125-02-9	AMMONIUMCHLORIDE
12125-01-8	AMMONIUMFLUORIDE
1336-21-6	AMMONIUMHYDBOXIDE
7773-06-0	AMMONIUMSULFAMATE
62-53-3	ANILINE
7440-36-0	ANTIMONY
01-00-0	ANTIMONY COMPOUNDS
1309-64-4	ANTIMONYTBIOXIDE
7440-39-3	BABIUM
01-00-2	BARIUM COMPOLINDS
94-36-0	BENZOYI PEBOXIDE
92-52-4	BIPHENYI
103-23-1	BISETHYI HEXVI
7726-95-6	BROMINE
353-59-3	BROMOCHLOBODIELLIOROMETHANE (HALON 1211)
74-83-9	BROMOMETHANE
110.10.0	
540-88-5	
123-86-4	
1/1-30-0	
75-65-0	
95 69 7	
00-00-7	
04-74-2	
123-72-8	
107-92-6	
75-20-7	
///8-54-3	
105-60-2	
106-47-8	CHLOROANILINE
108-90-7	CHLOROBENZENE
75-45-6	CHLORODIFLUOROMETHANE
74-87-3	CHLOROMETHANE
95-57-8	CHI OBOPHENOI

Appendix H: Category 3 Chemicals with CAS Numbers

Appendix H: Category 3 Chemicals with CAS Numbers

CAS Number	Chemical Name
7790-94-5	CHLOROSULFONIC ACID
1897-45-6	CHLOROTHALONIL
10101-53-8	CHROMIC SULFATE
28407-37-6	CI DIRECT BLUE 218
2832-40-8	CIDISPERSEYELLOW
81-88-9	CIFOODRED15
97-56-3	CISOLVENTYELLOWA
7440-48-4	COBALT
01-01-3	COBALT COMPOUNDS
7440-50-8	COPPER
01-01-5	COPPER COMPOUNDS
8001-58-9	CREOSOTE
108-39-4	CRESOLA
95-48-7	CRESOLB
1319-77-3	CRESOLMIXEDISOMER
98-82-8	CUMENE
3251-23-8	
7758-98-7	
110-82-7	CYCLOHEXANE
108-94-1	CYCLOHEXANONE
1163-19-5	
95-50-1	DICHLOBOBENZENEA
106-46-7	DICHLOBOBENZENEC
25321-22-6	DICHLOROBENZENEMIX
75-27-4	DICHLOROBROMOMETHANE
75-71-8	DICHLORODIFLUOROMETHANE
156-60-5	DICHLOROETHYLENE
75-09-2	DICHLOROMETHANE
76-14-2	DICHLOROTETRAFLUOROETHANE
111-42-2	DIETHANOLAMINE
109-89-7	DIETHYLAMINE
117-81-7	DIETHYLHEXYLPHT
84-66-2	DIETHYLPHTHALATE
822-06-0	DIISOCYANATES
124-40-3	DIMETHYLAMINE
121-69-7	DIMETHYLANILINE
131-11-3	DIMETHYLPHTALATE
117-84-0	DIOCTYLPHTHALATE
27176-87-0	DODECYLBENZENESULFONIC ACID
5952-26-1	ETHANOL, 2,2-OXYDI, DICARBAMATE
110-80-5	ETHOXYETHANOL
140-88-5	ETHYLACRYLATE
100-41-4	ETHYLBENZENE
74-85-1	ETHYLENE
142-59-6	ETHYLENE BIS DITHIOCARBAMATE
107-15-3	ETHYLENEDIAMINE
60-00-4	ETHYLENEDIAMINE-TETRAACETIC ACID (EDTA)
96-45-7	ETHYLENETHIOUREA
60-29-7	ETHYLETHER
133-07-3	FOLPET
64-18-6	FORMIC ACID

CAS Number	Chemical Name
1336-36-3	POLYCHLORINATEDBIPH
9016-87-9	POLYMERIC DIPHENYLMETHANE DIISOCYANATE
1310-58-3	POTASSIUMHYDROXIDE
7722-64-7	POTASSIUMPERMANGANATE
79-09-4	PROPIONICACID
107-12-0	PROPIONITRILE
110-86-1	PYRIDINE
108-46-3	RESORCINOL
7440-22-4	SILVER
01-03-7	SILVER AND COMPOUNDS
7761-88-8	SILVERNITRATE
7440-23-5	SODIUM
10588-01-9	SODIUM BICHROMATE
7631-90-5	SODIUM BISULFITE
128-04-1	SODIUM DIMETHYLDITHIOCARBAMATE
25155-30-0	SODIUM DODECYLBENZENESULFONATE
7681-49-4	SODIUM FLUORIDE
16721-80-5	SODIUM HYDROSULFIDE
1310-73-2	SODIUM HYDROXIDE
7681-52-9	SODIUM HYPOCHLORITE
10022-70-5	SODIUM HYPOCHLORITE
124-41-4	SODIUM METHYLATE
7632-00-0	SODIUM NITRITE
100-42-5	STYRENEMONOMER
108-98-5	THIOPHENOL
62-56-6	THIOUREA
108-88-3	TOLUENE
71-55-6	TRICHLOROETHANEA
75-69-4	TRICHLOROMONOFLUOROMETHANE
121-44-8	TRIETHYLAMINE
75-63-8	TRIFLUOROBROMOMETHANE
75-50-3	TRIMETHYLAMINE
95-63-6	TRIMETHYLBENZ /
108-05-4	VINYLACETATE
95-47-6	XYLENEB
106-42-3	XYLENEC
1330-20-7	XYLENEMIXEDISOMER
7440-66-6	ZINC
01-03-9	ZINC AND COMPOUNDS
1314-13-2	ZINC OXIDE FUME
14639-98-6	ZINCAMMONIUM CHLORIDE

Appendix H: Category 3 Chemicals with CAS Numbers

Note: This list of Category 3 chemicals does not contain the names of chemicals that have never been reported under TURA.

CAS Number	Chemical Name
76-13-1	EREON113
110-17-8	
100-00-0	FURAN TETRAHYDRO-
98-01-1	
01-02-2	GLYCOL ETHERS
422.56.0	
422-50-0	
7647 01 0	
7647-01-0	
74-90-8	
7664-39-3	HYDROGENFLUORIDE
78-59-1	
4098-71-9	
67-63-0	
80-05-7	
554-13-2	
14307-35-8	
110-16-7	MALEICACID
108-31-6	MALEICANHYDRIDE
7439-96-5	MANGANESE
01-02-7	MANGANESE COMPOUNDS
126-98-7	METHACRYLONITRILE
109-86-4	METHOXYETHANOL
96-33-3	METHYLACRYLATE
79-22-1	METHYLCHLOROFORMATE
5124-30-1	METHYLENE BIS(4-CYCLOHEXYLISOCYANATE)
101-14-4	METHYLENEBISCHLORO
108-10-1	METHYLISOBUTYLKETO
80-62-6	METHYLMETHACRYLATE
1634-04-4	METHYLTBUTYLET
75-79-6	METHYLTIRCHLOROSILANE
90-94-8	MICHLERSKETONE
75-04-7	MONOETHYLAMINE
91-20-3	NAPHTHALENE
7440-02-0	NICKEL
01-09-0	NITRATE COMPOUNDS
7697-37-2	NITRICACID
10102-43-9	NITRICOXIDE
10102-44-0	
88-75-5	NITROPHENOLA
99-55-8	NITBOTOLUIDINE
924-42-5	
20325-40-0	
30525-80-4	
504-42-2	
108-05 0	
106-90-2	
7004 00 0	
1004-38-2	
10025-87-3	
01-03-3	
85-44-9	
109-06-8	PICOLINE

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Appendix H: Category 3 Chemicals with CAS Numbers

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May 29, 2009

MEMORANDUM TO:

Standing Committee on General Government

FROM:

Marta Kennedy Research Officer

SUBJECT:

Follow-up to Background Information on Bill 167, the *Toxics Reduction Act, 2009*

This memo is a follow-up to the memo provided to the Standing Committee on General Government on May 28, 2009 regarding the exclusion of metals and alloys from the application of the Massachusetts *Toxics Use Reduction Act* (Massachusetts TURA).

That memo indicated that while alloys of copper, nickel, chromium, cobalt manganese and silver had been removed from the Massachusetts Toxic and Hazardous Substances List, these metals in non-alloy form may still be subject to the application of the Massachusetts TURA, and that we were awaiting confirmation of this.

The attached document, provided by the Massachusetts Toxic Uses Reduction Institute, indicates that nickel, chromium, cobalt and manganese in non-alloy form are reportable under the Massachusetts TURA. Copper and silver are substances reportable only to the U.S. federal Environmental Protection Agency. The document also indicates that other metals are subject to TURA reporting. Note that the document states that it applies to the calendar year 2002. However, it is the list that is currently in use.¹ The full alphabetical list of Massachusetts TURA reportable substances begins on page 19 of the document.

¹ Telephone interview, Policy Analyst, Massachusetts Toxics Use Reduction Institute, 28 May 2009.

<u>AppendixB</u>

WHAT TOXIC SUBSTANCES ARE SUBJECT TO TURA REPORTING?

All of the substances listed under Section 313 of EPCRA, and all of the substances that are on the federal Superfund (CERCLA) list of chemicals are reportable under TURA, except for those chemicals that have been delisted by the Administrative Council on Toxics Use Reduction. The list of TURA chemicals subject to reporting for reports due July 1, 2002 is located at the end of this Appendix. (The column entitled "ADD" refers to the year in which the chemical was added to the TURA list.) They include:

- □ EPCRA 313 and CERCLA chemicals and chemical categories sorted by chemical name
- □ EPCRA 313 and CERCLA chemicals and chemical categories sorted by CAS number

The following chemicals have been delisted from TURA (but in certain cases are reportable under EPCRA):

Reporting Year/Chemical Delisted	Reason/Note
1994	
Barium Sulfate	EPCRA delisting and not CERCLA reportable
All copper phthalocyanine compounds that are substituted with only hydrogen and/or chlorine and/or	EPCRA delisting and not CERCLA reportable
bromine (delisted from the copper compounds category)	
High molecular weight glycol ethers	EPCRA delisting and not CERCLA reportable
1995	
Certain Metal Alloys (refer to	TURA Administrative Council
Appendix F)	delisting
Chromium (III) Oxide from	TURA Administrative Council
chromium compounds category	delisting
Ammonium Sulfate Solution	EPCRA delisting and not CERCLA
(CAS#7783-20-2)	reportable
Ammonium Nitrate Solution	EPCRA delisting and not CERCLA
(CAS#6484-52-2)	reportable
1996	
Hydroquinone (except for the	TURA Administrative Council
manufacture of the chemical)	delisting
Acetic Acid at concentrations of 12%	TURA Administrative Council
or less	delisting
(CAS#103-23-1)	EPCKA delisting and not CERCLA reportable
1997	

The 1996 EPA TRI guidance document, "List of Toxic Chemicals within the Water Dissociable Nitrate Compounds Category and Guidance for Reporting" (EPA 745-R-96-004), provides a detailed description of how to report this chemical category. The following general overview is from the guidance document:

"Chemicals within the nitrate compounds category are only reportable when in aqueous solution. All water dissociable nitrate compounds are included in the nitrate compounds category, including ammonium nitrate. Specifically listed section 313 chemicals *are not* included in threshold determinations for chemical categories such as the water dissociable nitrate compounds category. Specifically listed toxic chemicals are subject to their own individual threshold determinations. As of December 1, 1994, ammonium nitrate (solution) is not an individually listed chemical on the EPCRA section 313 list. However, ammonium nitrate is still subject to reporting under the nitrate compounds category. In addition, the aqueous ammonia from the dissociation of ammonium nitrate when in aqueous solution is subject to reporting under the ammonia listing."

The following is an example from the TRI guidance document:

\checkmark Example

In a calendar year, a facility manufactures as byproducts, 20,000 pounds of sodium nitrate and 10,000 pounds of calcium nitrate, both in aqueous solutions, and releases these solutions to wastewater streams. The total quantity of nitrate compounds manufactured by the facility is the sum of the two chemicals, or 30,000 pounds, which exceeds the manufacturing threshold quantity of 25,000 pounds. The facility therefore is required to report for the nitrate compounds category.

There are three diisocyanates that are reported individually under EPCRA, and <u>not</u> as the diisocyanate chemical category:

• Toluene-2,4-diisocyanate , (584-84-9)

Toluene-2,6-diisocyanate (91-08-7)

Toluene diisocyanate (mixed isomers) (26471-62-5)

Rules for Reporting Glycol Ethers

There are separate categories for glycol ethers defined under EPCRA and CERCLA.

(Please refer to the TRI guidance document, "List of Toxic Chemicals within the Glycol Ethers Category and Guidance for Reporting," EPA 745-R-95-006.)

K Reporting Guidance

EPCRA: Glycol Ethers are reportable as the glycol ethers category, N230. The

"... The chemical ammonium hydroxide (NH₄OH) is a misnomer. It is a common name used to describe a solution of ammonia in water (i.e., aqueous ammonia), typically a concentrated solution of 28 to 30 percent ammonia. EPA has consistently responded to questions regarding the reportability of these purported ammonium hydroxide solutions under the EPCRA Section 313 ammonia listing by stating that these are 28 to 30 percent solutions of ammonia in water and that the solutions are reportable under EPCRA Section 313 ammonia listing. For a more detailed discussion, see page 34175 of the Federal Register final rule of June 30, 1995 (60 FR 34172).

Facilities should use the percent total ammonia specified on the label of ammonium hydroxide solutions they purchase to determine the total ammonia content in these solutions. Ammonium hydroxide has the chemical formula NH_4OH ; however, as mentioned above, strong evidence indicates that the species NH_4OH does not exist. Bottles of concentrated aqueous ammonia purchased from chemical supply companies are almost always labeled ammonium hydroxide. These solutions primarily consist of molecules of NH_3 dissolved in water (along with small amounts of ionized ammonia).

Please note that different chemical suppliers will reference the % of ammonia in different ways. Therefore, it is recommended that facilities contact their chemical suppliers to specify the amount of ammonia per gallon. This question could be specified in "percent by weight per gallon" or in "number of pounds per gallon" of solution, for example.

\checkmark Examples

Example 1:

Facility otherwise uses 1,000,000 pounds of 30% solution by weight of ammonium hydroxide (30% ammonia by weight).

1,000,000 pounds x 0.30 = 300,000 pounds of ammonia

Only 10% of ammonia is reportable on Form R per EPA <u>Guidance for Reporting</u> Aqueous Ammonia

300,000 x 0.10 = 30,000 pounds of ammonia

This facility would complete one Form S and one Form R for 30,000 pounds of reportable ammonia (see EPCRA Guidance for Reporting Aqueous Ammonia).

	 1		
✓ Examples			
			<u> </u>
Example 2:	 14 - A	·	

\checkmark Example

Hydrogen Cyanide: EPCRA category (Cyanide Compounds), EPCRA specifically list Report as: EPCRA specific chemical (Hydrogen Cyanide)

1,2,4 Trichlorobenzene: EPCRA specifically listed, CERCLA category (Chlorinated Benzenes)

Report as: EPCRA specific chemical (1,2,4 Trichlorobenzene)

Row 3:

When a specifically listed CERCLA chemical falls within an EPCRA listed category, it should be reported only under the EPCRA chemical category and not under the specific chemical name.

🖉 Example

Calcium Cyanide: EPCRA category (Cyanide Compounds), CERCLA specifically listed Report as: EPCRA category (Cyanide Compounds)

Row 4:

When a specifically listed CERCLA chemical falls within a CERCLA listed category, it should be reported **only as the specific CERCLA chemical**.

🖉 Example

• Benzenesulfonyl Chloride: CERCLA specifically listed, CERCLA category (Chlorinated Benzenes)

Report as: CERCLA specific chemical (Benzenesulfonyl Chloride)

Row 5:

Chemicals falling under an EPCRA chemical category or categories which are not specifically listed under EPCRA or CERCLA, should be reported under the EPCRA chemical category or categories.

\checkmark Example: Lead Chromate, Reporting Two Competing Compound Classes

If you are reporting lead chromate you need to report under the lead compounds and chromium compounds categories

• Report the SAME weight for Lead Compounds and Chromium Compounds use (total weight of the compound) on EACH Form S.

• For Byproduct tracking, report ONLY the weight of the reportable constituent for each category.

Section 2: Optional Questions

When the amounts reported in c, d and e in Section 1 are added together, the sum will in many cases equal the sum of f and g. In other words, lines c,d and e will often form a "materials balance." If lines c,d and e are not in approximate balance, you may use this section to explain why. Indicate all the reasons that apply by entering the number of pounds on the appropriate line below (e.g., 4,000 Chemical was held in inventory).

b. Chemical was consumed or transformed					
300; 33,000					
d. Chemical is a compound					
· · ·					
	b. Chemical was consumed or transformed 300; 33,000 d. Chemical is a compound				

f. Did anything non-routine occur at your facility during the reporting year which affected the data reported? Yes No I fyes, please explain.

Section 1: Facility-Wide Use of Listed Chemical

1012 Chromium Compounds a. CAS # b. Chemical Name (Dioxin wi

b. Chemical Name (Dioxin will be assumed to be grams, decimal points may be used)

Facility-wide use of chemical identified in a. Enter the total amount (in POUNDS, except for dioxin) for each applicable category. **NOTE:** 'Generated as byproduct' (item f.) generally means all waste containing the listed chemical before the waste is treated or recycled. Please refer to the reporting instructions before completing this section.

c. Manufactured e. Otherwise used 33,000

g. Shipped in or as product

d. Processed 350 f. Generated as Byproduct

Section 2: Optional Questions

When the amounts reported in c, d and e in Section 1 are added together, the sum will in many cases equal the sum of f and g. In other words, lines c,d and e will often form a "materials balance." If lines c,d and e are not in approximate balance, you may use this section to explain why. Indicate all the reasons that apply by entering the number of pounds on the appropriate line below (e.g., 4,000 Chemical was held in inventory).

a. Onemical was recycled on site	b. Chemical was consumed or transformed	
	See above	
c. Chemical was held in inventory	d. Chemical is a compound	
o Other (avalein helew)		

Summary of TURA Reportable Chemical Categories for Calendar Year 2002

Cb.

ulas Cata - an

TURA requires reporting on the chemical categories listed below. The individual chemicals included in each chemical category should always be reported in their parent chemical category (e.g., antimony compounds), and not as individual chemicals. Please note that this is not an exhaustive list of individual chemicals within the chemical categories.

Chemical Category	CAS NO.	Chemical Name
Antimony Compounds	1309-64-4	Antimony Trioxide
Includes any unique chemical	7647-18-9	Antimony Pentachloride
substance that contains antimony as	7783-56-4	Antimony Trifluoride
Includes but is not limited to:	7789-61-9	Antimony Tribromide
morades, but is not minica to,	10025-91-9	Antimony Trichloride
	28300-74-5	Antimony Potassium Tartrate
	· · · · · · · · · · · · · · · · · · ·	
Arsenic Componds	692-42-2	Diethylarsine
Includes any unique chemical	696-28-6	Dichlorophenylarsine
substance that contains arsenic	1303-28-2	Arsenic Pentoxide
as part of the chemical's	1303-32-8	Arsenic Disulfide
infrastructure. Includes, but is	1303-33-9	Arsenic Trisulfide
not limited to:	1327-52-2	Arsenic Acid
	1327-53-3	Arsenic Trioxide
	7631-89-2	Sodium Arsenate
х. 18 х	7645-25-2	Lead Arsenate
	7778-39-4	Arsenic Acid
	7778-44-1	Calcium Arsenate
	7784-34-1	Arsenous Trichloride
	7784-40-9	Lead Arsenate
	7784-41-0	Potassium Arsenate
	7784-46-5	Sodium Arsenite
	10102-48-4	Lead Arsenate
	10124-50-2	Potassium Arsenate
	52740-16-6	Calcium Arsenate
Barium Compounds	542-62-1	Barium Cyanide
substance that contains barium as part		
of the chemical's infrastructure. Does		
not include barium sulfate, CAS #		
7727-43-7. Includes, but is not		
limited to:		
Beryllium Compounds	7787-47-5	Bervllium Chloride
Includes any unique chemical	7787-49-7	Bervllium Fluoride
substance that contains	7787-55-5	Beryllium Nitrate
	1 1 1 0 1 0 0 0	

Copper Compounds		Copper,
Includes any unique chemical	:137-29-1	bis(dimethylcarbamodithioato-s-s)-
substance that contains copper	544-92-3	Copper Cyanide
as part of the chemical's	815-82-7	Cupric Tartrate
infrastructure. Does not	3251-23-8	Cupric Nitrate
include copper phthalocyanine	5893-66-3	Cupric Oxalate
compounds that are substituted	7447-39-4	Cupric Chloride
with only hydrogen, and/or	7758-98-7	Cupric Sulfate
chlorine, and/or bromine.		
Includes, but is not limited to:	10380-29-7	Cupric Sulfate, Ammoniated

Chemical Category

CAS No.

