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Thursday, December 4, 2003

Part IV

Environmental Protection Agency

40 CFR Parts 302 and 355 Reportable Quantity Adjustments for Carbamates and Carbamate-Related Hazardous Waste Streams; Reportable Quantity Adjustment for Inorganic Chemical Manufacturing Processes Waste (K178); Proposed Rule

ENVIRONMENTAL PROTECTION AGENCY

40 CFR Parts 302 and 355

[SW H-FRL-7594-4]

RIN 2050-AE12

Reportable Quantity Adjustments for Carbamates and Carbamate-Related Hazardous Waste Streams; Reportable Quantity Adjustment for Inorganic Chemical Manufacturing Processes Waste (K178)

AGENCY: Environmental Protection Agency (EPA).

ACTION: Proposed rule.

SUMMARY: The U.S. Environmental Protection Agency (EPA or "the Agency") is proposing reportable quantity (RQ) adjustments for 28 individual carbamates and five carbamate-related hazardous waste streams listed as hazardous wastes under the Resource Conservation and Recovery Act, and as hazardous substances with one-pound statutory RQs under the Comprehensive Environmental Response, Compensation, and Liability Act (CERCLA). In addition, EPA is proposing to adjust the one-pound statutory RQ of another hazardous waste stream, K178, which is unrelated to the carbamates addressed in this rule.

EPA thoroughly evaluated the intrinsic properties of these substances to assess the possibility of harm from the release of each substance into the environment and to determine the appropriate levels that require release notification. The proposed RQ adjustments will relieve the regulated community and emergency response personnel from the burden of making and receiving reports of releases that are unlikely to pose a threat to public health or welfare or the environment.

DATES: To make sure we consider your comments on this proposed rule, they must be postmarked on or before February 2, 2004. Comments postmarked after this date will be marked "late" and may not be considered.

ADDRESSES: Comments submitted by regular U.S. Postal Service mail should be sent to: Docket Coordinator, Superfund Docket Office, Mail Code 5202T, U.S. Environmental Protection Agency Headquarters, Ariel Rios Building, 1200 Pennsylvania Avenue, NW., Washington, DC 20460. Comments may also be submitted electronically, in person, or by special delivery. To ensure proper receipt by EPA, it is imperative that you identify the appropriate docket control number in the subject line on the first page of your comment. These docket control numbers, as well as detailed instructions on how to submit your comments, are provided in the section entitled "How and to Whom Do I Submit Comments?" in the supplemental information portion of this preamble.

FOR FURTHER INFORMATION CONTACT: For general information, contact the RCRA, Superfund, and EPCRA Hotline at 800/ 424-9346 or TDD 800/553-7672 (hearing impaired). In the Washington, DC metropolitan area, call 703/412-9810 or TDD 703/412-3323 (hearing impaired). For information on specific aspects of the rule, contact Lynn Beasley of the Office of Emergency and Remedial Response (5204G), U.S. Environmental Protection Agency, Ariel Rios Building, 1200 Pennsylvania Avenue, NW., Washington, DC 20460. Ms. Beasley's e-mail address is beasley.lynn@epa.gov, and her telephone number is 703/603-9086.

SUPPLEMENTARY INFORMATION:

I. General Information

A. Potentially Regulated Entities

Type of entity	Examples of affected entities
Industry	Manufacturers, handlers, transporters, and other users of carbamates. These substances are often used as insecticides, fungicides, herbicides, accelerators in the vulcanization of rubber, or as chemical intermediates in the manufacture of drugs, pesticides, or resins. In addition, entities that may release K178 waste streams will also be affected.
State, Local, or Tribal Governments Federal Government	State Emergency Response Commissions, and Local Emergency Planning Committees. National Response Center, and any Federal agency that may release these carbamates and waste streams.

This table is not intended to be exhaustive, but rather provides a guide for readers regarding entities likely to be regulated by this action. This table lists the types of entities that EPA is now aware could potentially be regulated by this action. Other types of entities not listed in the table could also be regulated. To determine whether your facility, company, business, or organization is regulated by this action, vou should carefully examine the proposed changes to 40 CFR parts 302 and 355. If you have questions regarding the applicability of this action to a particular entity, consult the person listed in the preceding FOR FURTHER INFORMATION CONTACT section.

B. How Can I Get Copies of Support Documents for This Rule?

1. *Docket.* EPA has established an official public docket for the Carbamates and Carbamate-Related Hazardous

Waste Streams (Docket ID No. SFUND-2002-0010) and an official public docket for the Inorganic Chemical Manufacturing Processes Waste (K178) (Docket ID No. SFUND-2002-0011). The official public docket consists of the documents specifically referenced in this action, any public comments received, and other information related to this action. Although a part of the official docket, the public docket does not include Confidential Business Information (CBI) or other information whose disclosure is restricted by statute. The official public docket is the collection of materials that is available for public viewing at the Superfund Docket in the EPA Docket Center, (EPA/ DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC. The EPA Docket Center Public Reading Room is open from 8:30 a.m. to 4:30 p.m., Monday through Friday,

excluding legal holidays. The telephone number for the Public Reading Room is (202) 566–1744, and the telephone number for the Superfund Docket is (202) 566–0270. You may copy a maximum of 100 pages from any regulatory docket at no cost. Additional copies cost \$0.15 per page. The Docket Office will mail copies of materials to you if you are located outside the Washington, DC metropolitan area.

2. *Electronic Access.* You may access this **Federal Register** document electronically through the EPA Internet under the **Federal Register** listings at *http://www.epa.gov/fedrgstr/.*

An electronic version of the public docket is available through EPA's electronic public docket and comment system, EPA Dockets. You may use EPA Dockets at http://www.epa.gov/edocket/ to submit or view public comments, access the index listing of the contents of the official public docket, and to access those documents in the public docket that are available electronically. Once in the system, select "search," then key in the appropriate docket identification number.

Certain types of information will not be placed in the EPA Dockets. Information claimed as CBI and other information whose disclosure is restricted by statute, which is not included in the official public docket, will not be available for public viewing in EPA's electronic public docket. EPA's policy is that copyrighted material will not be placed in EPA's electronic public docket but will be available only in printed, paper form in the official public docket. To the extent feasible, publicly available docket materials will be made available in EPA's electronic public docket. When a document is selected from the index list in EPA Dockets, the system will identify whether the document is available for viewing in EPA's electronic public docket. Although not all docket materials may be available electronically, you may still access any of the publicly available docket materials through the docket facility identified in Unit I.B.1. EPA intends to work towards providing electronic access to all of the publicly available docket materials through EPA's electronic public docket.

For public commenters, it is important to note that EPA's policy is that public comments, whether submitted electronically or in paper, will be made available for public viewing in EPA's electronic public docket as EPA receives them and without change, unless the comment contains copyrighted material, CBI, or other information whose disclosure is restricted by statute. When EPA identifies a comment containing copyrighted material, EPA will provide a reference to that material in the version of the comment that is placed in EPA's electronic public docket. The entire printed comment, including the copyrighted material, will be available in the public docket.

Public comments submitted on computer disks that are mailed or delivered to the docket will be transferred to EPA's electronic public docket. Public comments that are mailed or delivered to the Docket will be scanned and placed in EPA's electronic public docket. Where practical, physical objects will be photographed, and the photograph will be placed in EPA's electronic public docket along with a brief description written by the docket staff.

For additional information about EPA's electronic public docket, visit

EPA Dockets online or see 67 FR 38102, May 31, 2002.

C. How and to Whom Do I Submit Comments?

You may submit comments electronically, by mail, by facsimile, or through hand delivery/courier. To ensure proper receipt by EPA, identify the appropriate docket identification number in the subject line on the first page of your comment. Please ensure that your comments are submitted within the specified comment period. Comments received after the close of the comment period will be marked "late." EPA is not required to consider these late comments. However, late comments may be considered if time permits.