Chemical Name

Cyanide Compounds	57-12-5	Cyanides
X^+CN^- where $X = H^+$ or any other	143-33-9	Sodium Cyanide
group where a formal dissociation	151-50-8	Potassium Cyanide
may occur. For example KCN or $CA(Cn)_2$.	460-19-5	Cyanogen
Includes, but is not limited to:	506-61-6	Potassium Silver Cyanide
	506-64-9	Silver Cyanide
	506-68-3	Cynogen Bromide
	506-77-4	Cyanogen Chloride
	542-62-1	Barium Cyanide
	544-92-3	Copper Cyanide
	557-19-7	Nickel Cyanide
-	557-21-1	Zinc Cyanide
	592-01-8	Calcium Cyanide
	592-04-1	Mercuric Cyanide
	592-85-8	Mercuric Thiocyanate
•	592-87-0	Lead Thiocyanate
	1762-95-4	Ammonium Thiocyanate

Diisocyanates	91-93-0	3,3'-Dimethoxybenzidine-4,4 ¹ di-isocyanate
Includes only the chemicals listed		3-3'-Dimethyl-4,4'-diphenylene
here.	91-97-4	diisocyanate

	· ·			
Ethylenebisdithiocarbamic acid,	hiocarbamic acid, 111-54-6 Ethylenebisdithiocarbamic acid,			
salts, esters		· · · · ·		
Includes any unique chemical				
substance that contains an EBDC or				
an EBDC salt as part of that				
chemical's infrastructure. Includes,		•		
but is not limited to:				
Certain Glycol Ethers				
Please see guidance on page B-3 of				
this document.				

Lead Compounds	301-04-2	Lead Acetate
Includes any unique chemical	592-87-0	Lead Thiocyanate
substance that contains lead as	1072-35-1	Lead Stearate
part of the chemical's	1314-87-0	Lead Sulfide
infrastructure. Includes, but is	1335-32-6	Lead Subacetate
not limited to:	7428-48-0	Lead Stearate
	7446-27-7	Lead Phophate
	7446-14-2	Lead Sulfate
	7645-25-2	Lead Arsenate
· .	7758-95-4	Lead Chloride
	7783-46-2	Lead Fluoride
	10099-74-8	Lead Nitrate
	10101-63-0	Lead Iodide
	10102-48-4	Lead Arsenate
	13814-96-5	Lead Fluoborate
	15739-80-7	Lead Sulfate
	52652-59-2	Lead Stearate
	56189-09-4	Lead Stearate

Chemical Category

CAS No.

Chemical Name

Manganese Compounds	7722-64-7	Potassium Permanganate
Includes any unique chemical	· · · · ·	
substance that contains lead as		
part of the chemical's	-	
infrastructure. Includes, but is		
not limited to:		Manganese,
	15339-36-3	bis(dimethylcarbamodithiato-s-s)
Mercury Compounds	592-04-1	Mercuric Cyanide
Includes any unique chemical	592-85-8	Mercuric Thiocyanate
substance that contains	628-86-4	Mercury Fulminate
mercury as part of the	7782-86-7	Mercurous Nitrate
chemical's infrastructure.	7783-35-9	Mercuric sulfate
Includes, but is not limited to:	10415-75-5	Mercurous Nitrate
substance that contains lead as part of the chemical's infrastructure. Includes, but is not limited to: Mercury Compounds Includes any unique chemical substance that contains mercury as part of the chemical's infrastructure. Includes, but is not limited to:	15339-36-3 592-04-1 592-85-8 628-86-4 7782-86-7 7783-35-9 10415-75-5	Manganese, bis(dimethylcarbamodithiato-s-s) Mercuric Cyanide Mercuric Thiocyanate Mercury Fulminate Mercurous Nitrate Mercuric sulfate Mercurous Nitrate

	189-64-0	Benzo(a,h)pyrene
	191-30-0	Dibenzo(a,l)pyrene
	192-65-4	Dibenzo(a,e)pyrene
	193-39-5	Indeno[1,2,3-cd]pyrene
	194-59-2	7H-Dibenzo(c,g)carbazole
	205-99-2	Benzo(b)fluoranthene
	205-82-3	Benzo(j)fluoranthene
	206-44-0	Benzo(j.k)fluorene
	207-08-9	Benzo(k)fluoranthene
	218-01-9	Benzo(a)phenanthrene
	224-42-0	Dibenz(a,j)acridine
	226-36-8	Dibenz(a,h)acridine
	3697-24-3	5-Methylchrysene
	5385-75-1	Dibenzo(a,e)Fluoranthene
	5522-43-0	1-Nitropyrene
		•
Selenium Compounds		Carbamodithioic Acid, dimethyl-
Includes any unique chemical		,Tetraanhydrosulfid with
substance that contains	144-34-3	Orthothioselenious Acid
selenium as part of the	630-10-4	Selenourea.
chemical's infrastructure.	7446-08-4	Selenium Dioxide

		or mountoberonious riord				
selenium as part of the	630-10-4	Selenourea .				
chemical's infrastructure.	7446-08-4	Selenium Dioxide				
Includes, but is not limited to:	7488-56-4	Selenium sulfide				
	7782-82-3	Sodium Selenite				
	7783-00-8	Selenious Acid				
	10102-18-8	Sodium Selenite				
· · · · · · · · · · · · · · · · · · ·	12039-52-0	Selenious Acid, Dithallium (1+) Salt				

Silver Compounds	506-61-6	Potassium Silver Cyanide
Includes any unique chemical	506-64-9	Silver Cyanide
substance that contains silver		
as part of the chemical's		
infrastructure. Includes, but is		
not limited to:		
	7761-88-8	Silver Nitrate

Strychnine and Salts	57-24-9	Strychnine and Salts	
Includes any unique chemical	•		
substance that contains			
strychnine or a strychnine salt			
as part of the chemical's		-	
infrastructure. Includes, but is	•		
not limited to:	·	• *	



The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2002

Sorted Aphabetically by Name

CAS	<u>NAME</u>	С	313_	ADD		CAS	NAME	C	313	ADD
1341-49-7	Ammonium bifluoride	С		1992		7647-18-9	Antimony pentachloride	С	*	199:
10192-30-0	Ammonium bisulfite	С		1993	· .	28300-74-5	Antimony potassium tartrate	С	*	199:
1111-78-0	Ammonium carbamate	С		1992		7789-61-9	Antimony tribromide	С	*	199:
506-87-6	Ammonium carbonate	С		1992		10025-91-9	Antimony trichloride	С	*	199:
12125-02-9	Ammonium chloride	С		1993		7783-56-4	Antimony trifluoride	С	*	1993
7788-98-9	Ammonium chromate	С	*	1993		1309-64-4	Antimony trioxide	С	*	1992
3012-65-5	Ammonium citrate, dibasic	С		1992		86-88-4	Antu	C		1991
13826-83-0	Ammonium fluoborate	С		1993		12674-11-2	Aroclor 1016	С		1993
12125-01-8	Ammonium fluoride	С		1993		11104-28-2	Arocior 1221	С		1993
1336-21-6	Ammonium hydroxide	С	·	1992		11141-16-5	Aroclor 1232	С		1993
5972-73-6	Ammonium oxalate	С		1992		53469-21-9	Aroclor 1242	C		1993
6009-70-7	Ammonium oxalate	С		1992		12672-29-6	Aroclor 1248	С		1993
14258-49-2	Ammonium oxalate	С		1993		11097-69-1	Aroclor 1254	С	1	1993
131-74-8	Ammonium picrate	С		1991		11096-82-5	Aroclor 1260	C		1993
16919-19-0	Ammonium silicofluoride	С		1993		7440-38-2	Arsenic	С	313	1990
7773-06-0	Ammonium sulfamate	С		1993	-	1327-52-2	Arsenic acid	С	*	1992
12135-76-1	Ammonium sulfide	С		1993		7778-39-4	Arsenic acid	С	*	1993
10196-04-0	Ammonium sulfite	С		1993		1001	Arsenic Compounds	С	N02	
3164-29-2	Ammonium tartrate	С		1992					0	
14307-43-8	Ammonium tartrate	С		1993		1303-32-8	Arsenic disulfide	С	*	1992
1762-95-4	Ammonium thiocyanate	C	*	1992		1303-28-2	Arsenic pentoxide	С	*	1992
7803-55-6	Ammonium vanadate	C		1993		1327-53-3	Arsenic trioxide	С	*	1992
628-63-7	Amyl acetate	C		1992	•	1303-33-9	Arsenic trisulfide	С	*	1992
101-05-3	Anilazine	†	313	1995		1327-53-3	Arsenous oxide	С		1992
62-53-3	Aniline	c	313	1990		7784-34-1	Arsenous trichloride	С	*	199:
120-12-7	Anthracene	c	313	1990		1332-21-4	Asbestos (friable)	C	313	1990
7440-36-0	Antimony	c	313	1990		1912-24-9	Atrazine		313	199
1000	Antimony Compounds	С	N01			492-80-8	Auramine	С	X	1990
			0			71751-41-2	Avermectin B1		X	199

CAS: Chemical Abstract Service Registry Number

313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,* : member of a chemical category – should not be reported as an individual chemical.
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CAS	NAME	C	313	ADD
7787-47-5	Beryllium chloride	С	*	1993
1003	Beryllium Compounds	С	N05 0	
7787-49-7	Beryllium fluoride	С	*	1993
7787-55-5	Beryllium nitrate	С	*	1993
13597-99-4	Beryllium nitrate	С	*	1993
33213-65-9	beta - Endosulfan	С		1993
319-85-7	beta-BHC	С		1992
91-59-8	beta-Naphthylamine	С	313	1990
57-57-8	beta-Propiolactone	С	313	1990
82657-04-3	Bifenthrin		313	1995
1464-53-5	2,2'-Bioxirane	С	X	1990
92-52-4	Biphenyl	С	313	1990
108-60-1	Bis(2-chloro-1-methylethyl)ether	С	313	1990
111-91-1	Bis(2-chloroethoxy) methane	С	313	1991
111-44-4	Bis(2-chloroethyl) ether	С	313	1990
117-81-7	Bis(2-ethylhexyl)phthalate	С	X	1990
542-88-1	Bis(chloromethyl) ether	C ·	313	1990
97-74-5	Bis(dimethylthiocarbamoyl) sulfide (tetramethylthiurammonosulfide)	С		
38661-72-2	1,3-Bis(methylisocyanate)cyclohexane		313*	1995
10347-54-3	1,4-Bis(methylisocyanate)cyclohexane		313*	1995
56-35-9	Bis(tributyltin) oxide		313	1995
10294-34-5	Borane, trichloro-		X	1995
7637-07-2	Borane, trifluoro-		X 🕤	1995
10294-34-5	Boron trichloride		313	1995
7637-07-2	Boron trifluoride		313	1995
314-40-9	Bromacil		313	1995
53404-19-6	Bromacil, lithium salt		313	1995
7726-95-6	Bromine		313	1995

CAS	NAME	С	313 A	DD
598-31-2	Bromoacetone	C		1992
35691-65-7	1-Bromo-1-(bromomethyl)-1,3-		313	1995
	propanedicarbonitrile			
353-59-3	Bromochlorodifluoromethane		313	1992
75-25-2	Bromoform	C	313	1990
74-83-9	Bromomethane	C	313	1990
314-40-9	5-Bromo-6-methyl-3-(1-methylpropyl)-2,4-		Х	1995
404 55 0	(1H,3H)-pyrimidinedione		<u> </u>	
101-55-3	4-Bromophenyl phenyl ether	C		1991
75-63-8	Bromotrifluoromethane		313	1991
1689-84-5	Bromoxynil		313	1995
1689-99-2	Bromoxynil octanoate			313
52-51-7	Bronopol			X
357-57-3	Brucine	С	313	1992
106-99-0	1,3-Butadiene		C	313
78-79-5	1,3-Butadiene, 2-methyl-	C		1991
4170-30-3	2-Butenal	C	X	
123-73-9	2-Butenal, (e)-	C		199
764-41-0	2-Butene, 1,4-dichloro-	С	X	1992
123-86-4	Butyl acetate	С		199 [.]
141-32-2	Butyl acrylate		313	199(
85-68-7	Butyl benzyl phthalate	C		199(
109-73-9	Butylamine	C		199
106-88-7	1,2-Butylene oxide	С	313	199(
1114-71-2	Butylethylcarbamothioic acid S-propyl		X	199
·	ester	_		
123-72-8	Butyraldehyde		313	199(
107-92-6	Butyric acid	С		199 [,]
4680-78-8	C.I. Acid Green 3		313	199(
6459-94-5	C.I. Acid Red 114		313	199

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CAS	NAME	С	313	ADD
148-18-5	Carbamodithioic acid, diethyl-,sodium salt (sodium diethyldithiocarbamate)	C		1
144-34-3	Carbamodithioic acid, dimethyl-,	С	*	
	tetraanhydrosulfid with orthothioselenious			
	acid(selenium,			
	tetrakis(dimethyldithiocarbamate))			
2303-16-4	Carbamothioic acid, bis(1-methylethyl)-S-	C	X	
	(2,3-dichloro-2-propenyl)ester	ļ		
2008-41-5	Carbamothioic acid, bis(2-methylpropyl)-,	၂င		
	S-ethyl ester (butylate)	<u>.</u>		
52888-80-9	Carbamothioic acid, dipropyl-, S-	C		
4000 77 7	(pnenyimetnyi) ester (prosuitocarb)			
1929-77-7	Carbamothioic acid, dipropyl-, S-propyl	C		
62.25.2	Carboard		242	4000
03-23-2			313	1990
1563-66-2	Carbofuran	C	.313	1992
52888-80-9	Carbomothioic acid, dipropyl-, S-	C		
	(phenylmethyl) ester (prosulfocarb)	<u> </u>		
75-15-0	Carbon disulfide	С	313	1990
463-58-1	Carbon oxide sulfide (COS)	C	X	1990
56-23-5	Carbon tetrachloride	С	313	1990
75-44-5	Carbonic dichloride	С	X	1990
353-50-4	Carbonic difluoride	С		1992
79-22-1	Carbonochloridic acid, methylester	С	X	1991
463-58-1	Carbonyl sulfide	C	313	1990
5234-68-4	Carboxin		313	1995
120-80-9	Catechol	С	313	1990
75-69-4	CFC-11	С	X	1991
76-14-2	CFC-114		X	1991
76-15-3	CFC-115		X	1991
75-71-8	CFC-12	С	X	1991

AS	NAME	С	313 A	DD
75-72-9	9 CFC-13		X	199
2439-01-2	2 Chinomethionat		313	199
133-90-4	4 Chloramben	С	313	199(
305-03-3	3 Chlorambucil	С		199:
57-74-9	Chlordane	С	313	199(
100	Chlordane (Technical Mixture and Metabolites)	С		
115-28-0	6 Chlorendic acid		313	199!
90982-32-4	Chlorimuron ethyl	- ·	313	199
100	Chlorinated Benzenes	С		
100	Chlorinated Ethanes	С	1	
100	3 Chlorinated Naphthalene	С		
100	Chlophenols	C	N08 4	
7782-50-	5 Chlorine	С	313	199(
10049-04-4	4 Chlorine dioxide		313	199(
10049-04-4	4 Chlorine oxide (ClO2)		X	199(
494-03-	1 Chlornaphazine	C		1992
107-20-	Chloroacetaldehyde	С		199′
79-11-	3 Chloroacetic acid	C	313	199(
532-27-4	2-Chioroacetophenone	С	313	199(
101	1 Chloroalkyl Ethers	С		
4080-31-3	3 1-(3-Chloroallyl)-3,5,7-triaza-1- azoniaadamantane chloride		313	199!
51630-58-	4-Chloro-alpha-(1-		X	199!
	methylethyl)benzeneacetic acid cyano(3-			
400.00	phenoxyphenyl)methyl ester	-	240	4001
108-90-			313	1990
510-15-0			313	1990
66441-23-4	+/2-(4-((6-Chloro-2- benzovazolylen)ovy)nhenovy)nronanojo	· .	 x	199!
	INCLEVATORY CHIVAY DICHVAY DI VOGIULO		1 1	