1. Electronically. If you submit an electronic comment as prescribed below, EPA recommends that you include your name, mailing address, and an e-mail address or other contact information in the body of your comment. Also include this contact information on the outside of any disk or CD ROM you submit, and in any cover letter accompanying the disk or CD ROM. This ensures that you can be identified as the submitter of the comment and allows EPA to contact you in case EPA cannot read your comment due to technical difficulties or needs further information on the substance of your comment. EPA's policy is that EPA will not edit your comment, and any identifying or contact information provided in the body of a comment will be included as part of the comment that is placed in the official public docket, and made available in EPA's electronic public docket. If EPA cannot read your comment due to technical difficulties and cannot contact you for clarification, EPA may not be able to consider your comment.

i. EPA Dockets. Your use of EPA's electronic public docket to submit comments to EPA electronically is EPA's preferred method for receiving comments. Go directly to EPA Dockets at http://www.epa.gov/edocket, and follow the online instructions for submitting comments. Once in the system, select "search," and then key in Docket ID No. SFUND–2002–0010 for the Carbamates and Carbamate-Related Hazardous Waste Streams or Docket ID No. SFUND-2002-0011 for the Inorganic Chemical Manufacturing Processes Waste (K178). The system is an "anonymous access" system, which means EPA will not know your identity, e-mail address, or other contact information unless you provide it in the body of your comment.

ii. *E-mail.* Comments may be sent by electronic mail (e-mail) to

superfund.docket@epa.gov, Attention Docket ID No. SFUND-2002-0010 for Carbamates and Carbamate-Related Hazardous Waste Streams or Docket ID No. SFUND-2002-0011 for Inorganic **Chemical Manufacturing Processes** Waste (K178). In contrast to EPA's electronic public docket, EPA's e-mail system is not an "anonymous access" system. If you send an e-mail comment directly to the Docket without going through EPA's electronic public docket, EPA's e-mail system automatically captures your e-mail address. E-mail addresses that are automatically captured by EPA's e-mail system are included as part of the comment that is placed in the official public docket, and made available in EPA's electronic public docket.

iii. *Disk or CD ROM.* You may submit comments on a disk or CD ROM that you mail to the mailing address identified in Unit I.C.2. These electronic submissions will be accepted in WordPerfect or ASCII file format. Avoid the use of special characters and any form of encryption.

2. *By Mail.* Send an original and two copies of your comments to: Superfund Docket, Environmental Protection Agency, Mailcode: [5202T], 1200 Pennsylvania Ave., NW., Washington, DC, 20460, Attention Docket ID No. SFUND–2002–0010 for Carbamates and Carbamate-Related Hazardous Waste Streams or Docket ID No. SFUND–2002– 0011 for Inorganic Chemical Manufacturing Processes Waste (K178).

3. By Hand Delivery or Courier. Deliver your comments to: Public Reading Room, Room B102, EPA West Building, 1301 Constitution Avenue, NW., Washington, DC, Attention Docket ID No. SFUND–2002–0010 for Carbamates and Carbamate-Related Hazardous Waste Streams or Docket ID No. SFUND–2002–0011 for Inorganic Chemical Manufacturing Processes Waste (K178). Such deliveries are only accepted during the Docket's normal hours of operation as identified in Unit I.B.1.

4. *By Facsimile*. Fax your comments to: (202) 566–0272, Attention Docket ID. No. SFUND–2002–0010 for Carbamates and Carbamate-Related Hazardous Waste Streams or Docket ID No. SFUND–2002–0011 for Inorganic Chemical Manufacturing Processes Waste (K178).

D. What Should I Consider as I Prepare My Comments for EPA?

You may find the following suggestions helpful for preparing your comments:

1. Explain your views as clearly as possible.

2. Describe any assumptions that you used.

3. Provide any technical information and/or data you used that support your views.

4. If you estimate potential burden or costs, explain how you arrived at your estimate.

5. Provide specific examples to illustrate your concerns.

6. Offer alternatives.

7. Make sure to submit your comments by the comment period deadline identified.

8. To ensure proper receipt by EPA, identify the appropriate docket identification number in the subject line on the first page of your response. It would also be helpful if you provided the name, date, and **Federal Register** citation related to your comments.

II. Outline of Today's Preamble

A. Overview

- 1. Statutory Authority
- Does this proposed rule apply to me?
 What types of releases are exempt from
- these reporting requirements?
- B. Background
- C. Summary of Today's Action
- 1. What is the scope of today's rule?
- 2. What methodology is EPA using to adjust the RQs of the individual carbamates?
- 3. What RQs are proposed for the individual carbamates?
- 4. How is EPA adjusting the RQs for the carbamate-related waste streams?
- 5. What RQs are proposed for these carbamate-related waste streams?
- 6. What conforming changes are being made to Table 302.4 and its Appendix A?
- 7. What changes are being made to 40 CFR part 355?
- 8. What RQ is proposed for the K178 waste stream?
- D. Statutory and Regulatory Reviews
 - 1. Executive Order 12866: Regulatory Planning and Review
 - 2. Paperwork Reduction Act
 - 3. Regulatory Flexibility Act
 - 4. Unfunded Mandates Reform Act
 - 5. Executive Order 13132: Federalism
 - 6. Executive Order 13175: Consultation and Coordination with Indian Tribal Governments
 - 7. Executive Order 13045: Protection of Children from Environmental Risks and Safety Risks
 - 8. Executive Order 13211: Actions that Significantly Affect Energy Supply, Distribution or Use
 - 9. National Technology Transfer and Advancement Act of 1995

III. Preamble for Reportable Quantity Adjustments for Carbamates and Carbamate-Related Hazardous Waste Streams; Reportable Quantity Adjustment for Inorganic Chemical Manufacturing Processes Waste (K178)

A. Overview

1. Statutory Authority

The Comprehensive Environmental Response, Compensation, and Liability Act of 1980 (CERCLA), 42 U.S.C. 9601 et seq., as amended by the Superfund Amendments and Reauthorization Act of 1986, gives the Federal government broad authority to respond to releases or threats of releases of hazardous substances from vessels and facilities. The term "hazardous substance" is defined in section 101(14) of CERCLA by referencing various Federal environmental statutes. For example, the term includes "any hazardous waste having the characteristics identified under or listed pursuant to section 3001 of the Solid Waste Disposal Act * * *,' also known as the Resource Conservation and Recovery Act (RCRA).

Section 102(b) of CERCLA establishes reportable quantities (RQs) of one pound ("statutory RQs") for releases of most CERCLA hazardous substances. Under section 102(a) of CERCLA, the Administrator of EPA has the authority to adjust these RQs by regulation ("adjusted RQs").

Under CERCLA section 103(a), the person in charge of a vessel or facility from which a CERCLA hazardous substance has been released in a quantity that equals or exceeds its RQ must immediately notify the National Response Center (NRC) of the release. A release is reportable if an RQ or more is released within a 24-hour period (see 40 CFR 302.6). This reporting requirement serves as a trigger for informing the government of a release so that Federal personnel can evaluate the need for a Federal removal or remedial action and undertake any necessary action in a timely fashion.

In addition to the reporting requirements under CERCLA section 103, section 304 of the Emergency Planning and Community Right-to-Know Act of 1986 (EPCRA), 42 U.S.C. 11001 et seq., requires owners or operators of certain facilities to report releases of extremely hazardous substances (EHSs) and CERCLA hazardous substances to State and local authorities (see 40 CFR 355.40). After the release of a hazardous substance in a quantity equal to or greater than its RQ, facility owners or operators must immediately notify the community emergency coordinator for each local

emergency planning committee for any area likely to be affected by the release, and the State emergency response commission of any State likely to be affected by the release.