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11115-74-5 Chromic acid C * 1993 10101-53-8 Chromic sulfate C * 1993 7440-47-3 Chromium Compounds C 313 1990 1012 Chromium Compounds C N09 0 10049-05-5 Chromous chloride C X* 1993 218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C X* 1993 7440-48-4 Cobaltous bromide C * 1993 1013 Cobaltous formate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 14017-41-5 Cobaltous sulfamate C * 1992 14017-41-5 Cobaltous sulfamate C N10 0 544-92-3 Copper compounds C N10	CAS	NAME	C	313	ADD
10101-53-8 Chromic sulfate C * 1993 7440-47-3 Chromium Compounds C 313 1990 1012 Chromium Compounds C N09 0 10049-05-5 Chromous chloride C X 1993 218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 7789-43-7 Cobaltous bromide C N1993 14017-41-5 Cobaltous sulfamate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 1992 1015 Copper - reportable to EPA ONLY 313 1990 1015 Copper cysnide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S-S)-(copper dimethyldithiocarbamate) C * 56-72-4 Coumaphos C 313 1990 1319-77-3 <t< td=""><td>11115-74-5</td><td>Chromic acid</td><td>С</td><td>*</td><td>1993</td></t<>	11115-74-5	Chromic acid	С	*	1993
7440-47-3 Chromium Compounds C 313 1990 1012 Chromium Compounds C N09 0 10049-05-5 Chromous chloride C X 1993 218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 7789-43-7 Cobaltous bromide C N193 1993 544-18-3 Cobaltous formate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C V 1993 1015 Copper - reportable to EPA ONLY 313 1990 1015 Copper compounds C N10 0 544-92-3 Copper y bis(dimethylcarbamodithioato-S-S) C N10 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S) C 1991 1900 1319-77-3 Cresol (mixed isomers) C 313	10101-53-8	Chromic sulfate	С	*	1993
1012 Chromium Compounds C N09 0 10049-05-5 Chromous chloride C * 1993 218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 7789-43-7 Cobaltous bromide C * 1993 544-18-3 Cobaltous formate C * 1993 1014 Coke Oven Emissions C * 1993 1014 Coke Oven Emissions C * 1993 1014 Coke Oven Emissions C * 1993 1015 Copper - reportable to EPA ONLY 313 1990 1015 Copper compounds C N10 0 0 0 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S) C * 1991 3001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers)	7440-47-3	Chromium	С	313	1990
Image: 10049-05-5 Chromous chloride C * 1993 218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 6 6 7789-43-7 Cobaltous bromide C * 1993 544-18-3 Cobaltous formate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 1993 1015 Copper - reportable to EPA ONLY 313 1990 1015 Copper compounds C N10 0 0 0 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S-S)-(copper dimethyldithiocarbamate) C * 1991 56-72-4 Coumaphos C 1991 8001-58-9 Creosote C 313 1990 1319-77-3 Creosote C <	1012	Chromium Compounds	С	N09	
10049-05-5 Chromous chloride C * 1993 218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 6 1993 544-18-3 Cobaltous bromide C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 1993 1014 Coke Oven Emissions C N10 0 7440-50-8 Copper - reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 0 0 0 1 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C 1991 56-72-4 Coumaphos C 313 1990 1319-77-3 Cresol (mixed isomers) C	100 10 05 5			0	
218-01-9 Chrysene C X* 1992 7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 6 1993 7789-43-7 Cobaltous bromide C * 1993 544-18-3 Cobaltous formate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 1993 1015 Copper - reportable to EPA ONLY 313 1990 1015 Copper compounds C N10 544-92-3 Copper cyanide C N10 544-92-3 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C 1991 8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 14170-30-3	10049-05-5	Chromous chloride	С	*	1993
7440-48-4 Cobalt 313 1990 1013 Cobalt Compounds C N09 6 7789-43-7 Cobaltous bromide C * 1993 544-18-3 Cobaltous formate C * 1992 14017-41-5 Cobaltous sulfamate C * 1993 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 7440-50-8 Copper – reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * 1992 56-72-4 Coumaphos C 313 1990 1319-77-3 Cresote C 313 1990 1319-77-3 Cresote C 313 1990 1319-77-3 Cresote C 313 1990 1319-77-3 Cresol (mixed isomers) C	218-01-9	Chrysene	С	X*	1992
1013 Cobalt Compounds C N09 7789-43-7 Cobaltous bromide C * 1993 544-18-3 Cobaltous formate C * 1992 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 1993 1014 Coke Oven Emissions C V 313 1990 1015 Copper – reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * 1992 56-72-4 Coumaphos C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Crotonaldehyde, (E)- C 313 1990 1319-77-3 Crotonaldehyde, (E)- C 313 1990 132-	7440-48-4	Cobalt		313	1990
7789-43-7 Cobaltous bromide C * 1993 544-18-3 Cobaltous formate C * 1992 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 1993 1015 Copper - reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 544-92-3 Copper cyanide C * 1992 137-29-1 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C 1991 56-72-4 Coumaphos C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 14170-30-3 Crotona	1013	Cobalt Compounds	С	N09 6	
544-18-3 Cobaltous formate C * 1992 14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 7440-50-8 Copper – reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * 1992 56-72-4 Coumaphos C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Crotonaldehyde, (E)- C 313 1990 120-73-9 Crotonaldehyde, (E)- C 313 1990 135-20-6 Cupferron 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	7789-43-7	Cobaltous bromide	C	*	1993
14017-41-5 Cobaltous sulfamate C * 1993 1014 Coke Oven Emissions C * 7440-50-8 Copper – reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 0 0 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * * 56-72-4 Coumaphos C 313 1990 8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Crotonaldehyde, (E)- C 313 1992 123-73-9 Crotonaldehyde, (E)- C 313 1990 98-82-8 Cumene C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-18-3	Cobaltous formate	С	*	1992
1014 Coke Oven Emissions C I 7440-50-8 Copper – reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C 1991 56-72-4 Coumaphos C 313 1990 1319-77-3 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Crotonaldehyde, (E)- C 313 1990 123-73-9 Crotonaldehyde, (E)- C 313 1990 98-82-8 Cumene C 313 1990 135-20-6 Cupferron 313 1990 135-20-6 Cupferron 313 1991 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	14017-41-5	Cobaltous sulfamate	С	*	1993
7440-50-8 Copper – reportable to EPA ONLY 313 1990 1015 Copper Compounds C N10 0 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * 56-72-4 Coumaphos C 313 1990 1319-77-3 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 1319-77-3 Crotonaldehyde, (E)- C 313 1990 14170-30-3 Crotonaldehyde, (E)- C 313 1990 98-82-8 Cumene C 313 1990 135-20-6 Cupferron 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	1014	Coke Oven Emissions	С		
1015 Copper Compounds C N10 544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * 56-72-4 Coumaphos C 1991 8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 14170-30-3 Crotonaldehyde C 313 1990 98-82-8 Cumene C 313 1990 080-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	7440-50-8	Copper – reportable to EPA ONLY		313	1990
544-92-3 Copper cyanide C * 1992 137-29-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) C * 1991 56-72-4 Coumaphos C 1991 1991 8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 4170-30-3 Crotonaldehyde C 313 1992 123-73-9 Crotonaldehyde, (E)- C 313 1990 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	1015	Copper Compounds	С	N10	
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S)-(copper dimethyldithiocarbamate) Image: March and the symbol 56-72-4 Coumaphos C 1991 8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 4170-30-3 Crotonaldehyde C 313 1992 123-73-9 Crotonaldehyde, (E)- C 313 1990 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3	Copper cyanide	С	U *	1992
56-72-4 Coumaphos C 1991 8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 4170-30-3 Crotonaldehyde C 313 1992 123-73-9 Crotonaldehyde, (E)- C 313 1992 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1	Copper cyanide Copper, bis(dimethylcarbamodithioato-S-	C C	U * *	1992
8001-58-9 Creosote C 313 1990 1319-77-3 Cresol (mixed isomers) C 313 1990 4170-30-3 Crotonaldehyde C 313 1992 123-73-9 Crotonaldehyde, (E)- C 313 1991 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate)	C C	*	1992
1319-77-3 Cresol (mixed isomers) C 313 1990 4170-30-3 Crotonaldehyde C 313 1992 123-73-9 Crotonaldehyde, (E)- C 1991 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos	C C C	*	1992 1991
4170-30-3 Crotonaldehyde C 313 1992 123-73-9 Crotonaldehyde, (E)- C 1991 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote	C C C C	313	1992 1991 1990
123-73-9 Crotonaldehyde, (E)- C 1991 98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Cresol (mixed isomers)	C C C C C	313 313	1992 1991 1990 1990
98-82-8 Cumene C 313 1990 80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3 4170-30-3	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Cresol (mixed isomers) Crotonaldehyde	C C C C C C	0 * 313 313 313	1992 1991 1990 1990 1992
80-15-9 Cumene hydroperoxide C 313 1990 135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3 4170-30-3 123-73-9	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Cresol (mixed isomers) Crotonaldehyde Crotonaldehyde, (E)-	C C C C C C C C	313 313 313	1992 1991 1990 1990 1992 1991
135-20-6 Cupferron 313 1990 142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3 4170-30-3 123-73-9 98-82-8	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Cresol (mixed isomers) Crotonaldehyde Crotonaldehyde, (E)- Cumene	C C C C C C C C C C C	313 313 313 313 313	1992 1991 1990 1990 1992 1991 1990
142-71-2 Cupric acetate C 1991 12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3 4170-30-3 123-73-9 98-82-8 80-15-9	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Cresol (mixed isomers) Crotonaldehyde Crotonaldehyde, (E)- Cumene Cumene hydroperoxide	C C C C C C C C C C C C C C C C C C C	313 313 313 313 313 313 313	1992 1991 1990 1990 1992 1991 1990 1990
12002-03-8 Cupric acetoarsenite C 1993	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3 4170-30-3 123-73-9 98-82-8 80-15-9 135-20-6	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Cresol (mixed isomers) Crotonaldehyde Crotonaldehyde, (E)- Cumene Cumene hydroperoxide Cupferron	C C C C C C C C C	313 313 313 313 313 313 313 313	1992 1991 1990 1990 1992 1991 1990 1990
	544-92-3 137-29-1 56-72-4 8001-58-9 1319-77-3 4170-30-3 123-73-9 98-82-8 80-15-9 135-20-6 142-71-2	Copper cyanide Copper, bis(dimethylcarbamodithioato-S- S)-(copper dimethyldithiocarbamate) Coumaphos Creosote Creosote Cresol (mixed isomers) Crotonaldehyde Crotonaldehyde, (E)- Cumene Cumene hydroperoxide Cupferron Cupferron	C C C C C C C C C C C C C C C C C C C	313 313 313 313 313 313 313 313	1992 1991 1990 1990 1992 1991 1990 1990

CA	AS	NAME	С	313	ADD
L	7447-39-4	Cupric chloride	С		199
	3251-23-8	Cupric nitrate	C	*	199
	5893-66-3	Cupric oxalate	С	*	199
Γ	7758-98-7	Cupric sulfate	С	. *	1993
Γ	10380-29-7	Cupric sulfate, ammoniated	С	· *	199
÷ .	815-82-7	Cupric tartrate	С	*	1992
	21725-46-2	Cyanazine		313	199
	1016	Cyanide Compounds	С	N10 6	
	57-12-5	Cyanides (soluble salts and complexes)	С	*	199
L	460-19-5	Cyanogen	С	*	1992
	506-68-3	Cyanogen bromide	С	*	1992
	506-77-4	Cyanogen chloride	С	*	1992
	506-77-4	Cyanogen chloride ((CN)Cl)	С	*	1992
	11.34-23-2	Cycloate		313	199
	68-76-8	2,5-Cyclohexadiene-1,4-dione, 2,3,5-tris(1- aziridinyl)-		x	1990
	110-82-7	Cyclohexane	С	313	1990
	2556-36-7	1,4-Cyclohexane diisocyanate	1.	313*	199
., .	58-89-9	Cyclohexane, 1,2,3,4,5,6-hexachloro- ,(1.alpha.,2.alpha.,3.beta.,4.alpha.,5.alpha.,6 .beta.)-	С	x	199(
Ŀ	108-93-0	Cyclohexanol		313	199
	108-94-1	Cyclohexanone	С		199 [.]
	131-89-5	2-Cyclohexyl-4,6-dinitrophenol	С		199 [,]
	50-18-0	Cyclophosphamide	С		199 [,]
	68359-37-5	Cyfluthrin		313	199
	68085-85-8	Cyhalothrin		313	199!
	94-75-7	2,4-D	C	313	199(
	20830-81-3	Daunomycin	С		199:

CAS: Chemical Abstract Service Registry Number

313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,*: member of a chemical category – should not be reported as an individual chemical.
C: CERCLA Chemical – If a chemical is noted as ONLY a CERCLA chemical, then a STATE ONLY Form R as well as a Form S must be completed and submitted to DEP with your toxics use report.

NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

CAS	NAME	С	313	ADD
84-74-2	Dibutyl phthalate	С	313	1990
1918-00-9	Dicamba	С	313	1992
1194-65-6	Dichlobenil	С		1992
117-80-6	Dichlone	C		1991
99-30-9	Dichloran		313	1995
52645-53-1	3-(2,2-Dichloroethenyl)-2,2-		X	1995
	dimethylcyclopropane carboxylic acid, (3-			
00050 07 5	phenoxy-phenyl)methyl ester			
68359-37-5	3-(2,2-Dicnioroetnenyi)-2,2-	-1	X	
× .	cvano(4-fluoro-3-nhenovynhenyl)methyl			
	ester			
25321-22-6	Dichlorobenzene	С	X	1990
25321-22-6	Dichlorobenzene (mixed isomers)	С	313	1990
95-50-1	1,2-Dichlorobenzene	C	313	1990
541-73-1	1,3-Dichlorobenzene	С	313	1990
106-46-7	1,4-Dichlorobenzene	C	°313	1990
· 1018	Dichlorobenzidine	С		
91-94-1	3,3'-Dichlorobenzidine	С	313	1990
612-83-9	3,3'-Dichlorobenzidine dihydrochloride		313	1995
64969-34-2	3,3'-Dichlorobenzidine sulfate		313	1995
75-27-4	Dichlorobromomethane	С	313	1990
764-41-0	1,4-Dichloro-2-butene	C	313	1992
1717-00-6	1,1-Dichloro-1-fluoroethane		313	
1649-08-7	1,2-Dichloro-1,1-difluoroethane		313	1995
75-71-8	Dichlorodifluoromethane	С	313	1991
75-34-3	1,1-Dichloroethane	C	Х	1991
107-06-2	1,2-Dichloroethane	C	313	1990
75-35-4	1,1-Dichloroethylene	C	Х	1990
156-60-5	1,2-Dichloroethylene	С		1991

CAS	NAME	С	313 A	DD
540-59-0	1,2-Dichloroethylene		313	199(
111-44-4	Dichloroethyl ether	С	X	199(
75-43-4	Dichlorofluoromethane		313	
108-60-1	Dichloroisopropyl ether	С	X	199(
75-09-2	Dichloromethane	C	313	199(
1918-00-9	3,6-Dichloro-2-methoxybenzoic acid	С	X	1992
1982-69-0	3,6-Dichloro-2-methoxybenzoic acid, sodium salt		X	199:
542-88-1	Dichloromethyl ether	С	X	199(
19666-30-9	3-(2,4-Dichloro-5-(1-methylethoxy)phenyl)- 5-(1,1-dimethylethyl)-1,3,4-oxadiazol-2(3H)- one		X	199;
101-05-3	4,6-Dichloro-N-(2-chlorophenyl)-1,3,5- triazin-2-amine		x	199!
99-30-9	2,6-Dichloro-4-nitroaniline	_	X	199!
422-56-0	3,3-Dichloro-1,1,1,2,2-pentafluoropropane		313	199
507-55-1	1,3-Dichloro-1,1,2,2,3-pentafluoropropane		313	199!
136013-79-1	1,3-Dichloro-1,1,2,3,3-pentafluoropropane		313	199
422-48-0	2,3-Dichloro-1,1,1,2,3-pentafluoropropane		313	199!
120-83-2	2,4-Dichlorophenol	С	313	199(
50471-44-8	3-(3,5-Dichlorophenyl)-5-ethenyl-5-methyl- 2,4-oxazolidinedione		X	199
35554-44-0	1-(2-(2,4-Dichlorophenyl)-2-(2- propenyloxy)ethyl)-1H-imidazole		X	199!
75-99-0	2,2-Dichloropropionic acid	С		199 [.]
60207-90-1	1-(2-(2,4-Dichlorophenyl)-4-propyl-1,3- dioxolan-2-yl)-methyl-1H-1,2,4,-triazole		x	199!
127564-92-5	Dichloropentafluoropropane		313	199
13474-88-9	1,1-Dichloro-1,2,2,3,3-pentafluoropropane		313	199
111512-56-2	1,1-Dichloro-1,2,3,3,3-pentafluoropropane		313	199
422-44-6	1,2-Dichloro-1,1,2,3,3-pentafluoropropane		313	199!

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CAS	NAME	С	313	ADD
20325-40-0	3,3'-Dimethoxybenzidine dihydrochloride		313	1995
111984-09-9	3,3'-Dimethoxybenzidine hydrochloride		313	1995
91-93-0	3,3'-Dimethoxybenzidine-4,4'-diisocyanate		313*	1995
2300-66-5	Dimethylamine dicamba	·	313	1995
124-40-3	Dimethylamine	С	313	1991
60-11-7	4-Dimethylaminoazobenzene	С	313	1990
60-11-7	Dimethylaminoazobenzene	С	X	1990
57-97-6	7,12-Dimethylbenz[a]anthracene	С	313*	1991
119-93-7	3,3'-Dimethylbenzidine	С	313	1990
612-82-8	3,3'-Dimethylbenzidine dihydrochloride		313	1995
41766-75-0	3,3'-Dimethylbenzidine dihydrofluoride		313	1995
22781-23-3	2,2-Dimethyl-1,3-benzodioxol-4-ol		Х	1995
-	methylcarbamate	<u> </u> .		
79-44-7	Dimethylcarbamyl chloride	С	313	1990
2524-03-0	Dimethyl chlorothiophosphate		313	1995
91-97-4	3,3'-Dimethyl-4,4'-diphenylene diisocyanate		313*	1995
139-25-3	3,3'-Dimethyldiphenylmethane-4,4'-		313*	1995
	diisocyanate			
68-12-2	Dimethylformamide	С	Х	1995
57-14-7	Dimethylhydrazine	С	X	1990
57-14-7	1,1-Dimethyl hydrazine	С	313	1990
7696-12-0	2,2-Dimethyl-3-(2-methyl-1-	X	199	
	propenyl)cyclopropanecarboxylic acid		5	
	(1,3,4,5,6,7-hexahydro-1,3-dioxo-2H-		, I	
	isoindol-2-yl)methyl ester			
26002-80-2	2,2-Dimethyl-3-(2-methyl-1-		X	1995
	propenyl)cyclopropanecarboxylic acid (3-			
105.67.0	phenoxyphenyl)methyl ester		040	4000
105-07-9		C	313	1990
2524-03-0	Dimethyl phosphorochloridothioate		X	1995
131-11-3	Dimethyl phthalate	С	313	1990

CAS	NAME	С	313 A	DD
77-78-1	Dimethyl sulfate	С	313	1990
25154-54-5	Dinitrobenzene (mixed isomers)	C		1993
88-85-7	Dinitrobutyl phenol	C	313	1991
534-52-1	Dinitrocresol	C	X	1992
329-71-5	2,5-Dinitrophenol	С	<u>``</u>	1992
573-56-8	2,6-Dinitrophenol	С		1992
606-20-2	2,6-Dinitrotoluene	С	313	199(
534-52-1	4,6-Dinitro-o-cresol	С	313	199(
534-52-1	4,6-Dinitro-o-cresol and salts	С		1992
25550-58-7	Dinitrophenol	С		199:
51-28-5	2,4-Dinitrophenol	С	313	199(
121-14-2	2,4-Dinitrotoluene	С	313	199(
610-39-9	3,4-Dinitrotoluene	С		1992
25321-14-6	Dinitrotoluene (mixed isomers)	C	313	199(
117-84-0	Di-n-octyl phthalate	С		199(
39300-45-3	Dinocap	-	313	199
88-85-7	Dinoseb	С	X	199'
621-64-7	Di-n-propylnitrosamine	С	X	199(
123-91-1	1,4-Dioxane	С	313	199(
1060	Dioxin and Dioxin like Compounds		313	
957-51-7	Diphenamid		313	
122-39-4	Diphenylamine		313	199:
. 1019	Diphenylhydrazine	С		
122-66-7	1,2-Diphenylhydrazine	С	313	199(
152-16-9	Diphosphoramide, octamethyl-	С		199'
2164-07-0	Dipotassium endothall		313	199
136-45-8	Dipropyl isocinchomeronate		313	199
142-84-7	Dipropylamine	C		199 ⁻
19044-88-3	4-(Dipropylamino)-3,5-		X	199!