2. Does This Proposed Rule Apply to Me?

The person in charge of a vessel or facility from which a CERCLA hazardous substance is released in a quantity that equals or exceeds its RQ must notify appropriate authorities who can evaluate whether a government response is needed. Therefore, this proposed rule may affect the following entities: (1) Persons in charge of vessels or facilities that may release CERCLA hazardous substances (as identified in this proposal) and owners or operators of facilities that may release EHSs or CERCLA hazardous substances (as identified in this proposal) into the environment; and (2) entities that plan for or respond to such releases.

3. What Types of Releases Are Exempt From the Reporting Requirements?

In determining whether you must report the release of a carbamate that equals or exceeds its RQ, it should be noted that section 103(e) of CERCLA exempts from the notification provisions of CERCLA section 103(a): * * * the application of a pesticide product registered under the Federal Insecticide, Fungicide, and Rodenticide Act or * * * the handling and storage of such a pesticide product by an agricultural producer." The legislative history of CERCLA suggests that Congress intended this exemption to apply to the application of a pesticide generally in accordance with the pesticide's purpose.

In addition, if a release of a CERCLA hazardous substance meets the criteria under CERCLA section 103(e) for an exemption from reporting to the NRC, the same release is also exempt from reporting to State and local authorities under EPCRA section 304. In the context of today's proposed rule, EPA believes that the CERCLA section 103(e) reporting exemption provides a potential source of reporting relief under both CERCLA and EPCRA for certain releases of carbamate pesticides.

As EPA previously noted in an April 4, 1985 final rule (50 FR 13464), we do not consider the spill of a pesticide to be an application of the pesticide, nor do we consider a pesticide spill to be in accordance with the pesticide's purpose. Consequently, spills of a carbamate pesticide that equal or exceed an RQ must be reported to the NRC under CERCLA section 103 and to the appropriate State and local authorities under EPCRA section 304.

B. Background

In today's notice of proposed rulemaking (NPRM), EPA is proposing to adjust the statutory one-pound RQs for 28 individual carbamates and five carbamate-related waste streams. Today's rulemaking includes proposed RQ adjustments not only for individual carbamates, but also for thiocarbamates, dithiocarbamates, carbamoyl oximes, and several other individual substances that are closely related to carbamate production and/or waste generation. For purposes of simplicity, however, the preamble to today's proposed rule refers to all 28 individual substances for which RQ adjustments are being proposed as "carbamates," and to the five waste streams as "carbamaterelated" waste streams. In addition, EPA is proposing to adjust the one-pound statutory RQ of another hazardous waste stream, K178, which is unrelated to the carbamates addressed in this rule (see Section III.C.8 of today's preamble for information regarding K178). A summary of the developments leading up to today's proposed rulemaking as it relates to the carbamate-related substances is provided below.

On November 8, 1984, Congress amended RCRA by enacting the Hazardous and Solid Waste Amendments of 1984 (HSWA), 42 U.S.C. 6901 et seq. In one provision of HSWA—a newly added RCRA section 3001(e)(2)—Congress directed EPA to determine whether several wastes, including wastes generated from the production of carbamates, should be listed as RCRA hazardous wastes. Carbamates are widely used as active ingredients in pesticides, herbicides, insecticides, and fungicides, and in the production of synthetic rubber. Before Congress enacted HSWA in 1984, EPA already had regulated several carbamate substances under RCRA, CERCLA, and other statutes.

Based on our evaluation of the carbamate production wastes, we published on March 1, 1994 (59 FR 9808), a proposal to list 80 carbamaterelated substances as RCRA hazardous wastes and as CERCLA hazardous substances. These 80 substances included: (1) 70 individual carbamates; (2) six carbamate-related waste streams; and (3) four categories of carbamate substances. Subsequently, on February 9, 1995 (60 FR 7824), we finalized the listing of 64 of these 80 substances as RCRA hazardous wastes and CERCLA hazardous substances, deferring action on 12 individual substances and the four categories of carbamate substances

included in the proposed rule. Thus, EPA listed a total of 58 individual carbamates and six carbamate-related hazardous waste streams as RCRA hazardous wastes and CERCLA hazardous substances in the February 9, 1995 final rule. We published corrections to minor errors in these listings in the **Federal Register** on April 17, 1995 (60 FR 19165) and May 12, 1995 (60 FR 25619). We also modified our interpretation of the rule as it affected listings for K156 and K157 hazardous wastes on August 14, 1995 (60 FR 41817).

On November 1, 1996, the Court of Appeals (D.C. Circuit) ruled that EPA failed to follow proper rulemaking procedures in making some of the carbamate listing determinations in the February 9, 1995 rule. Dithiocarbamate Task Force v. EPA, 98 F.3d 1394 (D.C.Cir. 1996). As a result, the court vacated the RCRA hazardous waste and CERCLA hazardous substance listings for 24 of the 58 individual carbamates and one of the six carbamate-related waste streams (K160) included in that rule. In addition, the court vacated three other carbamate-related waste streams (K156, K157, and K158) only to the extent that they applied to the chemical 3-iodo-2-propynyl n-butylcarbamate. Under the court decision, the vacated carbamate listings are to be treated as though they had never been in effect.

To clarify the legal status of the vacated listings for the regulated community and the public, EPA, in a June 17, 1997 final rule (62 FR 32974), amended the lists of RCRA hazardous wastes and CERCLA hazardous substances (in 40 CFR parts 261 and 302 respectively) to remove the entries for the 24 individual carbamates and one carbamate-related waste stream (K160) that were vacated by the court, as well as revised the entries for K156, K157, and K158 to indicate that they do not apply to 3-iodo-2-propynyl nbutylcarbamate.

It is important to note, however, that the court's ruling did not change the February 9, 1995 listing of the 34 remaining individual carbamates as RCRA hazardous wastes; those listings remain in effect. Independent of the February 9, 1995 rule, EPA already has added six of these 34 individual carbamates to the CERCLA list of hazardous substances in Table 302.4 of 40 CFR 302.4, and developed adjusted RQs for these substances because of their listing under the Clean Air Act or Clean Water Act.¹ The six substances and their Chemical Abstracts Service Registry Numbers (CASRNs) are: carbaryl (CASRN 63–25–2); carbofuran (CASRN 1563–66–2); mercaptodimethur (CASRN 2032–65–7); mexacarbate (CASRN 315–18–4); propoxur (CASRN 114–26–1); and triethylamine (CASRN 121–44–8).² Thus, we are not proposing any RQ adjustments for these six substances today.

Upon the effective date of the February 9, 1995 final rule, the 28 remaining individual carbamates and the five carbamate-related hazardous waste streams became hazardous substances under CERCLA section 101(14)(C) and received one-pound statutory RQs. We are proposing today to adjust the statutory one-pound RQs for these 28 substances and five waste streams based on criteria that relate to the possibility of harm from the release of each hazardous substance into the environment. EPA will revise the List of Hazardous Substances and Reportable Quantities (Table 302.4 of 40 CFR 302.4) to reflect these proposed changes and other, conforming proposed changes, if they are finalized. However, until such time as we finalize the adjusted RQs proposed in today's rule, the statutory RO of one pound remains in effect for these substances.

Finally, eleven of the individual substances with proposed RQ adjustments in today's rule are also EPCRA section 302 EHSs. For the names of these 11 substances, see the proposed revisions to appendices A and B of 40 CFR part 355, included at the end of today's rule. In an August 30, 1989 rule (54 FR 35988), we proposed to adjust the RQs for all the EPCRA EHSs.³ We finalized adjustments to the RQs for all the EHSs, except the 11 included in today's rule (61 FR 20473, May 7, 1996). We are reproposing adjusted RQs for these 11 substances today, for reporting under both CERCLA and EPCRA.

C. Summary of Today's Action

1. What Is the Scope of Today's Rule?

In today's rule, we are proposing to adjust the one-pound statutory RQs for 28 individual carbamates (one of which is adjusted to a final RQ of one-pound) and five carbamate-related waste streams. In addition, EPA is proposing to adjust the one-pound statutory RQ of another hazardous waste stream, K178, which is unrelated to the carbamates

 $^{^1}$ We adjusted the RQs for five of these six substances in an April 4, 1985 final rule (50 FR 13456), and adjusted the RQ for the other substance,

propoxur, in a June 12, 1995 final rule (60 FR 30926).

² Although not a carbamate, triethylamine is used during the production of carbamates.

³We used the data from this August 30, 1989 proposed rulemaking, as well as more recent data, to support the RQ adjustments proposed for these 11 substances in today's rule.

addressed in this rule (see section III.C.8 of today's preamble for information regarding K178). We based these adjustments on specific scientific and technical criteria that relate to the possibility of harm from the release of a CERCLA hazardous substance in certain amounts. RQs are based, in part, on a determination of possible or potential harm, but they are not a determination that releases of a particular amount of a hazardous substance necessarily will harm the public health, welfare, or the environment. The quantity released is just one factor that the Federal government considers when it assesses the need to respond to such a release. Other factors include, but are not limited to, the location of the release, its proximity to drinking water supplies or other valuable resources, and the likelihood of exposure or injury to nearby populations. The RQ adjustments that EPA is proposing today would enable us to focus our resources on those releases that are most likely to pose potential threats to public health, welfare, or the environment. These RQ adjustments also would relieve the regulated community and emergency response personnel from the burden of making and receiving reports of releases that are unlikely to pose such threats.

2. What Methodology Is EPA Using To Adjust the RQs of the Individual Carbamates?

EPA has wide discretion in adjusting the statutory RQs for hazardous substances under CERCLA. Administrative feasibility and practicality are important considerations. Our methodology for adjusting the RQ of an individual hazardous substance begins with an evaluation of its intrinsic physical, chemical, and toxicological properties. These intrinsic properties—called "primary criteria"—are aquatic toxicity, mammalian toxicity (oral, dermal, and inhalation), ignitability, reactivity, chronic toxicity, and potential carcinogenicity.4

Generally, for each intrinsic property, EPA ranks hazardous substances on a five-tier scale, associating a specific range of values on each scale with an RQ value of 1, 10, 100, 1,000, or 5,000 pounds. Each hazardous substance may receive several tentative RQ values based on the primary criteria. The lowest of the tentative RQs becomes the "primary criteria RQ" for that substance.

When we find sufficient data in the scientific literature on the chronic toxicity and/or potential carcinogenicity (two of the six primary criteria) of a substance, we generally evaluate and summarize these data in a chemicalspecific profile. Following an extensive review of available scientific literature on the 28 individual carbamates addressed in today's proposed rule, we found that chronic toxicity profiles are warranted for nine of these 28 carbamates, and that potential carcinogenicity profiles are warranted for six of the 28 carbamates. EPA has placed these 15 draft chemical-specific profiles in the docket for this proposed rulemaking.⁵ Proposed RQs for several of the substances included in today's rule are based, at least in part, on the conclusions drawn in these profiles.

We are soliciting comments on these drafts. We will consider data that you submit, including any additional toxicity or carcinogenicity data that may be available on these substances or the other carbamates included in today's proposed rule. If the data are applicable, we will incorporate them into the draft profiles prior to their completion.

After assigning the primary criteria RQs, we further evaluate the substances for their susceptibility to certain degradative processes. These natural degradative processes, which we use as ''secondary RQ adjustment criteria,'' are biodegradation, hydrolysis, and photolysis (BHP). If a hazardous substance, when released into the environment, degrades relatively rapidly to a less hazardous form by one or more of the BHP processes, we generally increase its RQ (as determined by the primary RQ adjustment criteria) by one level.⁶ Conversely, if a hazardous substance degrades to a more hazardous product after its release, we assign an RQ equal to the RQ for the more hazardous substance, which may be one or more levels lower than the RQ for the original substance.

Three carbamates—bendiocarb, benomyl, and thiophanate-methylhave BHP data that are a sufficient basis for adjusting the primary criteria RQs for these substances. Although several other carbamates (e.g., propham) have BHP data that suggest rapid degradation, the evidence for most of these substances is not conclusive. Therefore, no adjustment to the RQs for these other carbamates was proposed on the basis of BHP.⁷ EPA is requesting that commenters submit additional degradation data (e.g., data on BOD₅ values and on half lives), if available, on these 28 individual substances.8

EPA could not locate acceptable data on any of the primary or secondary criteria for three of the 28 individual carbamates in today's proposed rule (see Table 1). In the past, when adjusting the statutory RQs of such data-poor hazardous substances, we have used data from chemically similar, surrogate substances.9 Therefore, to adjust the statutory RQs of the three data-poor carbamates in today's rule, we conducted an analysis of other carbamates to identify potential surrogate substances (*i.e.*, carbamates with primary criteria data that are chemically similar, based primarily on structural analogy, to the data-poor substances).

Table 1 lists the chemically similar carbamates EPA used as proposed surrogates, and the proposed RQs we assigned to each data-poor substance based on its chemically similar

⁸One or more of the following criteria must be met for a hazardous substance to qualify for further RQ adjustment based on BHP: (1) Biodegradation: the substance must have a five-day biochemical oxygen demand (BOD5) that equals or exceeds 50 percent of the theoretical oxygen demand as calculated based on stoichiometric oxidation; and (2) Hydrolysis/Photolysis: the half-life of the substance in the environment must be five days or less. For further information on the methodology for applying BHP, see the Technical Background Document to Support Rulemaking Pursuant to CERCLA section 102, Volume 1, March 1985, available for inspection at the Superfund Docket in the EPA Docket Center, (EPA/DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC.

⁹ For further information on, and examples of, EPA's use of surrogate data to adjust RQs of hazardous substances, see section 2 of the Technical Background Document to Support Rulemaking Pursuant to CERCLA section 102, Volume 8, available for inspection at the Superfund Docket in the EPA Docket Center, (EPA/DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC.

⁴ For further information on assigning adjusted RQs to hazardous substances under the primary criteria, see the Technical Background Document to Support Rulemaking Pursuant to CERCLA Section 102, Volume 2, August 1986 (for chronic toxicity), Volume 3, July 1989 (for potential carcinogenicity), and Volume 1, March 1985 (for the four other primary criteria), available for inspection at the Superfund Docket in the EPA Docket Center, (EPA/DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC.

⁵ You may inspect the preliminary draft Reportable Quantity documents and potential carcinogenicity evaluations at the Superfund Docket in the EPA Docket Center, (EPA/DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC.

⁶ We do not raise an RQ level based on BHP if the primary criteria RQ is already at its highest possible level (100 pounds for potential carcinogens and 5,000 pounds for all other types of hazardous substances). The secondary adjustment criteria of BHP are not applied to radionuclides.

⁷ To review a summary of the BHP data on the 28 carbamates included in today's rule, see Exhibit 4–3 of the Technical Background Document to Support Rulemaking Pursuant to CERCLA section 102, Volume 8, available for inspection at the Superfund Docket in the EPA Docket Center, (EPA/ DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC.

surrogate.¹⁰ We are requesting primary and secondary criteria data on these three data-poor substances. We also are soliciting comments from readers on our a choice of surrogate substances used to c

adjust the RQs for these three carbamates.