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The Massachusetts	Toxics	Use Reduc	ction Act	t, Reportab	le Chemical List f	or Calendar Year 2002
	•	~				

Sorted Aphabetically by Name

CAS	NAME	С	313	ADD
	methylethyl)amino]phosphinothioyl]oxy)			
	benzoic acid 1-methylethyl ester			1001
141-78-6	Ethyl acetate	C		1991
140-88-5	Ethyl acrylate	С	313	1990
31218-83-4	3-		X	1995
	((Ethylamino)methoxyphosphinothioyi)oxy)			
100-41-4	-2-butehold acid, 1-methylethyl ester		212	1000
51 70 6			313	1990
51-79-6			X	1990
/5-00-3	Ethyl chloride		X	1990
541-41-3	Ethyl chloroformate	L	313	1990
90982-32-4	Ethyl-2-((((4-chloro-6-methoxyprimidin-2-		X	1995
107 10 0	yl)-carbonyl)-amino]sulfonyl)benzoate			
107-12-0	Ethyl cyanide	C		1991
759-94-4	Ethyl dipropylthiocarbamate	Ľ	313	1995
74-85-1	Ethylene		313	1990
111-54-6	Ethylenebisdithiocarbamic acid, salts &	С	Х*	1991
	esters	L.,		
111-54-6	Ethylenebisdithiocarbamic acid, salts and		N17	
107.45.0	esters		1	1001
107-15-3	Ethylenediamine	C		1991
60-00-4	Ethylenediamine-tetraacetic acid (EDTA)	C		1991
106-93-4	Ethylene dibromide	C	X	1990
107-06-2	Ethylene dichloride	С	X	1990
107-21-1	Ethylene glycol	С	313	1990
75-21-8	Ethylene oxide	С	313	1990
96-45-7	Ethylene thiourea	С	313	1990
60-29-7	Ethyl ether	С		1991
97-63-2	Ethyl methacrylate	С		1991
62-50-0	Ethyl methanesulfonate	C	· ·	1991
151-56-4	Ethyleneimine	С	313	1990

CAS	NAME	С	313 A	DD
75-34-3	Ethylidene Dichloride	С	313	199 [,]
52-85-7	Famphur	С	313	1991
60168-88-9	Fenarimol	1	313	199!
13356-08-6	Fenbutatin oxide		313	199!
66441-23-4	Fenoxaprop ethyl		313	199!
72490-01-8	Fenoxycarb		313	199
39515-41-8	Fenpropathrin		313	199:
55-38-9	Fenthion		313	199:
51630-58-1	Fenvalerate		313	199
14484-64-1	Ferbam		313	199:
1185-57-5	Ferric ammonium citrate	С		199;
2944-67-4	Ferric ammonium oxalate	С		1992
55488-87-4	Ferric ammonium oxalate	С		199:
7705-08-0	Ferric chloride	С	·	199:
7783-50-8	Ferric fluoride	C		199:
10421-48-4	Ferric nitrate	С	*	199:
10028-22-5	Ferric sulfate	С		199:
10045-89-3	Ferrous ammonium sulfate	С		199:
7758-94-3	Ferrous chloride	С		199;
7720-78-7	Ferrous sulfate	С	*	199:
7782-63-0	Ferrous sulfate	С		199:
	Fine mineral fibers	С		-
	Fine mineral fibers (c)	C		
69806-50-4	Fluazifop butyl		313	199
2164-17-2	Fluometuron		313	199(
206-44-0	Fluoranthene	С	*	1992
86-73-7	Fluorene	С		199′
7782-41-4	Fluorine	С	313	199:
640-19-7	Fluoroacetamide	С		1992

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The Massachusetts	Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2002
•	Sorted Anhabetically by Name

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199! 1991 199 199' 199[.] 199(199' 1992 199! 199' 199' 199! 199' 199:

199(

CAS	NAME	C	313	ADD	È C	AS	NAME	C ·	313 A	ADD
1	Hexachlorocyclohexane (all isomers) CAS	С			Γ	7664-39-3	Hydrofluoric acid	С	X	19
	608-73-1					7664-39-3	Hydrofluoric acid (conc. 50% or greater)	С	X	19
58-89-9	Hexachlorocyclohexane (gamma isomer)	C	X	1990		7647-01-0	Hydrogen chloride (anhydrous)	С	X	19
77-47-4	Hexachlorocyclopentadiene	С	313	1990	ſ	7647-01-0	Hydrogen chloride (gas only)	C	X	19
67-72-1	Hexachloroethane	C	313	1990	· [74-90-8	Hydrogen cyanide	С	313	19
1335-87-1	Hexachloronaphthalene		313	1990		7664-39-3	Hydrogen fluoride	C	313	19
70-30-4	Hexachlorophene	C	313	1991		7664-39-3	Hydrogen fluoride (anhydrous)	C	X	19
1888-71-7	Hexachloropropene	С		1992	F	7783-06-4	Hydrogen sulfide	C	313	19
757-58-4	Hexaethyl tetraphosphate	С		1992	F	80-15-9	Hydroperoxide, 1-methyl-1-phenylethyl-	C	X	19
13356-08-6	Hexakis(2-methyl-2-		x	1995		123-31-9	Hydroquinone (manufactured only)	C	313	19
822-06-0	phenylpropyl)distannoxane		212*	1005	Ī	35554-44-0	Imazalil		313	19
620-31-0	Hexamethylphosphoramido		212	1995	Γ	193-39-5	Indeno(1,2,3-cd)pyrene	С	313*	19
110 54 2) 313 V	1990		13463-40-6	Iron carbonyl (Fe(CO)5), (TB-5-11)-		X	
F4005 04 0		<u> </u>	A 242	1995	Ē	13463-40-6	Iron, pentacarbonyl-		313	19
51235-04-2	A (411 211) Durimiding diang. 5 knows 6		313	1995	ľ	123-92-2	iso-Amyl acetate	C		19
55404-19-0	methyl-3-(1-methylpropyl), lithium salt		^	1995		110-19-0	iso-Butyl acetate	C		19
133-06-2	1H-Isoindole-1.3(2H)-dione. 3a.4.7.7a-	C	x	1990	· · [78-83-1	Isobutyl alcohol	C		19
	tetrahydro-2-[(trichloromethyl)thio]-				ľ	78-81-9	iso-Butylamine	С		19
67485-29-4	Hydramethylnon		313	1995	· . †	78-84-2	lsobutyraldehyde	·	313	19
302-01-2	Hydrazine	С	313	1990	ĥ	79-31-2	iso-Butyric acid	С	1	19
10034-93-2	Hydrazine sulfate		313	1990		465-73-6	Isodrin	C	313	19
57-14-7	Hydrazine, 1,1-dimethyl-	С	X	1990	ľ	25311-71-1	Isofenphos		313	19
1615-80-1	Hydrazine, 1,2-diethyl-	С		1992	ľ	55-91-4	Isofluorphate	С	1	19
540-73-8	Hydrazine, 1,2-dimethyl-	С	1.	1992	F	78-59-1	Isophorone	С	1.	19
122-66-7	Hydrazine, 1,2-diphenyl-	С	X	1990	t t	4098-71-9	Isophorone diisocyanate		313*	19
60-34-4	Hydrazine, methyl-	С	X	1990	-	78-79-5	Isoprene	C	<u>†</u>	19
122-66-7	Hydrazobenzene	С	X	1990	ŀ	42504-46-1	Isopropanolamine dodecylbenzene	С	1	19
7647-01-0	Hydrochloric acid	С		1990			sulfonate			
74-90-8	Hydrocyanic acid	С	X	1990	. • [67-63-0	Isopropyl alcohol (mfg-strong acid		313	19

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The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2002

Sorted Aphabetically by Name

CAS	NAME	С	313	ADD
7782-86-7	Mercurous nitrate	С	*	1993
10415-75-5	Mercurous nitrate	С	*	1993
7439-97-6	Mercury	С	313	1990
1028	Mercury Compounds	С	N45	
		_	8	
628-86-4	Mercury fulminate	С	*	1992
150-50-5	Merphos		313	1995
126-98-7	methacrylonitrile	С	313	1991
137-42-8	Metham sodium		313	1995
74-89-5	Methanamine	С		1991
75-50-3	Methanamine, N,N-dimethyl-	С		1991
124-40-3	Methanamine, N-methyl-	С	X	1991
62-75-9	Methanamine, N-methyl-N-nitroso-	С	X	1990
74-87-3	Methane, chloro-	С	X	1990
107-30-2	Methane, chloromethoxy-	С	X	1990
624-83-9	Methane, isocyanato-	С	X	
542-88-1	Methane, oxybis[chloro-	С	X	1990
509-14-8	Methane, tetranitro-	С	ı.	1992
67-66-3	Methane, trichloro-	С	X	1990
594-42-3	Methanesulfenyl chloride, trichloro-	С		1992
74-93-1	Methanethiol	C	X	1991
17702-57-7	Methanimidamide, N,N-dimethyl-N-[2-	С		·
	methýl-4-			
	[[(methylaino)carbonyl]oxy]phenol]-			
	(Formparanate)		ļ	
.23422-53-9	Methanimidamide, N,N-dimethyl-N-[3-	C		
	[[(methylamino)carbonyl]oxylphenyl]-			
	,mononydrochioride (formetanate			
07.50.4	nyarocnioriae)		240	1000
67-56-1	wetnanoi		313	1990
57-74-9	4,7-Methanoindan, 1,2,3,4,5,6,7,8,8-	C	X	1990

AS	NAME	С	313	ADD
	octachloro-2,3,3a,4,7,7a-hexahydro-			
91-80-5	Methapyrilene	С		199'
20354-26-1	Methazole		313	199
2032-65-7	Methiocarb	С	313	1992
16752-77-5	Methomyl	C		199:
94-74-6	Methoxone		313	199!
3653-48-3	Methoxone sodium salt		313	199!
72-43-5	Methoxychlor	С	313	199(
109-86-4	2-Methoxyethanol		313	199(
101200-48-0	2-(4-Methoxy-6-methyl-1,3,5-triazin-2-yl)-		x	199:
	metnylamino)carbonyl)amino)sulfonyl)-,			
96-33-3	Methyl acrylate		313	199(
74-83-9	Methyl bromide	С	X	1990
56-49-5	3-Methylcholanthrene	C	. *	199
74-87-3	Methyl chloride	С	X	199(
79-22-1	Methyl chlorocarbonate	С	313	199 ⁻
71-55-6	Methyl chloroform	С	X	199(
79-22-1	methyl chloroformate	С	X	199'
3697-24-3	5-Methylchrysene	•	313*	199!
75790-84-0	4-Methyldiphenylmethane-3,4-diisocyanate		313*	199
2439-01-2	6-Methyl-1,3-dithiolo[4,5-b]quinoxalin-2-one		X	199!
101-14-4	4,4'-Methylenebis(2-chloroaniline)	С	313	199(
97-23-4	2,2'-Methylenebis(4-chlorophenol)		X	199
101-61-1	4,4'-Methylenebis(N,N-	1	313	199(
	dimethyl)benzenamine			· ·
101-77-9	4,4'-Methylenedianiline	C	313	199(
78-93-3	Methyl ethyl ketone	С	313	199(
78-93-3	Methyl ethyl ketone (MEK)	С	X	1990
1338-23-4	Methyl ethyl ketone peroxide	С		1992

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CAS	NAME	C	313	ADD
	(trifluoromethyl) benzenamine			
117-84-0	n-Dioctylphthalate	С		1990
834-12-8	N-Ethyl-N'-(1-methylethyl)-6-(methylthio)- 1.3.5-triazine-2.4-diamine		X ·	1995
110-54-3	n-Hexane	c	313	1995
7440-02-0	Nickel	C	313	1990
15699-18-0	Nickel ammonium sulfate	C	*	1993
13463-39-3	Nickel carbonyl	С	*	1993
7718-54-9	Nickel chloride	C	*	1993
37211-05-5	Nickel chloride	С	*	1993
1029	Nickel Compounds	С	N49 5	
557-19-7	Nickel cyanide	С	*	1992
12054-48-7	Nickel hydroxide	С	*	1993
14216-75-2	Nickel nitrate	С	*	1993
7786-81-4	Nickel sulfate	С	*	1993
54-11-5	Nicotine	С	*	1991
1055	Nicotine and salts	7	N50 3	
1929-82-4	Nitrapyrin		313	1995
1090	Nitrate compounds (water dissociable)		N51 1	
7697-37-2	Nitric acid	С	313	1990
7697-37-2	Nitric acid (conc 80% or greater)	С	X	1990
10102-43-9	Nitric oxide	С		1993
139-13-9	Nitrilotriacetic acid		313	1990
98-95-3	Nitrobenzene	С	313	1990
92-93-3	4-Nitrobiphenyl	С	313	1990
1836-75-5	Nitrofen		313	1990
10102-44-0	Nitrogen dioxide	С		1993

The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Y	7ear 2002
Sorted Aphabetically by Name	

CAS	NAME	С	313	ADD
10544-72-6	Nitrogen dioxide	С		199
51-75-2	Nitrogen mustard		313	199
10102-43-9	Nitrogen oxide (NO)	С	· .	199
55-63-0	Nitroglycerin	С	313	199
25154-55-6	Nitrophenol (mixed isomers)	С		199
1030	Nitrophenols	C		
88-75-5	2-Nitrophenol	С	313	199
79-46-9	2-Nitropropane	С	313	199
5522-43-0	1-Nitropyrene		313*	199
1031	Nitrosamines	C		
62-75-9	Nitrosodimethylamine	С	X	199
1321-12-6	Nitrotoluene	С		199
872-50-4	N-Methyl-2-pyrrolidone		313	199
924-42-5	N-Methylolacrylamide		313	199
99-59-2	5-Nitro-o-anisidine		313	199
99-55-8	5-Nitro-o-toluidine	С	313	199
100-02-7	4-Nitrophenol	С	313	199
1116-54-7	N-Nitrosodiethanolamine	С		199
55-18-5	N-Nitrosodiethylamine	С	313	199
62-75-9	N-Nitrosodimethylamine	<u> </u>	313	199
924-16-3	N-Nitrosodi-n-butylamine	C	313	199
621-64-7	N-Nitrosodi-n-propylamine	С	313	199
86-30-6	N-Nitrosodiphenylamine	С	313	199
4549-40-0	N-Nitrosomethylvinylamine	С	313	199
59-89-2	N-Nitrosomorpholine	С	313	199
759-73-9	N-Nitroso-N-ethylurea	С	313	199
684-93-5	N-Nitroso-N-methylurea	С	313	199
615-53-2	N-Nitroso-N-methylurethane	С		199
16543-55-8	N-Nitrosonornicotine		313	199

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CAS	NAME	С	313	ADD
95-69-2	p-Chloro-o-toluidine		313	1995
104-12-1	p-Chlorophenyl isocyanate	-	313	1995
82-68-8	PCNB	С	X	1990
120-71-8	p-Cresidine		313	1990
106-44-5	p-Cresol	С	313	1990
87-86-5	PCP	С	X	1990
100-25-4	p-Dinitrobenzene	С	313	1990
1114-71-2	Pebulate		313	1995
40487-42-1	Pendimethalin		313	1995
608-93-5	Pentachlorobenzene	C		1992
76-01-7	Pentachloroethane	С	313	1991
82-68-8	Pentachloronitrobenzene	С	Х	1990
87-86-5	Pentachlorophenol	С	313	1990
504-60-9	1,3-Pentadiene	С		1992
1120-71-4	1,3-Propane sultone	С	X	1990
57-33-0	Pentobarbital sodium		313	1995
79-21-0	Peracetic acid		313	1990
127-18-4	Perchloroethylene	С	X	1990
594-42-3	Perchloromethyl mercaptan	С	313	1992
52645-53-1	Permethrin		313	1995
62-44-2	Phenacetin	С		1991
85-01-8	Phenanthrene	C .	313	1991
108-95-2	Phenol	С	313	1990
114-26-1	Phenol, 2-(1-methylethoxy)-,	С	Х	1990
	methylcarbamate	<u> </u>		
64-00-6	Phenol, 3-(1-methylethyl)-, methyl	C		
0001.07.0	carbamate (m-Cumenyi metnyicarbamate)	-		
2631-37-0	rnenoi, 3-metnyi-5-(1-metnyietnyi)-, methyi	C		
26002-80-2	Phenothrin		313	1995
			L	

CAS	NAME	С	313 A	DD
72490-01-8	(2-(4-Phenoxy-phenoxy)-ethyl)carbamic		X	199
	acid ethyl ester			·
696-28-6	Phenyl dichloroarsine	C		1992
23564-06-9	(1,2-Phenylenebis(iminocarbonothioyl))		X	199:
	biscarbamic acid diethyl ester			
95-54-5	1,2-Phenylenediamine		313	199
108-45-2	1,3-Phenylenediamine		313	199:
615-28-1	1,2-Phenylenediamine dihydrochloride		313	199
624-18-0	1,4-Phenylenediamine dihydrochloride		313	199
123-61-5	1,3-Phenylene diisocyanate		313*	199:
104-49-4	1,4-Phenylene diisocyanate		313*	199
62-38-4	Phenylmercuric acetate	С		1991
62-38-4	Phenylmercury acetate	С		199′
10453-86-8	5-(Phenylmethyl)-3-furanyl)methyl 2,2-		X	199
	dimethyl-3-(2-methyl-1-			
	propenyl)cyclopropanecarboxylate	_		
90-43-7	2-Phenylphenol		313	199(
103-85-5	Phenylthiourea	С	·	199 ⁻
57-41-0	Phenytoin	· · ·	313	199(
298-02-2	Phorate	С		1992
75-44-5	Phosgene	С	313	199(
7803-51-2	Phosphine	C.	313	199:
52-68-6	Phosphonic acid, (2,2,2-trichloro-1-	С	X	199(
	hydroxyethyl)-,dimethyl ester			
7664-38-2	Phosphoric acid	С	313	199(
961-11-5	Phosphoric acid, 2-chloro-1-(2,3,5-	``	X	199(
	trichlorophenyl) ethenyl dimethyl ester			
62-73-7	Phosphoric acid, 2-dichloroethenyl	С	X	199(
	dimethyl ester			
13194-48-4	Phosphorodithioic acid O-ethyl S,S-		X]	199
	dipropyl ester		1. · ·	

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115-07-1	1-Propene		X	1990
107-13-1	2-Propenenitrile	С	X	1990
126-98-7	2-Propenenitrile, 2-methyl-	С	X	1991
107-18-6	2-Propen-1-ol	С	X	1990
31218-83-4	Propetamphos		313	1995
60207-90-1	Propiconazole		313	1995
123-38-6	Propionaldehyde	С	313	1990
79-09-4	Propionic acid	С		1991
123-62-6	Propionic anhydride	С		1991
107-12-0	Propionitrile	С		1991
542-76-7	Propionitrile, 3-chloro-	С	X	1992
114-26-1	Propoxur	С	313	1990
115-07-1	Propylene		313	1990
75-56-9	Propylene oxide	С	313	1990
75-55-8	Propyleneimine	С	313	1990
106-49-0	p-Toluidine	С		1991
106-42-3	p-Xylene	C	313	1990
129-00-0	Pyrene	С		1991
121-21-1	Pyrethrins	С		1991
121-29-9	Pyrethrins	Ç		1991
8003-34-7	Pyrethrins	С		1993
110-86-1	Pyridine	C	313	1990
54-11-5	Pyridine, 3-(1-methyl-2-pyrrolidinyl)-,(S)-	С		1991
504-24-5	Pyridine, 4-amino-	С		1992
57-47-6	Pyrrolo[2,3-b] indol-5-ol, 1,2,3,3a,8,8a-	С		
-	hexahydro-1,3a,8-trimethyl-,			
	metnyicarbamate (ester), (3aS-cis)-			
91-22-5	Quinoline	- C	312	1990
106-51-4	Quinone		313	1990
			013	1000

CAS	NAME	С	313 A	DD
82-68	-8 Quintozene	С	313	199(
76578-14	-8 Quizalofop-ethyl		313	199
50-55	-5 Reserpine	C		199
10453-86	-8 Resmethrin		313	199
108-46	-3 Resorcinol	С		199
301-12	-2 S-(2-(Ethylsulfinyl)ethyl) O,O-dimethyl ester phosphorothioic acid		X	199
78-48	-8 S,S,S-Tributyltrithiophosphate		313	199
81-07	-2 Saccharin (manufacturing)	С	313	199
81-07	-2 Saccharin and salts	С		199(
94-59	-7 Safrole	С	313	199(
626-38	-0 sec-Amyl acetate	С		199;
105-46	-4 sec-Butyl acetate	С		199'
78-92	-2 sec-Butyl alcohol		313	199(
513-49	-5 sec-Butylamine	С		199:
13952-84	-6 sec-Butylamine	С		199:
7783-00	-8 Selenious acid	С	*	199:
12039-52	-0 Selenious acid, dithallium(1+) salt	С	*	199:
7782-49	-2 Selenium	С	313	199(
10	36 Selenium Compounds	С	N72 5	
7446-08	-4 Selenium dioxide	С	*	1992
7488-56	-4 Selenium sulfide	С	*	199;
630-10	-4 Selenourea	С	*	1992
74051-80	-2 Sethoxydim		313	199
75-77	-4 Silane, chlorotrimethyl-		X	199:
75-78	-5 Silane, dichlorodimethyl-		X	199
75-79	-6 Silane, trichloromethyl-		X	199
10	95 Silica, crystalline (respirable, < 10 microns)	<u> </u>		2002
7440-22	4 Silver – file to EPA ONLY		313	199(

CAS: Chemical Abstract Service Registry Number

313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,*: member of a chemical category – should not be reported as an individual chemical.