TABLE 1.—PROPOSED	RQS FOR THE	DATA-POOR	CARBAMATES
TABLE I. TROPODED	It do I of The	BARN FOOR	0/11/0/11/120

Data-poor carbamate	Proposed surrogate	Proposed RQ (pounds)
Bendiocarb phenol	Bendiocarb	1000
Carbofuran phenol	Carbofuran	10
Manganese dimethyldithiocarbamate	Ziram	10

Please note that in Table 2, below, we are assigning different RQs for the datapoor carbamate/surrogate pair of Bendiocarb phenol (data-poor carbamate) and Bendiocarb (its proposed surrogate) in shown in Table 1, above. In Table 2, EPA is applying the secondary criteria of BHP to adjust the RQ for bendiocarb to 100 pounds. However, due to structural differences between the two substances, we believe that it would be inappropriate to apply the BHP data for bendiocarb to bendiocarb phenol; hence, EPA is proposing a 1000-pound RQ for bendiocarb phenol (see Tables 1 and 2).

3. What RQs Are Proposed for the Individual Carbamates?

Table 2 lists the chemical names, CASRNs, and proposed RQs for the 28

individual carbamates included in today's proposed rule. The proposed RQs for 27 of the 28 individual carbamates would be raised from their one-pound statutory levels, while one of the 28 individual carbamates— Dimetilan—would be adjusted to its proposed, final RQ of one pound.

TABLE 2.—PROPOSED RQS FOR 28 INDIVIDUAL CARBAMATE	ΞS
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Chemical name	CASRN	Proposed RQ (pounds)
A2213	30558-43-1	5000
Aldicarb sulfone	1646-88-4	100
Barban	101–27–9	10
Bendiocarb	22781-23-3	100
Bendiocarb phenol	22961-82-6	1000
Benomyl	17804-35-2	10
Carbendazim	10605-21-7	10
Carbofuran phenol	1563–38–8	10
Carbosulfan	55285-14-8	1000
m-Cumenyl methylcarbamate	64–00–6	10
Diethylene glycol, dicarbamate	5952-26-1	5000
Dimetilan	644–64–4	1
Formetanate hydrochloride	23422-53-9	100
Formparanate	17702–57–7	100
Isolan	119–38–0	100
Manganese dimethyldithiocarbamate	15339–36–3	10
Metolcarb	1129–41–5	1000
Oxamyl	23135–22–0	100
Physostigmine salicylate	57–64–7	100
Physostigmine	57–47–6	100
Promecarb	2631–37–0	1000
Propham	122-42-9	1000
Prosulfocarb	52888-80-9	5000
Thiodicarb	59669–26–0	100
Thiophanate-methyl	23564-05-8	10
Tirpate	26419–73–8	100
Triallate	2303–17–5	100
Ziram	137–30–4	10

4. How is EPA Adjusting the RQs for the Carbamate-Related Waste Streams?

In addition to the 28 individual carbamate hazardous substances, we also are proposing to adjust the RQs of the five carbamate-related RCRA hazardous waste streams (K156, K157, K158, K159, and K161). The standard methodology used to adjust the RQs for RCRA hazardous waste streams differs from the methodology applied to individual hazardous substances. Our procedure for assigning RQs to RCRA waste streams is based on an analysis of the hazardous constituents of the waste

¹⁰ These three data-poor carbamates also are included in the list of 28 individual carbamates that appears in Table 2. For further information on the three data-poor carbamates and the chemically-

similar, surrogate substances that EPA has identified, see section 3 of the Technical Background Document to Support Rulemaking Pursuant to CERCLA section 102, Volume 8,

available for inspection at the Superfund Docket in the EPA Docket Center, (EPA/DC) EPA West, Room B102, 1301 Constitution Ave., NW., Washington, DC.

streams. Specifically, EPA identifies the constituents of each RCRA hazardous waste stream in 40 CFR part 261, appendix VII. We determine the RQ for each constituent within the waste stream and establish the lowest RQ value of these constituents as the adjusted RQ for the waste stream. We also apply this same methodology to adjust the RQ for K178 (see section III.C.8 for more information).

5. What RQs Are Proposed for These Carbamate-Related Waste Streams?

In the February 9, 1995 final rule, the five carbamate-related waste streams were assigned the statutory one-pound RQ required by CERCLA section 102(b). In today's rule, using the standard methodology for adjusting RQs for RCRA waste streams, EPA is proposing a one-pound adjusted RO for waste stream K161 and 10-pound adjusted RQs for the remaining four carbamaterelated waste streams (K156, K157, K158, and K159) based on the constituent(s) with the lowest RO within each of the waste streams. Table 3 lists the constituents and constituent ROs of each of the five carbamaterelated hazardous waste streams.

TABLE 3.—CONSTITUENTS OF FIVE CARBAMATE-RELATED WASTE STREAMS

10

Carbamate Waste Stream Con- stituents	RQ (pounds)
K156	1
benomyl	1
carbaryl	10
carbendazim	1
carbofuran	1
carbosulfan	100
formaldehyde	10
methylene chloride	100
triethylamine	500
K157	1
carbon tetrachloride	1
formaldehyde	10
methyl chloride	10
methylene chloride	100
pyridine	100
triethylamine	500
K158	1
benomyl	1
carbendazim	1
carbofuran	1
carbosulfan	100
chloroform	1
methylene chloride	100
K159	1
benzene	1
butylate	10
EPTC	100
molinate	1
pebulate	10
vernolate	10
K161	
antimony	500
arsenic	

TABLE	3.—CONS	TITUENTS	OF	FIVE
CARE	BAMATE-REL	ATED	W	/ASTE
STRE	AMS-Cont	tinued		

Carbamate Waste Stream Con-	RQ
stituents	(pounds)
metam sodium	10
ziram	10

6. What Conforming Changes Are Being Made to Table 302.4 and Its Appendix A?

EPA is proposing to modify the entries in Table 302.4 for the carbamates added by the February 9, 1995 final rule to the list of CERCLA hazardous substances. Specifically, we are proposing in today's rule to change the entries for the chemical names of the carbamates in the "Hazardous Substance" column in Table 302.4 to reflect more accurately the chemical names for these substances as they appear in the RCRA tables of hazardous wastes at 40 CFR 261.33(e) and (f).

For example, the February 9, 1995 final rule generally lists two names for each individual carbamate in Table 302.4—a chemical name and, in parentheses, a synonym. Thus, the February 9, 1995 final rule added one entry for each carbamate to the CERCLA list of hazardous substances. The same final rule alphabetically lists these two names as separate entries in the RCRA tables of hazardous wastes in 40 CFR 261.33.

10 Because each of the 28 individual 100 carbamates included in today's rule has 10 at least two separate entries in the RCRA 10 tables of hazardous wastes, we are 000 proposing to make the CERCLA table of 100 hazardous substances consistent by 000 000 listing the two (or more) synonymous 10 names as separate entries in Table 10 302.4. Thus, amendatory instruction 3, 100 which immediately precedes Table 100 302.4 in today's proposed rule, accounts 000 for the addition of the chemical names 000 and synonyms as separate entries in 000 10 Table 302.4, and amendatory instruction 10 2 accounts for the removal of the 10 previously listed names for these 10 substances. We believe that proposing 000 these changes to Table 302.4 is a 10 positive step toward ensuring that 000 chemical lists under RCRA and CERCLA 10 are more consistent and that carbamate 10 synonyms are easier to find in the table. 100 000 In addition, we are proposing

10 conforming changes to entries in
100 appendix A to Table 302.4 for the 28
1 carbamates added to the list of CERCLA
1 hazardous substances by the February 9,
1 1995 final rule.

7. What Changes Are Being Made to 40 CFR Part 355?

Appendices A and B of 40 CFR part 355, which list EHSs and their threshold planning quantities (TPQs) under EPCRA, also list the RQs for EHSs. Eleven of the individual carbamates for which EPA is proposing adjusted RQs are EHSs, as well as CERCLA hazardous substances. EPA today is proposing to revise appendices A and B of 40 CFR part 355 to include these adjusted RQs. For the names of these 11 substances, see the proposed revisions to appendices A and B included at the end of today's proposed rule.