C: CERCLA Chemical – If a chemical is noted as ONLY a CERCLA chemical, then a STATE ONLY Form R as well as a Form S must be completed and submitted to DEP with your toxics use report.

NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

6369-96-6 2,4,5-T amines C 1992 6369-97-7 2,4,5-T amines C 1992 34014-18-1 Tebuthiuron 313 1995 3383-96-8 Temephos 313 1995 3383-96-8 Temephos 313 1995 107-49-3 Tepp C 1991 5902-51-2 Terbacil 313 1995 625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl acetate C 1992 75-65-0 tert-Butyl alcohol 313 1990 75-64-9 tert-Butylamine C 1991 93-79-8 2,4,5-T esters C 1991 93-79-8 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrachlorobisphenol A 313 1990 95-94-3 1,2,4,5-Tetrachlorobenzene C 1992 <td< th=""><th>CAS</th><th>NAME</th><th>С</th><th>313</th><th>ADD</th></td<>	CAS	NAME	С	313	ADD
6369-97-7 2,4,5-T amines C 1992 34014-18-1 Tebuthiuron 313 1995 3383-96-8 Temephos 313 1995 3383-96-8 Temephos 313 1995 107-49-3 Tepp C 1991 5902-51-2 Terbacil 313 1995 625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl alcohol 313 1990 75-65-0 tert-Butylalcohol 313 1990 75-64-9 tert-Butylamine C 1991 93-79-8 2,4,5-T esters C 1992 2545-59-7 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrabromobisphenol A 313 1990 1746-01-6 2,3,7,8-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachloroethane 313 1990	6369-96-6	2,4,5-T amines	C		1992
34014-18-1 Tebuthiuron 313 1995 3383-96-8 Temephos 313 1995 107-49-3 Tepp C 1991 5902-51-2 Terbacil 313 1995 625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl acetate C 1992 75-65-0 tert-Butyl alcohol 313 1990 75-64-9 tert-Butylanine C 1991 93-79-8 2,4,5-T esters C 1991 1928-47-8 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrabromobisphenol A 313 1990 1746-01-6 2,3,7,8-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethane 313	6369-97-7	2,4,5-T amines	С		1992
3383-96-8 Temephos 313 1995 107-49-3 Tepp C 1991 5902-51-2 Terbacil 313 1995 625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl acetate C 1992 75-65-0 tert-Butyl alcohol 313 1990 75-64-9 tert-Butylanine C 1991 93-79-8 2,4,5-T esters C 1991 1928-47-8 2,4,5-T esters C 1992 2545-59-7 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1992 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrabromobisphenol A 313 1991 95-94-3 1,2,4,5-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachloroethane C 313 1992 630-20-6 1,1,1,2-Tetrachloroethane C 313 1991 127-18-4 Tetrachloroethylene C 313 1995 354-11-0 1,1,1,2-Tetrachl	34014-18-1	Tebuthiuron		313	1995
107-49-3 Tepp C 1991 5902-51-2 Terbacil 313 1995 625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl acetate C 1992 75-65-0 tert-Butyl alcohol 313 1990 75-64-9 tert-Butylanine C 1991 93-79-8 2,4,5-T esters C 1991 1928-47-8 2,4,5-T esters C 1992 2545-59-7 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1992 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrabromobisphenol A 313 1991 95-94-3 1,2,4,5-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachloroethane C 313 1992 630-20-6 1,1,1,2-Tetrachloroethane C 313 1991 127-18-4 Tetrachloroethlane C 313 1995 354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 358-90-2	3383-96-8	Temephos		313	1995
5902-51-2 Terbacil 313 1995 625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl acetate C 1992 75-65-0 tert-Butyl alcohol 313 1990 75-64-9 tert-Butyl alcohol 313 1990 75-64-9 tert-Butylamine C 1991 93-79-8 2,4,5-T esters C 1992 2545-59-7 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrabromobisphenol A 313 1992 95-94-3 1,2,4,5-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachloroethane C 313 1992 630-20-6 1,1,2.2-Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethylene C 313 1995 354-11-0 1,1,2.2-Tetrachloro-2-fluoroethane 313 1995 <td< td=""><td>107-49-3</td><td>Терр</td><td>С</td><td></td><td>1991</td></td<>	107-49-3	Терр	С		1991
625-16-1 tert-Amyl acetate C 1992 540-88-5 tert-Butyl acetate C 1992 75-65-0 tert-Butyl alcohol 313 1990 75-64-9 tert-Butylamine C 1991 93-79-8 2,4,5-T esters C 1991 1928-47-8 2,4,5-T esters C 1992 2545-59-7 2,4,5-T esters C 1992 25168-15-4 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 61792-07-2 2,4,5-T esters C 1993 79-94-7 Tetrabromobisphenol A 313 1990 95-94-3 1,2,4,5-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachloroethane C 313 1992 630-20-6 1,1,2-Tetrachloroethane C 313 1992 127-18-4 Tetrachloroethylene C 313 1995 354-11-0 1,1,2-Tetrachloro-2-fluoroethane 313 1995 354-11-0 1,1,2-Tetrachloro-2-fluoroethane 313 1995	5902-51-2	Terbacil		313	1995
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79-94-7 Tetrabromobisphenol A 313 95-94-3 1,2,4,5-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachlorodibenzo-p-dioxin (TCDD) C 1992 630-20-6 1,1,1,2-Tetrachloroethane C 313 1992 79-34-5 1,1,2,2-Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethylene C 313 1990 354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 354-11-0 1,1,2,2-Tetrachloro-2-fluoroethane 313 1995 354-11-0 1,1,2,2-Tetrachloro-2-fluoroethane 313 1995 354-11-0 1,1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetractholorvinphos 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	61792-07-2	2,4,5-T esters	С.		1993
95-94-3 1,2,4,5-Tetrachlorobenzene C 1991 1746-01-6 2,3,7,8-Tetrachlorodibenzo-p-dioxin (TCDD) C 1992 630-20-6 1,1,1,2-Tetrachloroethane C 313 1992 79-34-5 1,1,2,2-Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethylene C 313 1990 354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 354-14-3 1,1,2,2-Tetrachloro-2-fluoroethane 313 1995 354-14-3 1,1,2,2-Tetrachloro-2-fluoroethane 313 1995 354-11-0 1,1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	79-94-7	Tetrabromobisphenol A		313	
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630-20-6 1,1,2-Tetrachloroethane C 313 1992 79-34-5 1,1,2,2-Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethylene C 313 1990 354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 354-11-0 1,1,2,2-Tetrachloro-2-fluoroethane 313 1995 354-11-0 1,1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	1746-01-6	2,3,7,8-Tetrachlorodibenzo-p-dioxin (TCDD)	С		1992
79-34-5 1,1,2,2-Tetrachloroethane C 313 1990 127-18-4 Tetrachloroethylene C 313 1990 354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 354-11-0 1,1,2,2-Tetrachloro-2-fluoroethane 313 1995 354-11-0 1,1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	630-20-6	1,1,1,2-Tetrachloroethane	С	313	1992
127-18-4 Tetrachloroethylene C 313 1990 354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 354-11-0 1,1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	79-34-5	1,1,2,2-Tetrachloroethane	С	313	1990
354-14-3 1,1,2,2-Tetrachloro-1-fluoroethane 313 1995 354-11-0 1,1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	127-18-4	Tetrachloroethylene	C	313	1990
354-11-0 1,1,2-Tetrachloro-2-fluoroethane 313 1995 58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	354-14-3	1,1,2,2-Tetrachloro-1-fluoroethane		313	1995
58-90-2 2,3,4,6-Tetrachlorophenol C * 1991 961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	354-11-0	1,1,1,2-Tetrachloro-2-fluoroethane	*	313	1995
961-11-5 Tetrachlorvinphos 313 1990 64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	58-90-2	2,3,4,6-Tetrachlorophenol	C	*	1991
64-75-5 Tetracycline hydrochloride 313 1995 78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	961-11-5	Tetrachlorvinphos		313	1990
78-00-2 Tetraethyl lead C 1991 107-49-3 Tetraethyl pyrophosphate C 1991 3689-24-5 Tetraethyldithiopyrophosphate C 1992	64-75-5	Tetracycline hydrochloride		313	1995
107-49-3Tetraethyl pyrophosphateC19913689-24-5TetraethyldithiopyrophosphateC1992	78-00-2	Tetraethyl lead	C	· .	1991
3689-24-5 Tetraethyldithiopyrophosphate C 1992	107-49-3	Tetraethyl pyrophosphate	С		1991
	3689-24-5	Tetraethyldithiopyrophosphate	С		1992

CAS	NAME	С	313 A	DD
533-74-4	Tetrahydro-3,5-dimethyl-2H-1,3,5-		X	199
	thiadiazine-2-thione		· •	
53404-60-7	Tetrahydro-3,5-dimethyl-2H-1,3,5-		X	199
0710500	thiadiazine-2-thione, ion(1-), sodium			
67485-29-4	Tetrahydro-5,5-dimethyl-2(1H)-	X	199	
	pyrimidinone(3-(4-(trifluoromethyl)phenyl)-		5	
	1-(2-(4-(triffuorometnyi)pnenyi)etnenyi)-2-			
7696-12-0	Tetramethrin		212	100/
20515 41 9	2.2.2.2. Totromothylovolonronono			400/
33313-41-0	carboxylic acid evano(3-		^	199:
· ·	phenoxynhenvi)methyl ester			
509-14-8	Tetranitromethane	C		1992
1314-32-5	Thallic oxide	C	*	199;
7440-28-0	Thallium	C	313	199(
7791-12-0	Thallium chloride TICI	C	*	199;
1038	Thallium Compounds	C	N76	
			0	
10031-59-1	Thallium sulfate	С	*	199:
563-68-8	Thallium(I) acetate	С	*	1992
6533-73-9	Thallium(I) carbonate	С	*	1992
10102-45-1	Thallium(I) nitrate	С	*	199:
- 7446-18-6	Thallium(I) sulfate	С	*	1992
6533-73-9	Thallous carbonate	С	*	1992
7791-12-0	Thallous chloride	С	*	199:
7446-18-6	Thallous sulfate	С	*	1992
148-79-8	Thiabendazole	·	313	199:
148-79-8	2-(4-Thiazolyl)-1H-benzimidazole		X	199!
62-55-5	Thioacetamide	С	313	199(
28249-77-6	Thiobencarb		313	199!
139-65-1	4,4'-Thiodianiline		313	199(

CAS: Chemical Abstract Service Registry Number

313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,*: member of a chemical category – should not be reported as an individual chemical.

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NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

CAS	NAME	C	313	ADD
1582-09-8	Trifluralin	С	313	1990
26644-46-2	Triforine		313	1995
75-50-3	Trimethylamine	С		1991
95-63-6	1,2,4-Trimethylbenzene		313	1990
16938-22-0	2,2,4-Trimethylhexamethylene diisocyanate		313*	1995
15646-96-5	2,4,4-Trimethylhexamethylene diisocyanate		313*	1995
540-84-1	2,2,4-Trimethylpentane	С		
540-84-1	2,2,4-Trimethylpentane	С		
2655-15-4	2,3,5-Trimethylphenyl methylcarbamate	•	313	1995
99-35-4	1,3,5-Trinitrobenzene	С		1991
639-58-7	Triphenyltin chloride		313	1995
76-87-9	Triphenyltin hydroxide	<u> </u>	313	1995
126-72-7	Tris(2,3-dibromopropyl) phosphate	С	313	1990
14484-64-1	Tris(dimethylcarbamodithioato-S,S')iron		X	1995
72-57-1	Trypan blue	С	313	1991
13560-99-1	2,4,5-T salts	С		1993
66-75-1	Uracil mustard	С		1991
541-09-3	Uranyl acetate	С		1992
10102-06-4	Uranyl nitrate	С	. *	1993
36478-76-9	Uranyl nitrate	С	*	1993
2164-17-2	Urea, N,N-dimethyl-N'-[3-		X	
•	(trifluoromethyl)phenyl]-			
51-79-6	Urethane	C	313	1990
7440-62-2	Vanadium (except when in alloy)		313	1990
1314-62-1	Vanadium pentoxide	С		1992
1065	Vanadium Compounds		N77 0	
27774-13-6	Vanadyi sulfate	С		1993
2699-79-8	Vikane		X	1995
50471-44-8	Vinclozolin		313	1995

AS	NAME	С	313 A	ADD
108-05-4	Vinyl acetate	С	313	199
108-05-4	Vinyl acetate monomer	С	X	199
593-60-2	Vinyl bromide	С	313	199
75-01-4	Vinyl chloride	С	313	199
75-35-4	Vinylidene chloride	C	313	199
81-81-2	Warfarin	С	X*	199 [,]
1075	Warfarin and salts		N87 4	-
81-81-2	Warfarin, & salts, conc.>0.3%	С	X*	199 [.]
1330-20-7	Xylene (mixed isomers)	С	313	1990
1300-71-6	Xylenol	С		1992
87-62-7	2,6-Xylidine		313	199(
7440-66-6	Zinc (fume or dust)	С	313	1990
557-34-6	Zinc acetate	С	*	1992
14639-97-5	Zinc ammonium chloride	С	*	199:
14639-98-6	Zinc ammonium chloride	С	*	199:
52628-25-8	Zinc ammonium chloride	С	. *	1993
1332-07-6	Zinc borate	С	*	1992
7699-45-8	Zinc bromide	С	*	199:
3486-35-9	Zinc carbonate	С	*	1992
7646-85-7	Zinc chloride	С	*	199:
1039	Zinc Compounds	С	N98 2	
557-21-1	Zinc cyanide	С	*	1992
7783-49-5	Zinc fluoride	C	*	199:
557-41-5	Zinc formate	С	*	199;
7779-86-4	Zinc hydrosulfite	С	*	199:
7779-88-6	Zinc nitrate	С	*	199:
127-82-2	Zinc phenolsulfonate	С	*	199'
1314-84-7	Zinc phosphide	С		1992

CAS: Chemical Abstract Service Registry Number

313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,*: member of a chemical category - should not be reported as an individual chemical.

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NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2001

Sorted Numerically by CAS

CAS

NAME

313 ADD

С

CAS

NAME

 $C \leftarrow 313 \ ADD$

1000 Antimony Compounds	C	N010	
1001 Arsenic Compounds	С	N020	
1002 Barium Compounds		N040	
1003 Beryllium Compounds	C	N050	
1004 Cadmium Compounds	C	N078	
1005 Chlordane (Technical Mixture and M	Aetabolites) C		
1006 Chlorinated Benzenes	C		
1007 Chlorinated Ethanes	С		
1008 Chlorinated Naphthalene	C		
1011 Chioroalkyl Ethers	С		
1009 Chlorophenols	С	N084	
1012 Chromium Compounds	С	N090	
1013 Cobalt Compounds	C	N096	
1014 Coke Oven Emissions	C		
1015 Copper Compounds	C	N100	
1016 Cyanide Compounds	С	N106	
1017 DDT and Metabolites	С		
1018 Dichlorobenzidine	С		
1019 Diphenylhydrazine	C		
1020 Endosulfan and Metabolites	C		
1021 Endrin and Metabolites	C		
Fine mineral fibers	C		
Fine mineral fibers (c)	C		
1022 Glycol Ethers	C	N230	
1023 Haloethers	C		
1024 Halomethanes	С		
HCFC-141b	X		
1025 Heptachlor and Metabolites	C		
1026 Lead Compounds	C	N420	
1027 Manganese Compounds	С	N450	
1028 Mercury Compounds	C	N458	
1029 Nickel Compounds	C	N495	

	•			
1030	Nitrophenols	C		
1031	Nitrosamines	С		
1033	Phthalate Esters	С		
1034	Polybrominated Biphenyls (PBBs)		N575	
1035	Polynuclear Aromatic Hydrocarbons	C		
1036	Selenium Compounds	С	N725	
1037	Silver Compounds	С	N740	
1038	Thallium Compounds	C	N760	-
1039	С	N982		
1040	Polycyclic aromatic compounds (includes only 21 chemicals)		N590	
	Polycyclic organic matter	С		
	Polycyclic Organic Matter (e)	С		
1045	Polychlorinated alkanes (C10 to C13)		N583	
1050	Diisocyanates (includes only 20 chemicals)		N120	
1055	Nicotine and salts	С	N503	19
1060	Dioxin and Dioxin like Compounds		N150	
1065	Vanadium Compounds		N770	
1070	Strychnine, and salts	C	N746	19
1075	Warfarin and salts		N874	
1090	Nitrate compounds (water dissociable)		N511	
1095	Silica, crystalline (respirable, < 10 microns)			20
6-60-7	Picloram	Ì	313	19
30-59-3	Tributyltin fluoride		313	19
50-00-0	Formaldehyde	С	313	19
50-00-0	Formaldehyde (solution)	C	X	1
50-07-7	Mitomycin C	С		19
50-18-0	Cyclophosphamide	С		19
50-29-3	DDT	С	,	
50-32-8	Benzo[a]pyrene	С	313*	19
50-55-5	Reserpine	С		19
51-03-6	Piperonyl butoxide		313	19
51-21-8	5-Fluorouracil		X	

CAS: Chemical Abstract Service Registry Number

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NOTE: Some EPA Chemicals may have been deslisted from the US EPA's EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

APPENDIX B - Num - 1

The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar	Year 2001
Sorted Numerically by CAS	

CAS	NAME	C 313 ADD		CAS		NAME	C 3	13 AD	D	
60-29-7	Ethane, 1,1'-oxybis-	С		1991		65-85-0	Benzoic acid	С		19
60-29-7	Ethyl ether	С		1991		66-75-1	Uracil mustard	C		19
60-34-4	Hydrazine, methyl-	С	Х			67-56-1	Methanol	C	313	3 19
60-34-4	Methyl hydrazine	C	313	1990		67-63-0	Isopropyl alcohol (mfg-strong acid process)		31:	3 19
60-35-5	Acetamide	С	313	1990		67-64-1	Acetone	C	1	\top
60-51-5	Dimethoate	С	313	1991		67-66-3	Chloroform	C	313	3 19
60-57-1	Dieldrin	С		1991		67-66-3	Methane, trichloro-	С	X	19
61-82-5	Amitrole	С	313	1991		67-72-1	Hexachloroethane	С	313	3 19
62-38-4	Phenylmercuric acetate	C		1991		68-12-2	Dimethylformamide	С	X	1
62-38-4	Phenylmercury acetate	С		1991		68-12-2	N,N-Dimethylformamide	С	313	319
62-44-2	Phenacetin	С		1991		68-76-8	2,5-Cyclohexadiene-1,4-dione, 2,3,5-tris(1-		X	19
62-50-0	Ethyl methanesulfonate	C		1991		-	aziridinyl)-			
62-53-3	Aniline	С	313	1990		68-76-8	Triaziquone		313	3 19
62-55-5	Thioacetamide	С	313	1990		70-25-7	Guanidine, N-methyl-N'-nitro-N-nitroso-	С		19
62-56-6	Thiourea	С	313	1990		70-30-4	Hexachlorophene	С	313	3 19
62-73-7	Dichlorvos	С	313	1990		71-36-3	n-Butyl alcohol	С	313	3 19
62-73-7	Phosphoric acid, 2-dichloroethenyl dimethyl	С	X	1990	·	71-43-2	Benzene	С		3
	ester					71-55-6	1,1,1-Trichloroethane	С		3
62-74-8	Fluoroacetic acid, sodium salt	С	Х	1991		71-55-6	Methyl chloroform	С	X	19
62-74-8	Sodium fluoroacetate	С	313	1991		72-20-8	Endrin	С		19
62-75-9	Methanamine, N-methyl-N-nitroso-	С	Х	1990		72-43-5	Benzene, 1,1'-(2,2,2-trichloroethylidene)bis [4-	С	X	19
62-75-9	Nitrosodimethylamine	С	Х	1990			methoxy-			
62-75-9	N-Nitrosodimethylamine	С	313	1990		72-43-5	Methoxychlor	C	313	3 19
63-25-2	1-Naphthalenol, methylcarbamate	С	Х	1990	,	72-54-8	DDD	С		19
63-25-2	Carbaryl	С	313	1990	-	72-55-9	DDE	C		19
64-00-6	Phenol, 3-(1-methylethyl)-, methyl carbamate	C				72-57-1	Trypan blue	С	313	3 19
	(m-Cumenyl methylcarbamate)					74-83-9	Bromomethane	С	313	19
64-18-6	Formic acid	С	313	1991		74-83-9	Methyl bromide	С	X	19
64-19-7	Acetic acid (concentrations of 12% or less are	С		1991		74-85-1	Ethene	· ·	X	19
	NOT reportable)	·			•	74-85-1	Ethylene		313	19
64-67-5	Diethyl sulfate	C	313	1990		74-87-3	Chloromethane	C	313	19
64-75-5	Tetracycline hydrochloride		313	1995		74-87-3	Methane, chloro-	C	X	19

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NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

APPENDIX B – NUM - 3

The Massachusetts	Toxics Use Reduction A	ct, Reportable	Chemical List fo	r Calendar Year 2001
		/ 1		
	Sorted N	umerically by (TAS	

CAS	NAME	C 3	13 AD	D	CAS	0110	NAME	C 3	13 ADD
75-86-5	Acetone cyanohydrin	С	X	· ·		78-88-6	2,3-Dichloropropene	C	313 19
75-87-6	Acetaldehyde, trichloro-	С			1	78-92-2	sec-Butyl alcohol		313 19
75-88-7	2-Chloro-1,1,1-trifluoroethane		313	1995		78-93-3	Methyl ethyl ketone	С	313
75-88-7	HCFC-133a		X	1995		78-93-3	Methyl ethyl ketone (MEK)	С	X. 19
75-99-0	2,2-Dichloropropionic acid	C		1991	-	78-99-9	1,1-Dichloropropane	С	19
76-01-7	Pentachloroethane	С	313	1991	-	79-00-5	1,1,2-Trichloroethane	С	313 19
76-02-8	Trichloroacetyl chloride		313	1995	1	79-01-6	Trichloroethylene	С	313
76-06-2	Chloropicrin		313	1995	-	79-06-1	Acrylamide	С	313 19
76-13-1	Ethane, 1,1,2-trichloro-1,2,2,-trifluoro-		X	1990		79-09-4	Propionic acid	С	19
76-13-1	Freon 113		313	1990		79-10-7	Acrylic acid	С	313 19
76-14-2	CFC-114		X	1991		79-11-8	Chloroacetic acid	C	313 19
76-14-2	Dichlorotetrafluoroethane		313	1991		79-19-6	Thiosemicarbazide	С	313 19
76-15-3	CFC-115		· X			79-21-0	Ethaneperoxoic acid		X 19
76-15-3	Monochloropentafluoroethane		313	1991		79-21-0	Peracetic acid		.313 19
76-44-8	1,4,5,6,7,8,8-Heptachloro-3a,4,7,7a-tetrahydro-	C	X	1990		79-22-1	Carbonochloridic acid, methylester	С	X 19
	4,7-methano-1H-indene	<u> </u>				79-22-1	Methyl chlorocarbonate	С	313 19
76-44-8	Heptachlor	C	313	1990		79-22-1	methyl chloroformate	С	X 19
76-87-9	Triphenyltin hydroxide		313	1995		79-31-2	iso-Butyric acid	C	
77-47-4	Hexachlorocyclopentadiene	C	313	1990		79-34-5	1,1,2,2-Tetrachloroethane	С	313
77-73-6	Dicyclopentadiene		313	1995		79-44-7	Dimethylcarbamyl chloride	С	313 19
77-78-1	Dimethyl sulfate	C	313	1990		79-46-9	2-Nitropropane	C	313 19
78-00-2	Tetraethyl lead	C		1991	7	79-94-7	Tetrabromobisphenol A		
78-48-8	DEF		<u> </u>		8	30-05-7	4,4'-Isopropylidenediphenol		313 19
78-48-8	S,S,S-Tributyltrithiophosphate		313	1995	8	30-15-9	Cumene hydroperoxide	C	313 19
78-59-1	Isophorone	<u>с</u>	, ,		8	30-15-9	Hydroperoxide, 1-methyl-1-phenylethyl-	С	X 19
78-79-5	1,3-Butadiene, 2-methyl-	С		1991	8	30-62-6	Methyl methacrylate	С	313 19
78-79-5	Isoprene	С		1991	8	31-07-2	Saccharin (manufacturing)	С	313 19
78-81-9	iso-Butylamine	С		1991	8	31-07-2	Saccharin and salts	С	
78-83-1	Isobutyl alcohol	<u>с</u>			8	31-81-2	Warfarin	С	X* 19
78-84-2	Isobutyraldehyde		313	1990	8	31-81-2	Warfarin, & salts, conc.>0.3%	С	N874 19
78-87-5	1,2-Dichloropropane	C	313	1990	8	31-88-9	C.I. Food Red 15		313 19
78-87-5	Propane 1,2-dichloro-	С	X	1990	. 8	32-28-0	1-Amino-2-methylanthraquinone		313 19

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APPENDIX B – NUM - 5

The Massachusetts Toxics Use Re	eduction Act, I	Reportable Chemical I	list for Calendar Y	(ear 2001
-	Sorted Nume	rically by CAS	-	

CAS	NAME	C 3	13 AD	D	CAS	NAME	C 3	13 AD	D
94-79-1	2,4-D Esters	С		1991	1	(Disulfiram)			T
94-80-4	2,4-D butyl ester	С	313	1991	1 98-01-1	Furfural	•	C	<u> </u>
94-80-4	2,4-D Esters	С	X	1991	1 98-07-	Benzoic trichloride	С	313	3
94-82-6	2,4-DB		313	1995	5 98-07-1	Benzotrichloride	С	X	19
95-06-7	Carbamodithioic acid, diethyl-, 2-chloro-2-	С			98-09-9	Benzenesulfonyl chloride	C		19
	propenyl ester(sulfallate)				98-82-8	Cumene	С	313	19
95-47-6	Benzene, o-dimethyl-	C	X	1990	98-86-2	Acetophenone	C	313	19
95-47-6	o-Xylene	C	313	1990	98-87-3	Benzal chloride	С	313	19
95-48-7	o-Cresol	C	313	1990	98-88-4	Benzoyl chloride	C	313	119
95-50-1	1,2-Dichlorobenzene	С	313	1990	98-95-	Nitrobenzene	С	313	3 19
95-50-1	o-Dichlorobenzene	С	X	1990	99-08-	m-Nitrotoluene	С		19
95-53-4	o-Toluidine	C	313	1990	99-30-9	2,6-Dichloro-4-nitroaniline		-x-	19
95-54-5	1,2-Phenylenediamine	313	1995		99-30-9	Dichloran		313	3119
95-57-8	2-Chlorophenol	С	*	1991	99-35-4	1,3,5-Trinitrobenzene	С		19
95-63-6	1,2,4-Trimethylbenzene		313	1990	99-55-8	5-Nitro-o-toluidine	С	313	119
95-69-2	p-Chloro-o-toluidine		313	1995	5 99-59-2	5-Nitro-o-anisidine		313	119
95-80-7	2,4-Diaminotoluene	С	313	1990	99-65-0	m-Dinitrobenzene	- C	313	st
95-94-3	1,2,4,5-Tetrachlorobenzene	С		·	99-99-0	p-Nitrotoluene	c		19
95-95-4	2,4,5-Trichlorophenol	С	313	1990	100-01-6	p-Nitroaniline	C	313	119
96-09-3	Styrene oxide	С	313		100-02-7	4-Nitrophenol	C	313	19
96-12-8	1,2-Dibromo-3-chloropropane	С	313		100-02-7	p-Nitrophenol	C	X	-
96-12-8	DBCP	С	X	1990	100-25-4	p-Dinitrobenzene	C	313	19
96-18-4	1,2,3-Trichloropropane		313	1995	100-41-4	Ethylbenzene	c	313	19
96-33-3	Methyl acrylate		313	1990	100-42-	Styrene	C	313	;
96-45-7	Ethylene thiourea	С	313	1990	100-44-7	Benzyl chloride	c	313	119
97-23-4	2,2'-Methylenebis(4-chlorophenol		Х	1995	100-47-0	Benzonitrile	c	-	19
97-23-4	Dichlorophene		313	1995	100-75-4	N-Nitrosopiperidine	С	313	19
97-56-3	C.I. Solvent Yellow 3		313	1990	101-05-3	4.6-Dichloro-N-(2-chlorophenvl)-1.3.5-triazin-2-		X	19
97-63-2	Ethyl methacrylate	C				amine			
97-74-5	Bis(dimethylthiocarbamoyl) sulfide	С			101-05-3	Anilazine		313	19
	(tetramethylthiurammonosulfide)				101-14-4	4,4'-Methylenebis(2-chloroaniline)	C	313	, †
97-77-8	Thioperoxydicarbonic diamide, tetraethyl	C		·	101-14-4	MBOCA	С	X	19

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NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

APPENDIX B – NUM - 7
			Sorte	d Num	iericali	y by CAS				•
CAS	NAME	<u>C 31</u>	<u>3 AD</u>	D	<u> </u>	CAS	NAME	<u>`C 3</u> 1	<u>13 AD</u>	<u>D</u>
108-39-4	m-Cresol	С	313	1990		111-42-2	Diethanolamine	С	313	19
108-45-2	1,3-Phenylenediamine		313	1995	e.	111-44-4	Bis(2-chloroethyl) ether	С	- 313	19
108-46-3	Resorcinol	С		1991		111-44-4	Dichloroethyl ether	С	X	19
108-60-1	Bis(2-chloro-1-methylethyl)ether	С	313		1 · · [111-54-6	Ethylenebisdithiocarbamic acid, salts & esters	С	N171	19
108-60-1	Dichloroisopropyl ether	С	Х	1990		111-91-1	Bis(2-chloroethoxy) methane	С	313	19
108-88-3	Toluene	С	313	1990		114-26-1	Phenol, 2-(1-methylethoxy)-, methylcarbamate	С	X	19
108-90-7	Chlorobenzene	С	313	1990		114-26-1	Propoxur	С	313	,
108-93-0	Cyclohexanol		313	1995		115-02-6	Azaserine	C		19
108-94-1	Cyclohexanone	С		1991		115-07-1	1-Propene		X	19
108-95-2	Phenol	С	313	1990	I. [115-07-1	Propene		X	19
108-98-5	Benzenethiol	С		1991		115-07-1	Propylene		313	19
108-98-5	Thiophenol	С		1991		115-28-6	Chlorendic acid		313	19
109-06-8	2-Methylpyridine	С	313			115-29-7	Endosulfan	C ·		19
109-06-8	2-Picoline	С	X	1991		115-32-2	Benzenemethanol, 4-chloroalpha4-	С	X	19
109-73-9	Butylamine	С		1991			chlorophenyl)alpha(trichloromethyl)-			
109-77-3	Malononitrile	С	313	1991		115-32-2	Dicofol	C	313	19
109-86-4	2-Methoxyethanol		313	1990		116-06-3	Aldicarb	С	313	19
109-89-7	Diethylamine	С	-	1991		117-79-3	2-Aminoanthraquinone	· .	313	19
109-99-9	Furan, tetrahydro-	С		1991		117-80-6	Dichlone	С	<u> </u>	19
110-00-9	Furan	С		1991		117-81-7	Bis(2-ethylhexyl)phthalate	C	X	19
110-16-7	Maleic acid	С				117-81-7	DEHP	С	X .	19
110-17-8	Fumaric acid	С				117-81-7	Di(2-ethylhexyl) phthalate	C	313	19
110-19-0	iso-Butyl acetate	С		1991	ŀ	117-84-0	Di-n-octyl phthalate	C		19
110-54-3	Hexane	C	Х	1995		117-84-0	n-Dioctylphthalate	С		19
110-54-3	n-Hexane	С	313	1995		118-74-1	Hexachlorobenzene	C ·	313	19
110-57-6	trans-1,4-Dichloro-2-butene		313	1995		119-38-0	Carbamic acid, dimethyl-, 3-methyl-1-(1-	C		
110-57-6	trans-1,4-Dichlorobutene	 	X	1995			methylethyl)-1H-pyrazol-5-yl ester (isolan)			1
110-75-8	2-Chloroethyl vinyl ether	c	· ·	1991		119-90-4	3,3'-Dimethoxybenzidine	C	313	19
110-80-5	2-Ethoxyethanol	С	313	1990		119-93-7	3,3'-Dimethylbenzidine	C	313	19
110-80-5	Ethanol, 2-ethoxy-	C	Х	1990		119-93-7	o-Tolidine			X
110-82-7	Cyclohexane	С	313	1990		120-12-7	Anthracene		313	19
110-86-1	Pyridine	C	313	1990		120-36-5	2,4-DP		313	19

The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2001

CAS: Chemical Abstract Service Registry Number

313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,*: member of a chemical category – should not be reported as an individual chemical.

C: CERCLA Chemical – If a chemical is noted as ONLY a CERCLA chemical, then a STATE ONLY Form R as well as a Form S must be completed and submitted to DEP with your toxics use report.

NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

The Massachusetts Toxics Use Reduction Act	, Reportable Chemical List for Calendar Year 2001
Soutod Norm	anias line has CAS