8. What RQ Is Proposed for the K178 Waste Stream? $^{\scriptscriptstyle 11}$

As noted in section III.C.4 of the preamble, the Agency's standard methodology for adjusting the RQs for RCRA waste streams is based on an analysis of the hazardous constituents of each waste, as identified in 40 CFR part 261, appendix VII. We determine an RQ for each constituent and establish the lowest RO value of these constituents as the adjusted RQ for the waste stream. When there are hazardous constituents identified for a waste stream that are not individual CERCLA hazardous substances, EPA develops an RQ for these constituents in order to assign an appropriate RQ to the waste stream (see 48 FR 23565, May 25, 1983). In other words, we derive the RQ for an RCRA waste stream based on the lowest RQ of all of the hazardous constituents identified for that waste in appendix VII of 40 CFR part 261, regardless of whether the constituents are CERCLA hazardous substances.

On September 14, 2000, EPA published a proposed rule to list three waste streams from inorganic chemical manufacturing processes as RCRA hazardous wastes in 40 CFR 261.32 and as CERCLA hazardous substances in 40 CFR 302.4 (65 FR 55684). In that rule, we proposed to adjust the one-pound statutory RQ for two of the three waste streams, K176 and K177. For the third waste stream, K178 (nonwastewaters from the production of titanium dioxide by the chloride-ilmenite process), EPA identified two hazardous constituents. The two hazardous constituents identified in the proposed rule were: thallium, which is a CERCLA hazardous substance with a 1,000-pound RQ; and manganese, which does not appear on the CERCLA hazardous substance list in 40 CFR 302.4 and, therefore, has not been assigned an RQ. Because EPA had

67922

¹¹HSWA also directed EPA to determine whether wastes from the Inorganic Chemical Industry should be listed as RCRA hazardous wastes.

not yet developed an RQ for manganese at that time, we did not propose to adjust the RQ for K178 in the September 14, 2000 proposed rule.

Numerous commenters to the proposed rule objected to using manganese as a basis for listing K178 wastes, citing potential adverse impacts to many industries. Although EPA continues to believe that manganese poses significant issues that ultimately should be resolved, the court-ordered schedule for the hazardous waste listings provided no flexibility to address those issues fully before finalizing the listings. For that reason, in the final rule, EPA deferred final action on adding manganese to appendix VII of 40 CFR part 261 as a basis for listing K178 wastes (66 FR 58258; November 20, 2001). The final hazardous waste listing for K178 is based solely on thallium. As a result, we are proposing an RQ of 1,000 pounds for the K178 waste stream, based on the constituent RQ for thallium, the sole hazardous constituent identified for the waste stream

As stated in the section of the preamble entitled "How and to Whom Do I Submit Comments?," it is important to identify docket control number SFUND–2002–0011 in the subject line on the first page of your correspondence if you are submitting comments on the proposed 1,000-pound RQ for K178.

D. Statutory and Regulatory Reviews

1. Executive Order 12866: Regulatory Planning and Review

Under Executive Order 12866 (58 FR 51735), the Agency must determine whether this regulatory action is "significant" and therefore subject to formal review by the Office of Management and Budget (OMB) and to the requirements of the Executive Order, which include assessing the costs and benefits anticipated as a result of the proposed regulatory action. The Order defines "significant regulatory action" as one that is likely to result in a rule that may: (1) Have an annual effect on the economy of \$100 million or more or adversely affect in a material way the economy, a sector of the economy, productivity, competition, jobs, the environment, public health or safety, or state, local, or tribal governments or communities; (2) create a serious inconsistency or otherwise interfere with an action taken or planned by another agency; (3) materially alter the budgetary impact of entitlements, grants, user fees, or loan programs or the rights and obligations of recipients thereof; or (4) raise novel legal or policy

issues arising out of legal mandates, the President's priorities, or the principles set forth in the Executive Order.

It has been determined that this proposed rule is not a "significant regulatory action" under the terms of Executive Order 12866 and is therefore not subject to OMB review. EPA performed an economic analysis, which shows that this proposed rule will result in an annual cost savings of approximately \$90,640 to the regulated community and to Federal, State, and local governments, and does not result in any of the other effects that define a significant regulatory action. In this proposed rule, EPA would raise the RQs for 27 of the 28 individual substances and five of the six waste streams (including K178) from their current statutory one-pound levels. The remaining individual carbamate substance and carbamate-related waste stream will remain subject to an RQ of one pound.

We have estimated that these adjustments from the statutory onepound RQs will reduce by approximately 176 the number of reportable releases for these hazardous substances each year (see the economic analysis mentioned above). The estimated \$90,640 cost savings reflects only those effects of the RQ adjustments that are readily quantifiable in dollars and are associated with the release notification requirements under CERCLA section 103 and EPCRA section 304, including the associated activities of recordkeeping and notification processing

A detailed presentation of EPA's methodology, data sources, and computations applied for estimating the number of affected entities (industrial facilities) and economic impacts attributable to today's proposal is provided in the "Economic Impact Analysis" to this proposal.

2. Paperwork Reduction Act

The information collection requirements in this proposed rule have been submitted for approval to the Office of Management and Budget (OMB) under the Paperwork Reduction Act, 44 U.S.C. 3501 et seq. Information Collection Request (ICR) documents have been prepared by EPA (ICR Nos. 1049.09 and 1395.04). A copy of these ICRs may be obtained from Susan Ambry by mail at Collection Strategies Division; U.S. Environmental Protection Agency (2822), 1200 Pennsylvania Avenue, NW., Washington, DC 20460, or by calling (202) 566-1676, and by email at *ambry.susan@epamail.epa.gov*. A copy also may be downloaded off the Internet at http://www.epa.gov/icr.

EPA proposes the following conditions for reporting and recordkeeping: mandatory reporting requirements (CERCLA section 103(a) and EPCRA section 304) serve as triggers for informing the government of a release so that Federal, State, or local personnel can evaluate the need for removal or remedial actions, and undertake any necessary action in a timely fashion.

We estimate that the public reporting burden for collecting information required under CERCLA section 103 averages 4.1 labor hours (i.e., combined managerial, technical, and clerical hours) per response; the reporting burden under EPCRA section 304 averages approximately 5 labor hours per response. This estimate includes the time required to: make a determination whether a release requires a report to the NRC, the State, and local agencies; make the call(s); maintain a log of any calls made to government organizations; and make a follow-up written notification (if required under EPCRA section 304). The average burden estimates of 4.1 and 5 hours are provided only for the purpose of calculating the labor costs associated with the entire release reporting and recordkeeping process under CERCLA and EPCRA. Thus, these burden estimates should not be misinterpreted as reflecting the amount of time an individual has before he or she must call the NRC. Rather, CERCLA and EPCRA require that persons in charge of vessels or facilities immediately notify the NRC, the State, and local agencies of releases that equal or exceed an RQ.

Because we are proposing to raise the RQs for all but two of the substances included in today's rule, we expect the net reporting and recordkeeping burden associated with reporting releases of these substances under CERCLA section 103 to decrease. As noted in the economic impact analysis supporting today's proposed rule (and in Section III.D.2 of this preamble), we estimate that the annual reporting and recordkeeping burdens associated with reports to the NRC will be reduced by approximately 720 hours, and to SERCs and LEPCs by 880 hours.

An Agency may not conduct or sponsor, and a person is not required to respond to a collection of information unless it displays a currently valid OMB control number. The OMB control numbers for EPA's regulations are listed in 40 CFR part 9 and 48 CFR chapter 15.

3. Regulatory Flexibility Act

The Regulatory Flexibility Act (RFA), as amended by the Small Business Regulatory Enforcement Fairness Act of 1996 (SBREFA), 5 U.S.C. 601 *et. seq.*, generally requires an agency to prepare a regulatory flexibility analysis of any rule subject to notice and comment rulemaking requirements under the Administrative Procedure Act or any other statute, unless the agency certifies that the rule will not have a significant economic impact on a substantial number of small entities. Small entities include small businesses, small organizations, and small governmental jurisdictions.