NAME	C 3	813 AD	D	(CAS	NAME	C 3	13 AI	DD
9-2 o-Anisidine hydrochloride	· ·	313	1990	ר ה	145-73-3	Endothall	С	T	T
2-7 alpha-Naphthylamine	С	313	1990	列 [148-18-5	Carbamodithioic acid, diethyl-,sodium salt	С		\top
J-6 Benzeneamine, N-hydroxy-N-nitroso,		X	1990	ז L		(sodium diethyldithiocarbamate)			
ammonium salt					148-79-8	2-(4-Thiazolyl)-1H-benzimidazole		Х	1!
J-6 Cupferron		313	1990	\mathbf{D}	148-79-8	Thiabendazole		31	3 1
0-1 Carbamodithioic acid, dibutyl, sodium salt	С] [148-82-3	Melphalan	C		1
(Sodium dibutyldithiocarbamate)] [149-30-4	2-Mercaptobenzothiazole		31	3 1!
5-8 Dipropyl isocinchomeronate		313	1995	3 [149-30-4	MBT		Х	1!
3-8 Thiram	С	313			150-50-5	Merphos		31	3 1
J-1 Copper, bis(dimethylcarbamodithioato-S-S)-	C	*			150-68-5	Monuron	1		
(copper dimethyldithiocarbamate)				-	151-50-8	Potassium cyanide	C		* 19
J-4 Zinc, bis(dimetylcarbomodithioato-S,S)-, (ziram)	С	. *			151-56-4	Aziridine	C	X	1
1-7 Potassium N-methyldithiocarbamate		313	1995		151-56-4	Ethyleneimine	C	31	311
2-8 Metham sodium		313	1995		152-16-9	Diphosphoramide. octamethyl-			1!
2-8 Sodium methyldithiocarbamate		X	1995		156-10-5	p-Nitrosodiphenylamine		31	3 1
3-2 Disodium cyanodithioimidocarbonate		313	1995	<u>i</u> -	156-60-5	1.2-Dichloroethvlene	C		1!
3-9 Nitrilotriacetic acid		313	1990	<u>)</u> -	156-62-7	Calcium cvanamide	Ċ	31	3
5-3 3,3'-Dimethyldiphenylmethane-4,4'-diisocyanate		313*	1995	5	189-55-9	Benzo(rst)pentaphene	Ċ	313'	
5-1 4,4'-Thiodianiline		313	1990	7 F	189-55-9	Dibenzla ilpyrene		X*	1
3-5 Ethyl acrylate	С	313	1990	ŋ -	189-64-0	Dibenzo(a,h)pyrene		313'	- 19
2-2 Butyl acrylate		313	1990	7 F	191-24-2	Benzolghilpervlene	C	1.	+
3-6 Ethyl acetate	С		1991	1 -	191-30-0	Dibenzo(a.l)pyrene		313'	- 19
3-9 1,3-Dichloropropane	С		1991	1 -	192-65-4	Dibenzo(a e)pyrene		313'	19
J-6 Nabam		313	1995		193-39-5	Indeno(1 2 3-cd)pyrene	C	313'	1
I-2 Cupric acetate	С		1991	1 -	194-59-2	7H-Dibenzo(c g)carbazole		313'	-10
1-7 Dipropylamine	С		1991	1 -	196-86-9	6-Methyl-1 3-dithiolo[4 5-b]quinoxalin-2-one	1	Y	-10
3-9 Sodium cyanide (Na(CN))	С	*		1 -	196-86-9	Chinomethionat	_	31	3 10
)-0 Kepone	С			1 -	107-14-3	Dodecy/guaniding monoacotato			10
1-3 Carbamodithioic acid, dimethyl-,	С	*		1 -	197-14-3	Dodino		121	2/10
tetraanhydrosulfid with orthothioselenious					205_82.2	Benzo(i)fluoranthana		2127	
acid(selenium,					205-02-3	Benzolbifluoranthone		212*	1
tetrakis(dimethyldithiocarbamate))				」┝	200-99-2	Eluoranthono		*	+
	NAME 9-2 o-Anisidine hydrochloride 2-7 alpha-Naphthylamine 0-6 Benzeneamine, N-hydroxy-N-nitroso, ammonium salt 0-6 Cupferron 0-1 Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate) 5-8 Dipropyl isocinchomeronate 6-8 Thiram 9-1 Copper, bis(dimethylcarbamodithioato-S-S)-(copper dimethyldithiocarbamate) 0-4 Zinc, bis(dimethylcarbamodithioato-S,S)-, (ziram) 1-7 Potassium N-methyldithiocarbamate 2-8 Metham sodium 2-8 Sodium methyldithiocarbamate 3-9 Nitrilotriacetic acid 5-3 3,3'-Dimethyldiphenylmethane-4,4'-diisocyanate 5-4 4,4'-Thiodianiline 8-5 Ethyl acrylate 8-6 Ethyl acetate 8-9 1,3-Dichloropropane 9-6 Nabam 1-2 Cupric acetate 4-7 Dipropylamine 3-9 Sodium cyanide (Na(CN)) 0-0 Kepone 4-3 Carbamodithioic acid, dimethyl-, tetraanhydrosulfid with orthothioselenious acid(selenium, tetrakis(dimethyldithiocarba	NAME C 3 9-2 o-Anisidine hydrochloride 2-7 alpha-Naphthylamine C 0 0-6 Benzeneamine, N-hydroxy-N-nitroso, ammonium salt C 0 0 1 Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate) C 0 0 1 Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate) C 0 1 Carbamodithioato-Scape C 0 0 1 Carbamodithioato-Scape C 0 1 1 0 1 Copper, bis(dimethylcarbamodithioato-Scape C 0 1	NAMEC313AII9-2o-Anisidine hydrochloride3132-7alpha-NaphthylamineC3132-7alpha-NaphthylamineC3130-6Benzeneamine, N-hydroxy-N-nitroso, ammonium saltX0-6Cupferron3130-1Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate)C3-1Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate)C3-8ThiramC3139-1Copper, bis(dimethylcarbamodithioato-S-S)- (copper dimethyldithiocarbamate)C0-4Zinc, bis(dimethylcarbomodithioato-S,S)-, (ziram)C**3132-8Metham sodium3132-8Sodium methyldithiocarbamateX3-2Disodium cyanodithioimidocarbonate3133-3DimethyldithiocarbamateX3-2Disodium cyanodithioimidocarbonate3133-3Dimethyldiphenylmethane-4,4'-diisocyanate3133-1AttrakeC3133-2Butyl acrylate3133-4Ethyl acetateC8-6Ethyl acetateC3-9Sodium cyanide (Na(CN))C*3-9Sodium cyanide (Na(CN))C*3-9Sodium cyanide (Na(CN))C*3-9Sodium cyanide (Na(CN))C*3-9Sodium cyanide (Na(CN))C*3-1Carbamodithioic acid, dimethyl-, tetraanhydrosulfid with	NAME C 313 ADD 9-2 o-Anisidine hydrochloride 313 1990 2-7 alpha-Naphthylamine C 313 1990 0-6 Benzeneamine, N-hydroxy-N-nitroso, ammonium salt X 1990 0-6 Cupferron 313 1990 0-1 Carbamodithioic acid, dibutyl, sodium salt C 313 1990 0-1 Carbamodithioc acid, dibutyl, sodium salt C 313 1990 0-1 Carbamodithiocarbamate) C 313 1990 0-4 Zinc, bis(dimethylcarbamodithioato-S-S)- (copper dimethyldithiocarbamate) C * 1992 0-4 Zinc, bis(dimethylcarbamodithioato-S,S)-, (ziram) C * 1992 2-8 Metham sodium 313 1992 313 1992 2-8 Sodium methyldithiocarbamate X 1992 313 1992 3-9 Nitrilotriacetic acid 313 1992 313 1992 3-9 Disodium cyanodithioimidocarbonate 313 <td>NAMEC313ADD9-2o-Anisidine hydrochloride31319902-7alpha-NaphthylamineC31319900-6Benzeneamine, N-hydroxy-N-nitroso, ammonium saltX19900-6Cupferron31319900-1Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate)C3130-6Eupferron31319900-1Carbamodithioic acid, dibutyl, sodium salt (Sodium dibutyldithiocarbamate)C3135-8Dipropyl isocinchomeronate31319956-8ThiramC31319956-8ThiramC31319959-1Copper, bis(dimethylcarbamodithioato-S,S)-, (copper dimethyldithiocarbamate)C*0-4Zinc, bis(dimetylcarbomodithioato-S,S)-, (ziram)C*1-7Potassium N-methyldithiocarbamate31319952-8Sodium methyldithiocarbamate31319953-9Nitrilotriacetic acid31319953-9Nitrilotriacetic acid31319905-33,3'-Dimethyldiphenylmethane-4,4'-diisocyanate31319905-44,4'-Thiodianiline31319908-5Ethyl acetateC19918-91,3-DichloropropaneC19919-12-DipropylamineC19913-9Sodium cyanide (Na(CN))C*1-2Cupric acetateC19913-9Sodium cya</td> <td>NAMEC313ADDCAS9-2 o-Anisidine hydrochloride313 1990145-73-32-7 alpha-NaphthylamineC313 1990148-78-80-6 Benzeneamine, N-hydroxy-N-nitroso, armonium saltX19900-6 Cupferron3131990148-79-80-6 Cupferron3131990148-79-80-7 Carbamodithioic acid, dibutyl, sodium saltC148-82-3(Sodium dibutyldithiocarbamate)31319950-8 Dipropyl isocinchomeronate31319950-4 Zinc, bis(dimethylcarbamodithioato-S,S)- (copper dimethyldithiocarbamate)*0-4 Zinc, bis(dimethylcarbamodithioato-S,S)- (copper dimethyldithiocarbamate)*0-4 Zinc, bis(dimethyldithiocarbamate)3130-4 Zinc, bis(dimethyldithiocarbamate)3130-4 Zinc, bis(dimethyldithiocarbamate)3130-5 2.33.3'Dimethyldithiocarbamate3132.4 Metham sodium3132.5 Ethyl acrylate3133-9 Dickiur cyanodithioimidocarbonate3133-9 Nitrilotriacetic acid3133-9 Sodium cyanodithioimidocarbonate3133-14, 4'-Thiodianilline3133-5 Ethyl acrylate3133-6 Ethyl acetateC19913-9 Sodium cyanide (Na(CN))C2-17 DipropylamineC3-19 Sodium cyanide (Na(CN))C2-17 DipropylamineC3-19 Sodium cyanide (Na(CN))19913-19 Sodium cyanide (Na(CN))C3-19 Sodium cyanide (Na(CN))-<td>NAMEC313ADD9-2 b-Anisidine hydrochloride31319902-7 Japha-NaphthylamineC31319900-6 Benzeneamine, N-hydroxy-N-nitroso, ammonium saltX19900-6 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o-Anisidine hydrochloride 313 1990145-73-32-7 alpha-NaphthylamineC 313 1990148-78-80-6 Benzeneamine, N-hydroxy-N-nitroso, armonium saltX19900-6 Cupferron3131990148-79-80-6 Cupferron3131990148-79-80-7 Carbamodithioic acid, dibutyl, sodium saltC148-82-3(Sodium dibutyldithiocarbamate)31319950-8 Dipropyl isocinchomeronate31319950-4 Zinc, bis(dimethylcarbamodithioato-S,S)- (copper dimethyldithiocarbamate)*0-4 Zinc, bis(dimethylcarbamodithioato-S,S)- (copper dimethyldithiocarbamate)*0-4 Zinc, bis(dimethyldithiocarbamate)3130-4 Zinc, bis(dimethyldithiocarbamate)3130-4 Zinc, bis(dimethyldithiocarbamate)3130-5 2.33.3'Dimethyldithiocarbamate3132.4 Metham sodium3132.5 Ethyl acrylate3133-9 Dickiur cyanodithioimidocarbonate3133-9 Nitrilotriacetic acid3133-9 Sodium cyanodithioimidocarbonate3133-14, 4'-Thiodianilline3133-5 Ethyl acrylate3133-6 Ethyl acetateC19913-9 Sodium cyanide (Na(CN))C2-17 DipropylamineC3-19 Sodium cyanide (Na(CN))C2-17 DipropylamineC3-19 Sodium cyanide (Na(CN))19913-19 Sodium cyanide (Na(CN))C3-19 Sodium cyanide (Na(CN))- <td>NAMEC313ADD9-2 b-Anisidine hydrochloride31319902-7 Japha-NaphthylamineC31319900-6 Benzeneamine, N-hydroxy-N-nitroso, ammonium saltX19900-6 Cupferron3131990148-18-50-6 Cupferron31319900-1 Carbarnodithioic acid, dibutyl, sodium saltC148-29-81-7 botassium controlithic acid, dibutyl, sodium saltC148-29-82-8 Dipropyl isocinchomeronate31319950-8 ThiramC3139-1 Copper, bis(dimethylcarbamodithioato-S,S)-, (ziram)C*1-7 Potassium N-methyldithiocarbamate31319952-8 Sodium methyldithiocarbamate31319952-8 Sodium methyldithiocarbamate31319952-9 Disolum cyanodithioimidocarbonate31319952-14,4'-Thiodianiline31319952-14,4'-Thiodianiline31319903-3 Nitrilotriacetic acid31319903-5Bityl acylate31319903-6Ethyl acylate31319903-6Ethyl acylate31319903-7C*19913-8Schur acylate31319903-14,4'-Thiodianiline31319903-14,4'-Thiodianiline31319903-14,4'-Thiodianiline31319903-14,4'-Thiodianiline31319903-14,4'-Thiodianiline31319903-14,4'-Thiodianiline313</td> <td>NAMEC313ADD2-2 o-Anisidine hydrochloride31319902-7 Jalpha-NaphthylamineC31319902-7 Jalpha-NaphthylamineC31319900-6 Benzeneamine, N-hydroxy-N-nitroso, armonium sattX19900-6 Cupferron313148-73-8Codum diethyldithiocarbamate)0-6 Cupferron313148-79-8C/abendithioic acid, diethyl-, sodium salt0-6 Cupferron31319900-1 Carbamodithiocarbamate)31319950-2 Copper dimethyldithiocarbamate)31319950-3 ThiramC3139-1 Copper, bis(dimethylcarbamodithiocarbasmate)C1-7 Potassium N-methyldithiocarbamate3139-2 Budi acrylate3133-2 Disodium cyanodithioid carbonate3133-3 DichloropropaneC0-4 Cuppic scienceC2-5 Litry acrylate3133-1 Siposi19953-2 Dicholamiline3133-3 DichloropropaneC3-4 Sodium cyanide (Na(CN))C1-7 DipopylamineC3-8 Sodium cyanide (Na(CN))C1-2 Oupric acetateC1-3 Carbamodithioic acid, dimethyl,C1-3 Carbamodithioic acid, dimethyl,C1-4 Copper dimineC1-3 Carbamodithioic acid, dimethyl,C1-7 Potassium N-methyldithocarbamate3131-7 Potassium CyanadeC2-8 Sodium methyldithocarbamate3132-9 Solium cyanideC<td< 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dimethyl,C1-3 Carbamodithioic acid, dimethyl,C1-4 Copper dimineC1-3 Carbamodithioic acid, dimethyl,C1-7 Potassium N-methyldithocarbamate3131-7 Potassium CyanadeC2-8 Sodium methyldithocarbamate3132-9 Solium cyanideC <td< td=""><td>NAME C 313 ADD 2-2 o-Anisidine hydrochloride 313 1990 145-73-3 Endothall C 313 AI 2-7 Jalpha-Naphthylarnine C 313 1990 145-73-3 Endothall C Statum C 313 AI 148-73-3 Endothall C Statum Statum C 313 1990 145-73-3 Endothall C Statum Statum<!--</td--></td></td<>	NAME C 313 ADD 2-2 o-Anisidine hydrochloride 313 1990 145-73-3 Endothall C 313 AI 2-7 Jalpha-Naphthylarnine C 313 1990 145-73-3 Endothall C Statum C 313 AI 148-73-3 Endothall C Statum Statum C 313 1990 145-73-3 Endothall C Statum Statum </td

 313: EPCRA 313 Reportable Chemical, X: EPCRA 313 Synonym,* : member of a chemical category – should not be reported as an individual chemical.
 C: CERCLA Chemical – If a chemical is noted as ONLY a CERCLA chemical, then a STATE ONLY Form R as well as a Form S must be completed and submitted to DEP with your toxics use report.

NOTE: Some EPA Chemicals may have been delisted from the EPCRA 313 list, but the chemical MAY STILL be listed as a CERCLA chemical.

T1 M	T. ' TI T	7.1 1 [*] A	D / 11	C1 · 1	T C	C 1 1 1	17 0001
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			Sorte	ed Nume	erically by CAS				
CAS	NAME	<u>C</u> 3	313 AI	DD	CAS	NAME	C· 3	13 ADI	D
460-35-5	3-Chloro-1,1,1-trifluoropropane		313	1995	533-74-4	Tetrahydro-3,5-dimethyl-2H-1,3,5-thiadiazine-2-	1	X	19
460-35-5	HCFC-253fb		X	1995		thione			
463-58-1	Carbon oxide sulfide (COS)	С	X	1990	534-52-1	4,6-Dinitro-o-cresol	C	313	19
463-58-1	Carbonyl sulfide	С	313	1990	534-52-1	4,6-Dinitro-o-cresol and salts	С		19
465-73-6	Isodrin	С	313	1992	534-52-1	Dinitrocresol	C	X	19
492-80-8	Auramine	С	X	1990	540-59-0	1,2-Dichloroethylene		313	19
492-80-8	C.I. Solvent Yellow 34	С	313	1990	540-73-8	Hydrazine, 1,2-dimethyl-	С		
494-03-1	Chlomaphazine	C	-		540-84-1	2,2,4-Trimethylpentane	C		
496-72-0	Diaminotoluene	C	1		540-84-1	2,2,4-Trimethylpentane	С		
504-24-5	4-Aminopyridine	С		1992	540-88-5	tert-Butyl acetate	C	•	19
504-24-5	Pyridine, 4-amino-	С	1	1992	541-09-3	Uranyi acetate	С		19
504-60-9	1,3-Pentadiene	С	· · ·	1992	541-41-3	Ethyl chloroformate		313	19
505-60-2	Ethane, 1,1'-thiobis[2-chloro-		X	1995	541-53-7	2,4-Dithiobiuret	С	313	19
505-60-2	Mustard gas		X	1995	541-53-7	Dithiobiuret	С	X	19
506-61-6	Potassium silver cyanide	С	*	1992	541-73-1	1,3-Dichlorobenzene	С	313	19
506-64-9	Silver cyanide	C	*	1992	542-62-1	Barium cyanide	C	*	19
506-68-3	Cyanogen bromide	С	*	1	542-75-6	1,3-Dichloropropene	С	X	19
506-77-4	Cyanogen chloride	c	*	11	542-75-6	1,3-Dichloropropylene	С	313	19
506-77-4	Cyanogen chloride ((CN)Cl)	С	*	1992	542-76-7	3-Chloropropionitrile	C ·	313	19
506-87-6	Ammonium carbonate	С	1	1992	542-76-7	Propionitrile, 3-chloro-	С	Х	19
506-96-7	Acetyl bromide	С	1	1992	542-88-1	Bis(chloromethyl) ether	C .	313	19
507-55-1	1,3-Dichloro-1,1,2,2,3-pentafluoropropane		313	1995	542-88-1	Chloromethyl ether	С	X	19
507-55-1	HCFC-225cb		X	1995	542-88-1	Dichloromethyl ether	С	Х	
509-14-8	Methane, tetranitro-	С	-	1992	542-88-1	Methane, oxybis[chloro-	С	X	19
509-14-8	Tetranitromethane	c		1992	543-90-8	Cadmium acetate	С	*	
510-15-6	Benzeneacetic acid, 4-chloro-alpha(4-	C	x		544-18-3	Cobaltous formate	С	*	Γ
	chlorophenyl)alphahydroxy-, ethyl ester				544-92-3	Copper cyanide	С	*	19
510-15-6	Chlorobenzilate	С	313	ŀ	554-13-2	Lithium carbonate		313	19
513-49-5	sec-Butylamine	·C	1.	1992	554-84-7	m-Nitrophenol	С		19
528-29-0	o-Dinitrobenzene	C	313	1990	556-61-6	Isothiocyanatomethane		X	19
532-27-4	2-Chloroacetophenone	С	313	1990	556-61-6	Methyl isothiocyanate		313	19
533-74-4	Dazomet		313		557-19-7	Nickel cyanide	С	*	\square

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The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2001	
Sorted Newsonia 11rs has CAS	

CAS	NAME	C 31	13 AD	D	crical.	CAS	NAME	C 31	13 AD	D
764-41-0	1,4-Dichloro-2-butene	С	313	1992	-	1116-54-7	N-Nitrosodiethanolamine	C		19
764-41-0	2-Butene, 1,4-dichloro-	С	X	1992		1120-71-4	1,3-Propane sultone	С	X	19
765-34-4	Glycidylaldehyde	С		1992		1120-71-4	Propane sultone	С	313	3 19
812-04-4	1,1-Dichloro-1,2,2-trifluoroethane		313			1129-41-5	Carbamic acid, methyl- 3-methylphenyl ester	С		T
812-04-4	HCFC-123b		Х				(metolcarb)			
815-82-7	Cupric tartrate	С	*			1134-23-2	Cycloate		313	3 19
822-06-0	Hexamethylene-1,6-diisocyanate	С	313*	1995		1163-19-5	Decabromodiphenyl oxide		313	3 19
823-40-5	Diaminotoluene	С				1185-57-5	Ferric ammonium citrate	C		19
834-12-8	Ametryn		313	1995		1194-65-6	Dichlobenil	С		19
834-12-8	N-Ethyl-N'-(1-methylethyl)-6-(methylthio)-1,3,5,-		Х	1995		1300-71-6	Xylenol	C		19
	triazine-2,4-diamine					1303-28-2	Arsenic pentoxide	C ·	*	
842-07-9	C.I. Solvent Yellow 14		313	1990		1303-32-8	Arsenic disulfide	С	*	
872-50-4	N-Methyl-2-pyrrolidone		313	1995		1303-33-9	Arsenic trisulfide	С	*	' 19
924-16-3	N-Nitrosodi-n-butylamine	С	313	1990		1309-64-4	Antimony trioxide	С	*	' 19
924-42-5	N-Methylolacrylamide		313	1995		1310-58-3	Potassium hydroxide	C		19
930-55-2	N-Nitrosopyrrolidine	C				1310-73-2	Sodium hydroxide	С		19
933-75-5	2,3,6-Trichlorophenol	С	*			1313-27-5	Molybdenum trioxide		313	3 19
933-78-8	2,3,5-Trichlorophenol	С	• *	1992		1314-20-1	Thorium dioxide		313	3 19
957-51-7	Diphenamid	313				1314-32-5	Thallic oxide	С	*	19
959-98-8	alpha - Endosulfan	С		1992		1314-62-1	Vanadium pentoxide	С		19
961-11-5	Phosphoric acid, 2-chloro-1-(2,3,5-trichlorophenyl)	X	1990		1314-80-3	Sulfur phosphide	С		19
	ethenyl dimethyl ester				Γ	1314-84-7	Zinc phosphide	С	*	19
961-11-5	Tetrachlorvinphos		313	1990		1314-84-7	Zinc phosphide (conc. <= 10%)	С	*	19
989-38-8	C.I. Basic Red 1	r	313	1990		1314-84-7	Zinc phosphide (conc. > 10%)	С		19
1024-57-3	Heptachlor epoxide	С				1314-87-0	Lead sulfide	С	*	í 19
1031-07-8	Endosulfan sulfate	С		1992		1319-72-8	2,4,5-T amines	С		19
1066-30-4	Chromic acetate	С	*	1992		1319-77-3	Cresol (mixed isomers)	С	313	19
1066-33-7	Ammonium bicarbonate	С		1992	·	1320-18-9	2,4-D Esters	С	X	19
1072-35-1	Lead stearate	C	*	1992		1320-18-9	2,4-D propylene glycol butyl ether ester	C	313	\$ 19
1111-78-0	Ammonium carbamate	С		1992		1321-12-6	Nitrotoluene	С		19
1114-71-2	Butylethylcarbamothioic acid S-propyl ester		Х	1995		1327-52-2	Arsenic acid	С	*	19
1114-71-2	Pebulate		313			1327-53-3	Arsenic trioxide	С	*	19

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The Massachusetts T	oxics Use Reduction Act, Reportable Chemical List fo	r Calendar Year 2001
	Sorted Numerically by CAS	•