For purposes of assessing the impacts of today's rule on small entities, small entity is defined as: (1) A small business that has fewer than 1000 or 100 employees per firm depending upon the SIC code the firm primarily is classified; (2) a small governmental jurisdiction that is a government of a city, county, town, school district or special district with a population of less than 50,000; and (3) a small organization that is any not-for-profit enterprise which is independently owned and operated and is not dominant in its field.

After considering the economic impacts of today's proposed rule on small entities, I hereby certify that this proposal will not have a significant economic impact on a substantial number of small entities. In determining whether a rule has a significant economic impact on a substantial number of small entities, the impact of concern is any significant adverse economic impact on small entities, since the primary purpose of the regulatory flexibility analyses is to identify and address regulatory alternatives "which minimize any significant economic impact of the proposed rule on small entities" (5 U.S.C. 603 and 604). Thus, an agency may certify that a rule will not have a significant economic impact on a substantial number of small entities if the rule relieves regulatory burden, or otherwise has a positive economic effect on small entities subject to the rule. For more information regarding the economic impact of this proposed rule, please refer to the economic background document to this proposal.

We have therefore concluded that today's proposed rule will relieve regulatory burden for small entities. We continue to be interested in the potential impacts of the proposed rule on small entities and welcome comments on issues related to such impacts.

4. Unfunded Mandates Reform Act

Title II of the Unfunded Mandates Reform Act of 1995 (UMRA), Public Law 104–4, establishes requirements for Federal Agencies to assess the effects of

their regulatory actions on State, local, and tribal governments and the private sector. Under section 202 of the UMRA, EPA must prepare a written analysis, including a cost-benefit analysis, for proposed and final rules with "Federal mandates" that may result in expenditures to State, local, and tribal governments, in the aggregate, or to the private sector, of \$100 million or more in any one year. Before promulgating an EPA rule for which a written statement is needed, section 205 of the UMRA requires EPA to identify and consider a reasonable number of regulatory alternatives and adopt the least costly, most cost-effective, or least burdensome alternative that achieves the objectives of the rule. The provisions of section 205 do not apply when they are inconsistent with applicable law. Moreover, section 205 allows EPA to adopt an alternative other than the least costly, most cost-effective or least burdensome alternative if the Administrator publishes with the final rule an explanation why that alternative was not adopted.

Before EPA establishes any regulatory requirements that may significantly or uniquely affect small governments, including tribal governments, it must have developed under section 203 of the UMRA a small government agency plan. The plan must provide for notifying potentially affected small governments, enabling officials to have meaningful and timely input in the development of regulatory proposals, and informing, educating, and advising small governments on compliance with the regulatory requirements.

EPA has determined that this rule does not include a Federal mandate that may result in expenditures of \$100 million or more for State, local, or tribal governments, in the aggregate, or the private sector in any one year. This is because this proposed rule imposes no enforceable duty on any State, local, or tribal governments. EPA also has determined that this rule contains no regulatory requirements that might significantly or uniquely affect small governments. In addition, as discussed above, the private sector is not expected to incur costs exceeding \$100 million. Therefore, today's proposed rule is not subject to the requirements of sections 202 and 205 of UMRA.

5. Executive Order 13132: Federalism

Executive Order 13132, entitled "Federalism" (64 FR 43255, August 10, 1999), requires EPA to develop an accountable process to ensure "meaningful and timely input by State and local officials in the development of regulatory policies that have federalism implications." "Policies that have federalism implications" is defined in the Executive Order to include regulations that have "substantial direct effects on the States, on the relationship between the national government and the States, or on the distribution of power and responsibilities among the various levels of government."

This proposal does not have federalism implications. It will not have substantial direct effects on the States, on the relationship between the national government and the States, or on the distribution of power and responsibilities among the various levels of government, as specified in Executive Order 13132. This rule directly affects manufacturers, handlers, transporters, and other users of carbamates; in addition, entities that may release K178 waste streams will also be affected. There are no State and local government bodies that incur direct compliance costs by this rulemaking. State and local government implementation expenditures are expected to be less than \$500,000 in any one year. Thus, the requirements of section 6 of the Executive Order do not apply to this proposal.

In the spirit of Executive Order 13132, and consistent with EPA policy to promote communications between EPA and State and local governments, EPA specifically solicits comment on this proposed rule from State and local officials.

6. Executive Order 13175: Consultation and Coordination With Indian Tribal Governments

Executive Order 13175, entitled "Consultation and Coordination with Indian Tribal Governments" (65 FR 67249, November 9, 2000), requires EPA to develop an accountable process to ensure "meaningful and timely input by tribal officials in the development of regulatory policies that have tribal implications." This proposed rule does not have tribal implications, as specified in Executive Order 13175. Today's rule does not significantly or uniquely affect the communities of Indian tribal governments, nor would it impose substantial direct compliance costs on them. Thus, Executive Order 13175 does not apply to this rule.

7. Executive Order 13045: Protection of Children From Environmental Risks and Safety Risks

The Executive Order 13045, entitled "Protection of Children from Environmental Health Risks and Safety Risks (62 FR 19885, April 23, 1997) applies to any rule that EPA determines (1) is "economically significant" as

67924

defined under Executive Order 12866, and (2) the environmental health or safety risk addressed by the rule has a disproportionate effect on children. If the regulatory action meets both criteria, the Agency must evaluate the environmental health or safety effects of the planned rule on children; and explain why the planned regulation is preferable to other potentially effective and reasonably feasible alternatives considered by the Agency.

This proposal is not subject to the Executive Order because it is not economically significant as defined in Executive Order 12866, and because the Agency does not have reason to believe the environmental health or safety risks addressed by this proposed rule present a disproportionate risk to children.

8. Executive Order 13211: Actions That Significantly Affect Energy Supply, Distribution or Use

This proposed rule is not a "significant energy action" as defined in Executive Order 13211, "Actions Concerning Regulations That Significantly Affect Energy Supply, Distribution, or Use" (66 FR 28355 (May 22, 2001)) because it is not likely to have a significant adverse effect on the supply, distribution, or use of energy. This proposed rule reduces regulatory burden. It thus should not adversely affect energy supply, distribution, or use.

9. National Technology Transfer and Advancement Act of 1995

Section 12(d) of the National Technology Transfer and Advancement Act of 1995 ("NTTAA"), Public Law 104-113, section 12 (d) (15 U.S.C. 272 note) directs EPA to use voluntary consensus standards in its regulatory activities, unless to do so would be inconsistent with applicable law or otherwise impractical. Voluntary consensus standards are technical standards (e.g., materials specifications, test methods, sampling procedures, and business practices) that are developed or adopted by voluntary consensus standards bodies. The NTTAA directs us to provide Congress, through OMB, explanations when we decide not to use available and applicable voluntary consensus standards.

The proposed rule does not involve technical standards. Therefore, EPA is

not considering the use of any voluntary consensus standards.

List of Subjects

40 CFR Part 302

Environmental protection, Air pollution control, Chemicals, Hazardous substances, Hazardous wastes, Intergovernmental relations, Natural resources, Reporting and recordkeeping requirements, Superfund, Water pollution control, Water supply.

40 CFR Part 355

Air pollution control, Chemicals, Hazardous substances, Intergovernmental relations, Natural resources, Reporting and recordkeeping requirements, Superfund, Water pollution control, Water supply.

Dated: November 25, 2003.

Michael O. Leavitt,

Administrator.

For the reasons set out in the preamble, it is proposed to amend title 40, chapter I of the Code of Federal Regulations as follows:

PART 302—DESIGNATION, REPORTABLE QUANTITIES, AND NOTIFICATION

1. The authority citation for part 302 continues to read as follows:

Authority: 42 U.S.C. 9602, 9603, 9604; 33 U.S.C. 1321 and 1361.