CAS -	NAME	C 3	13 AD	D	CA	S	NAME	C 3	13 AD	D
1929-82-4	1 Nitrapyrin		313	1995	2	2303-16-4	Diallate	С	31:	3
1937-37-7	C.I. Direct Black 38		313	1990	2	2303-17-5	Triallate		31:	3
1982-69-0	3,6-Dichloro-2-methoxybenzoic acid, sodium salt		X	1995	2	312-35-8	Propargite	C	31:	3 1
1982-69-0) Sodium dicamba		313	1995	2	2524-03-0	Dimethyl chlorothiophosphate		31:	3 19
2008-41-5	Carbamothioic acid, bis(2-methylpropyl)-, S-	С			2	2524-03-0	Dimethyl phosphorochloridothioate		X	1
	ethyl ester (butylate)		_		2	2545-59-7	2,4,5-T esters	С		1
2008-46-0	2,4,5-T amines	С	_	1992	2	2556-36-7	1,4-Cyclohexane diisocyanate		313*	1
2025-85-2	2 Selenium dioxide	C		1992	2	2602-46-2	C.I. Direct Blue 6	•	31:	3 1
2032-65-7	Mercaptodimethur	С	Х		2	2631-37-0	Phenol, 3-methyl-5-(1-methylethyl)-, methyl	C		
2032-65-7	/ Methiocarb	С	313	1992		.1	carbamate (promecarb)			
2095-58-1	Borane, trifluoro-		Х	1995	. 2	2655-15-4	2,3,5-Trimethylphenyl methylcarbamate	_	313	3 1
2095-58-1	Boron trifluoride		313	1995	2	2699-79-8	Sulfuryl fluoride		313	3 1
2125-68-3	Phosphorous trichloride	С		1993	2	2699-79-8	Vikane			ব
2125-68-3	Phosphorus trichloride	С	1		2	2702-72-9	2,4-D sodium salt		31:	3 1
2139-59-4	Potassium bromate		313	1995	2	2763-96-4	5-(Aminomethyl)-3-isoxazolol	С		1
2146-10-8	Sodium chromate	С	*		2	2763-96-4	Muscimol	С	1	1
2148-87-8	Hydrogen sulfide	С	313		2	2764-72-9	Diquat	C	1.	1
2151-06-8	Strontium chromate	C.	*		2	2832-40-8	C.I. Disperse Yellow 3		31:	3 1
2151-16-3	Ammonium bichromate	С	*		2	2837-89-0	2-Chloro-1,1,1,2-tetrafluoroethane		313	3
2155-70-6	Tributyltin methacrylate	-	313	1995	2	2837-89-0	HCFC-124		>	
2164-07-0	7-Oxabicyclo(2.2.1)heptane-2,3-dicarboxylic acid	I, · · ·	Х	1995	2	921-88-2	Chlorpyrifos	С		1
	dipotassium salt				2	944-67-4	Ferric ammonium oxalate	С		1
2164-07-0	Dipotassium endothall		313	1995	2	971-38-2	2,4-D chlorocrotyl ester	С		-
2164-17-2	Pluometuron		313	1990	2	971-38-2	2,4-D Esters	C	X	1
2164-17-2	Urea, N,N-dimethyl-N'-[3-(trifluoromethyl)phenyl]		X		3	012-65-5	Ammonium citrate, dibasic	С	-	1
2212-67-1	1H-Azepine-1 carbothioic acid, hexahydro-S-ethy	/!	Х	1995	3	3118-97-6	C.I. Solvent Orange 7		31:	3 1
	ester				3	164-29-2	Ammonium tartrate	С	-	1
2212-67-1	Molinate		313	1995	3	165-93-3	4-Chloro-o-toluidine, hydrochloride	c	1	1
2234-13-1	Octachloronaphthalene		313	1990	3	173-72-6	1,5-Naphthalene diisocyanate	L	313*	11
2300-66-5	Dimethylamine dicamba		313	1995	3	251-23-8	Cupric nitrate	С	-	* 1
2303-16-4	Carbamothioic acid, bis(1-methylethyl)-S-(2,3-	C	X		3	288-58-2	O,O-Diethyl S-methyl dithiophosphate	. c	1	1
	dichloro-2-propenyl)ester				3	383-96-8	Temephos		31:	3

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The Massachusetts	Toxics U	se Reduction	Act, Reportable	Chemical List for Calendar Year 2001

			Sorte	ed Nume	rically by CAS				
CAS	NAME	C 3	513 AI	DD	CAS	NAME	C	313 AD	D
7440-66-	6 Zinc (fume or dust)	C	313	3 1990	7696-12-0	2,2-Dimethyl-3-(2-methyl-1-		X	119
7446-08-4	4 Selenium dioxide	C	1	*		propenyl)cyclopropanecarboxylic acid (1,3,	4,5,6,7-		
7446-14-2	2 Lead sulfate	C		* 1992		hexahydro-1,3-dioxo-2H-isoindol-2-yl)meth	yl ester		
7446-18-	6 Thallium(I) sulfate	С	,	* 1992	7696-12-0	Tetramethrin		31:	3 19
7446-18-	6 Thallous sulfate	С	3	* 1992	7697-37-2	Nitric acid	С	31:	3 19
7446-27-	7 Lead phosphate	C.	7	1992	7697-37-2	Nitric acid (conc 80% or greater)	С	X	19
7447-39-4	4 Cupric chloride	С		1992	7699-45-8	Zinc bromide	C		* 19
7488-56-4	4 Selenium sulfide	С	,	1992	7705-08-0	Ferric chloride	С		19
7550-45-0	0 Titanium chloride (TiCl4) (T-4)-	С	X	1990	7718-54-9	Nickel chloride	С		* 19
7550-45-0	0 Titanium tetrachloride	С	313	3 1990	7719-12-2	Phosphorous trichloride	C		19
7558-79-4	4 Sodium phosphate, dibasic	С		1992	7720-78-7	Ferrous sulfate	C.	•	* 19
7601-54-9	9 Sodium phosphate, tribasic	С		1993	7722-64-7	Potassium permanganate	С		* 19
7631-89-2	2 Sodium arsenate	С	, ,	1993	7723-14-0	Phosphorus	C		19
7631-90-	5 Sodium bisulfite	c	1	1993	7723-14-0	Phosphorus (yellow or white)	C	313	3 19
7632-00-0	Sodium nitrite	С	313	1993	7726-95-6	Bromine		313	3 19
7645-25-2	2 Lead arsenate	С	*	1993	7733-02-0	Zinc sulfate	С		19
7646-85-	Zinc chloride	C	*	1993	7738-94-5	Chromic acid	С		* 19
7647-01-0	Hydrochloric acid	c		1990	7758-01-2	Potassium bromate		313	3 19
7647-01-0	Hydrogen chloride (anhydrous)	C	X	1990	7758-29-4	Sodium phosphate, tribasic	С		19
7647-01-0	Hydrogen chloride (gas only)	С	X	1990	7758-94-3	Ferrous chloride	С		19
7647-18-9	Antimony pentachloride	c	*	1993	7758-95-4	Lead chloride	С		* 19
7664-38-2	2 Phosphoric acid	С	313	1990	7758-98-7	Cupric sulfate	С		* 19
7664-39-3	B Hydrofluoric acid	С	x		7761-88-8	Silver nitrate	С		* 19
7664-39-3	Hydrofluoric acid (conc. 50% or greater)	С	X		7773-06-0	Ammonium sulfamate	С		19
7664-39-3	B Hydrogen fluoride	С	313	1990	7778-39-4	Arsenic acid	С		* 19
7664-39-3	B Hydrogen fluoride (anhydrous)	c	X	1990	7778-44-1	Calcium arsenate	C	4	* 19
7664-41-7	Ammonia	c	313	1990	7778-50-9	Potassium bichromate	С	4	* 19
7664-93-9	Sulfuric acid (aerosol)	C	313	1995	7778-54-3	Calcium hypochlorite	C		19
7681-49-4	Sodium fluoride	С		1993	7779-86-4	Zinc hydrosulfite	С	+	* 19
7681-52-9	Sodium hypochlorite	Ċ.		1993	7779-88-6	Zinc nitrate	С	4	* 19
				1	7782-41-4	Fluorine	С	313	3 19
					7782-49-2	Selenium	С	313	3 19

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The Massachusetts	Foxics Use Reduction Act, Reportable Chemical List for Calendar Year 2001
	Control Nisses and a line loss CAC

CAS	NAME	C _3	<u>13 AD</u>	D	C	AS	NAME	C 313	AD	D
10102-43-9	Nitric oxide	С	2 199			12002-03-8	Paris green	C		19
10102-43-9	Nitrogen oxide (NO)	С	\.	1993		12039-52-0	Selenious acid, dithallium(1+) salt	С	*	19
10102-44-0	Nitrogen dioxide	С		1993		12054-48-7	Nickel hydroxide	С	*	19
10102-45-1	Thallium(I) nitrate	С	*	1993		12122-67-7	Carbamodithioic acid, 1,2-ethanediylbis-, zinc co	mplex	X	19
10102-48-4	Lead arsenate	С	*	1993		12122-67-7	Zineb		313	3 19
10108-64-2	Cadmium chloride	С	*			12125-01-8	Ammonium fluoride	C.	[· ·	19
10124-50-2	Potassium arsenite	С	. *	1993		12125-02-9	Ammonium chloride	C		19
10124-56-8	Sodium phosphate, tribasic	С		1993		12135-76-1	Ammonium sulfide	С		19
10140-65-5	Sodium phosphate, dibasic	С		1993		12427-38-2	Carbamodithioic acid, 1,2-ethanediylbis-, manga	nese	Х	1
10192-30-0	Ammonium bisulfite	С		1993			complex		L	
10196-04-0	Ammonium sulfite	С		1	Ĺ	12427-38-2	Maneb		313	<u>; 19</u>
10222-01-2	2,2-Dibromo-3-nitrilopropionamide		313	1995	· .	12672-29-6	Aroclor 1248	С	Ľ.	
10294-34-5	Borane, trichloro-		Х	1995		12674-11-2	Aroclor 1016	С	L	19
10294-34-5	Boron trichloride		313	1995		12771-08-3	Sulfur monochloride	С	· .	19
10347-54-3	1,4-Bis(methylisocyanate)cyclohexane		313*	1995		13194-48-4	Ethoprop		313	19
10361-89-4	Sodium phosphate, tribasic	С		1993		13194-48-4	Ethoprophos		X	19
10380-29-7	Cupric sulfate, ammoniated	С	*			13194-48-4	Phosphorodithioic acid O-ethyl S,S-dipropyl este	r	X	19
10415-75-5	Mercurous nitrate	C	*			13356-08-6	Fenbutatin oxide		313	; 19
. 10421-48-4	Ferric nitrate	С	*			13356-08-6	Hexakis(2-methyl-2-phenylpropyl)distannoxane	-	Х	19
10453-86-8	5-(Phenylmethyl)-3-furanyl)methyl 2,2-dimethyl-3-	-(2-	X	1995		13463-39-3	Nickel carbonyl	С	*	19
	methyl-1-propenyl)cyclopropanecarboxylate	`	ŀ			13463-40-6	Iron carbonyl (Fe(CO)5), (TB-5-11)-		X	· ·
10453-86-8	Resmethrin		313	1995		13463-40-6	Iron, pentacarbonyl-		313	19
10544-72-6	Nitrogen dioxide	С				13474-88-9	1,1-Dichloro-1,2,2,3,3-pentafluoropropane		313	19
10588-01-9	Sodium bichromate	С	*			13474-88-9	HCFC-225cc		X	(
10605-21-7	Carbamic acid, 1H-benzimidazol-2-yl,methyl	C				13560-99-1	2,4,5-T salts	С		19
	ester (carbendazim)	<u> </u>	ļ			13597-99-4	Beryllium nitrate	C	*	19
11096-82-5	Aroclor 1260	C				13684-56-5	Desmedipham	313	1995	<i>i</i>
11097-69-1	Aroclor 1254	<u>с</u>		1993		13746-89-9	Zirconium nitrate	С	•*	19
11104-28-2	Aroclor 1221	C		1993		13765-19-0	Calcium chromate	С	*	19
11115-74-5	Chromic acid	C	*			13814-96-5	Lead fluoborate	С.,	*	
11141-16-5	Aroclor 1232	С		1993		13826-83-0	Ammonium fluoborate	С		19
12002-03-8	Cupric acetoarsenite			1993		13952-84-6	sec-Butylamine			19

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The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2001 Sorted Mr

			Solle	a nume	encal	IY DY CAS				
CAS	NAME	C 3	813 AE	D		CAS	NAME	C 31	3 AD	D
23950-58-5	Benzamide, 3,5-dichloro-N-(1,1-dimethyl-2-	С	X	1993		26952-23-8	Dichloropropene	С		
	propynyl					27176-87-0	Dodecylbenzenesulfonic acid	С		1
23950-58-5	Pronamide	C	313			27314-13-2	4-Chloro-5-(methylamino)-2-[3-		X	19
25154-54-5	Dinitrobenzene (mixed isomers)	С					(trifluoromethyl)phenyl]-3(2H)-pyridazinone			
25154-55-6	Nitrophenol (mixed isomers)	С		1993		27314-13-2	Norflurazon		313	3 19
25155-30-0	Sodium dodecylbenzenesulfonate	С		1993		27323-41-7	Triethanolamine dodecylbenzene sulfonate	С		19
25167-82-2	Trichlorophenol	С	*			27774-13-6	Vanadyl sulfate	С		19
25168-15-4	2,4,5-T esters	С				28057-48-9	d-trans-Allethrin		313	3 19
25168-26-7	2,4-D Esters	С			_	28057-48-9	d-trans-Chrysanthemic acid of d-allethrone		X	19
25311-71-1	2-((Ethoxyl((1-methylethyl)amino]phosphinothioy	l]oxy)	X	1995		28249-77-6	Carbamic acid, diethylthio-, S-(p-chlorobenzyl)		X	19
-	benzoic acid 1-methylethyl ester					28249-77-6	Thiobencarb		313	319
25311-71-1	Isofenphos		313	1995	. [28300-74-5	Antimony potassium tartrate	С		19
25321-14-6	Dinitrotoluene (mixed isomers)	С	313			28407-37-6	C.I. Direct Blue 218	· · ·	313	3 19
25321-22-6	Dichlorobenzene	С	. X - 2°			29082-74-4	Octachlorostyrene		313	+
25321-22-6	Dichlorobenzene (mixed isomers)	С	313			29232-93-7	O-(2-(Diethylamino)-6-methyl-4-pyrimidinyl)-O,C)-	X	19
25376-45-8	Diaminotoluene (mixed isomers)	C	313	1990			dimethyl phosphorothioate			
25376-45-8	Toluenediamine	C	X	1990		29232-93-7	Pirimiphos methyl		313	3 19
25550-58-7	Dinitrophenol	С		1993		30525-89-4	Paraformaldehyde	C		
26002-80-2	2,2-Dimethyl-3-(2-methyl-1-		X	1995		30558-43-1	Ethanimidothioci acid, 2-(dimethylamino-n-	С		
	propenyl)cyclopropanecarboxylic acid (3-			-			hydroxy-2-oxo-, methyl ester (A2213)			
	phenoxyphenyl)methyl ester		_			30560-19-1	Acephate		313	19
26002-80-2	Phenothrin			313		30560-19-1	Acetylphosphoramidothioic acid O,S-dimethyl es	ster	X	19
26264-06-2	Calcium dodecylbenzenesulfonate	С				31218-83-4	3-((Ethylamino)methoxyphosphinothioyl)oxy)-2-		X	19
26419-73-8	1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-	С				×	butenoic acid, 1-methylethyl ester			
	,O- [(methylamino)carbonyl]oxime (tripate)				L	31218-83-4	Propetamphos		313	5 19
26471-62-5	Benzene, 1,3-diisocyanatomethyl-	C	X	1990		32534-95-5	2,4,5-TP esters		С	Γ
26471-62-5	Toluene diisocyanate (unspecified isomer)	C	X			33089-61-1	Amitraz			3
26471-62-5	Toluenediisocyanate (mixed isomers)	C	313		Г	33213-65-9	beta - Endosulfan		С	1
26628-22-8	Sodium azide (Na(N3))	С	313		Г	34014-18-1	N-(5-(1,1-Dimethylethyl)-1,3,4-thiadiazol-2-yl)-N	,N'-	X	19
26638-19-7	Dichloropropane	С					dimethylurea			
26644-46-2	N,N'-(1,4-Piperazinediylbis(2,2,2-trichloroethylide	ne))	X	1995	· · L	34014-18-1	Tebuthiuron		313	, 19
	bisformamide					34077-87-7	Dichlorotrifluoroethane		313	<u>ا</u>
26644-46-21	Triforine		1 313							

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The Massachusetts Toxics Use Reduction Act, Reportable Chemical List for Calendar Year 2001 Sorted Numerically by CAS

CAS	NAME	C 🗄	313 AD	D
55290-64-7	Dimethipin			313
55406-53-6	3-lodo-2-propynyl butylcarbamate		313	1995
55488-87-4	Ferric ammonium oxalate	C	·····	· ·
56189-09-4	Lead stearate	С	*	·
57213-69-1	Triclopyr triethylammonium salt		313	1995
59669-26-0	Thiodicarb			313
60168-88-9	.alpha(2-Chlorophenyl)alpha4-chlorophenyl){ pyrimidinemethanol	5-	X	1995
60168-88-9	Fenarimol			313
60207-90-1	1-(2-(2,4-Dichlorophenyl)-4-propyl-1,3-dioxolan-2- methyl-1H-1,2,4,-triazole	-yl)-	X	1995
60207-90-1	Propiconazole		313	1995
61792-07-2	2,4,5-T esters	С		(
62476-59-9	5-(2-Chloro-4-(trifluoromethyl)phenoxy)-2-nitrober acid, sodium salt	izoic	Х	1995
62476-59-9	Acifluorfen, sodium salt		313	1995
63938-10-3	Chlorotetrafluoroethane		313	
64902-72-3	2-Chloro-N-(((4-methoxy-6-methyl-1,3,5-triazin-2- yl)aminolcarbonyl)benzenesulfonamide		X	1995
64902-72-3	Chlorsulfuron		313	1995
64969-34-2	3 3'-Dichlorobenzidine sulfate		313	1995
66441-23-4	2-(4-((6-Chloro-2-		X	1995
00441 20 4	benzoxazolylen)oxy)phenoxy)propanoic acid, ethy ester	/1		
66441-23-4	Fenoxaprop ethyl	,	313	1995
67485-29-4	Hydramethylnon		313	1995
67485-29-4	Tetrahydro-5,5-dimethyl-2(1H)-pyrimidinone(3-(4- (trifluoromethyl)phenyl)-1-(2-(4-		Х	1995
	(trifluoromethyl)phenyl)ethenyl)-2- propenylidene)hydrazone			
68085-85-8	3-(2-Chloro-3,3,3-trifluoro-1-propenyl)-2,2-		X	1995
•	Dimethylcyclopropanecarboxylic acid cyano(3-			
	phenoxyphenyl) methyl ester			

CAS	NAME C 31	3 ADI)
68085-85-8	3 Cyhalothrin	313	19
68359-37-5	3-(2,2-Dichloroethenyl)-2,2-	X	
	dimethylcyclopropanecarboxylic acid, cyano(4-fluoro-		
•	3-phenoxyphenyl)methyl ester		
68359-37-5	5 Cyfluthrin		3
69409-94-5	Fluvalinate		3
69409-94-5	N-(2-Chloro-4-(trifluoromethyl)phenyl)-DL-valine(+)-	X	19
·	cyano(3-phenoxyphenyl)methyl ester		
69806-50-4	2-(4-(5-(Trifluoromethyl)-2-pyridinyl]oxy]-	X	19
•	phenoxy)propanoic acid, butyl ester		
69806-50-4	Fluazifop butyl	313	19
71751-41-2	Abamectin		3
71751-41-2	2 Avermectin B1	X	19
72178-02-0	5-(2-Chloro-4-(trifluoromethyl)phenoxy)-N-	X	19
	methylsulfonyl)-2-nitrobenzamide		
72178-02-0	Fomesafen	313	19
72490-01-8	2 (2-(4-Phenoxy-phenoxy)-ethyl)carbamic acid ethyl	X	19
	ester		
72490-01-8	Fenoxycarb	313	19
74051-80-2	2-(1-(Ethoxyimino) butyl)-5-(2-(ethylthio)propyl)-3-	Х	19
	hydroxyl-2-cyclohexen-1-one		
74051-80-2	Sethoxydim	313	19
75790-84-0	4-Methyldiphenylmethane-3,4-diisocyanate	313*	19
75790-87-3	2,4'-Diisocyanatodiphenyl sulfide	313*	19
76578-14-8	2-(4-((6-Chloro-2-quinoxalinyl)oxy]phenoxy) propanoic	X	19
	acid ethyl ester		
76578-14-8	Quizalofop-ethyl	313	19
77501-63-4	5-(2-Chloro-4-(trifluoromethyl)phenoxy)-2-nitro-2-	Х	19
	ethoxy-1-methyl-2-oxoethyl ester		
77501-63-4	Lactofen	313	19
82657-04-3	Bifenthrin	313	19
88671-89-0	alphaButylalpha(4-chlorophenyl)-1H-1,2,4-	X	19
· · · ·	triazole-1-propanenitrile		

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	The Massachu	isens TOXICS	OSC K	eduction Act,	repor	ladic CI	lenne	ar Lis	l IOI Ca	nenuar	1 ear 2001	
Sorted Numerically by CAS												
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NAM	E			С	313	ADD						

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