2. Table 302.4 in § 302.4 is amended by removing the entries for "1,3-Benzodioxol-4-ol, 2,2-dimethyl-, (Bendiocarb phenol)", "1,3-Benzodioxol-4-ol, 2,2-dimethyl-, methyl carbamate (Bendiocarb)", "7-Benzofuranol, 2,3-dihydro-2,2-dimethyl-(Carbofuran phenol)", "Benzoic acid, 2hydroxy-, compd. with (3aS-cis)-1,2,3,3a,8,8a-hexahydro-1,3a,8trimethylpyrrolo[2,3b]indol-5-yl methylcarbamate ester (1:1) (Physostigmine salicylate)", "Carbamic acid, 1H-benzimidazol-2-yl, methyl ester (Carbendazim)", "Carbamic acid, [1-[(butylamino)carbonyl]-1Hbenzimidazol-2-yl, methyl ester (Benomyl)", "Carbamic acid, (3chlorophenyl)-, 4-chloro-2-butynyl ester (Barban)", "Carbamic acid, [(dibutylamino)thio]methyl-, 2,3dihydro-2,2-dimethyl-7benzofuranyl ester (Carbosulfan)", "Carbamic acid, dimethyl-,1[(dimethylamino)carbonyl]-5-methyl-1H-pyrazol-3-yl ester

(Dimetilan)", "Carbamic acid, dimethyl-, 3-methyl-1-(1methylethyl)-1H-pyrazol-5-yl ester (Isolan)", "Carbamic acid, methyl-, 3methylphenyl ester (Metolcarb)", "Carbamic acid, [1,2 phenylenebis(iminocarbonothiovl)]bis-. dimethyl ester (Thiophanate-methyl)", "Carbamic acid, phenyl-, 1-methylethyl ester (Propham)", "Carbamothioic acid, bis(1-methylethyl)-, S-(2,3,3-trichloro-2propenyl) ester (Triallate)", "Carbamothioic acid, dipropyl-, S-(phenylmethyl) ester (Prosulfocarb)" "1,3-Dithiolane-2-carboxaldehyde, 2,4dimethyl-, O-[(methylamino)carbonyl]oxime (Tirpate)", "Ethanimidothioci acid, 2-(dimethylamino-N-hydroxy-2-oxo-, methyl ester (A2213)" "Ethanimidothoic acid, 2-(dimethylamino)-N-[[(methylamino)carbonyl]oxy]-2-oxo-, methyl ester (Oxamyl)", "Ethanimidothioic acid, N,N'-[thiobis[(methylimino) carbonyloxy]]bis-, dimethyl ester (Thiodicarb)", "Ethanol, 2,2'oxybis-, dicarbamate (Diethylene glycol, dicarbamate)", "Manganese, bis(dimethylcarbamodithioato-S,S')-(Manganese dimethyldithiocarbamate)", "Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)carbonyl]oxy]phenyl]-, monohydrochloride (Formetanate hydrochloride)", "Methanimidamide, N,N-dimethyl-N'-[2-methyl-4-[[(methylamino)carbonyl]oxy]phenyl]-(Formparanate)", "Phenol, 3-(1methylethyl)-, methyl carbamate (m-Cumenyl methylcarbamate)", "Phenol, 3-methyl-5-(l-methylethyl)-, methyl carbamate (Promecarb)", "Propanal, 2methyl-2-(methylsulfonyl)-, O-[(methylamino)carbonyl] oxime (Aldicarb sulfone)", "Pyrrolo[2,3blindol-5-ol, 1,2,3,3a,8,8a-hexahydro-1,3a,8-trimethyl-, methylcarbamate (ester), (3aS-cis)-(Physostigmine)", "Zinc, bis(dimethylcarbamodithioato-S,S')-(Ziram)", "K156", "K157", "K158", "K159", "K161", and "K178" and adding the following new entries in alphabetical order to read as follows (applicable footnotes have been republished without change):

§ 302.4 Designation of hazardous substances.

* * * * *

TABLE 302.4.—LIST OF HAZARDOUS SUBSTANCES AND REPORTABLE QUANTITIES [Note: All Comments/Notes Are Located at the End of This Table]

Aldicarb sulfore 1646884 4 P203 100 (45.4 Barban 101279 4 U280 10 (45.4 Bardingarb phenol 22861826 4 U364 1000 (45.6 Bennory 1.3.Berzodowal-4.al, 2.2-dimethyl. 100 (45.4 1000 (45.6 J.3.Berzodowal-4.al, 2.2-dimethyl. 100 (45.4 1000 (45.6 1000 (45.6 J.3.Berzodowal-4.al, 2.2-dimethyl. 100 (45.4 1000 (45.6 1000 (45.6 J.3.Berzodowal-4.al, 2.2-dimethyl. 100 (45.4 1000 (45.6 1000 (45.6 J.Berzodowal-4.al, 2.2-dimethyl. 100 (45.4 1000 (45.4 1000 (45.6 J.3.Berzodowal-2.2-bindlo5-2.0 100 (45.4 1007 (45.4 100 (45.4 Carbanic acid, (1) (1) bunylamicy acid withyl.scher 1563388 100 (45.4 100 (45.4 Carbanic acid, (1) (1) bunylamicy acid withyl - benzindkz02-2.9/1, methyl ester 10808217 10277 10 (45.6 Carbanic acid, (10 bunylamicy) - Admityl - 2.2 dimethyl-1.7-benzoluramyl ester 10280 100 (45.4 11281 10278 10277 10 (45.6 Carbanic acid, (10 bunylamicy) - Admityl-2.2 dimethyl-1.1-byrzzol-3.9/1 ester 11380 4 11281 100 (45.4<	Hazardous substance	CASRN	Statutory code†	RCRA waste number	Final RQ pounds (Kg)
Barban 101279 4 U280 10 (454 Bendiocarb 22781233 4 U278 100 (454 Bendiocarb 2281826 4 U278 100 (454 Ja-Benzdovol-4-0; 2.2-dimethyl 22861826 4 U271 10 (454 Ja-Benzdovol-4-0; 2.2-dimethyl 100 (454 1000 (454 1000 (454 Ja-Benzdovol-4-0; 2.2-dimethyl 10 (454 1000 (454 1000 (454 Ja-Benzdovol-4-0; 2.2-dimethyl 10 (454 1000 (454 1000 (454 Carbanic acd, 11-Houriyamico-2-dy methyl ester 1780/4352 10 (454 Carbanic acd, 11-Houriyamico-2-dy methyl ester 1780/3352 10 (454 Carbanic acd, 12-shoroparenyl, -1-Hoeralidazo-2-yll, methyl ester 1780/3352 10 (454 Carbanic acd, 12-shoroparenyl, -1-choro-2-bruryl ester 1780/3352 10 (454 Carbanic acd, 11-Houriyamico-2-dinocarbanicyl Practa-3-yl ester 113280 10 (454 Carbanic acd, 11-Houriyamico-2-dinocarbanicyl Practa-3-yl ester 113281 10 (454 Carbanic acd, 12-phenyleestelylenyle ster 12364068 U271 10 (454 <		30558431	4 *	U394	5000 (2270)
Bendiocarb 22781233 4 U278 100 (45.4 Bendiocarb 22861826 4 U271 100 (45.4 1.3.Bencodxox1-4.0, 2.2-dimethyt- 22861826 4 U278 100 (45.4 1.3.Bencodxox1-4.0, 2.2-dimethyt- 22661826 4 U278 100 (45.4 1.3.Bencodxox1-2, 2.3-dimethyt- 1563388 4 U367 10 (45.4 Benzolc acid, 2-hydroxy- compd. with (3aS-cis)-1,2,3,3,8,8,a-haxahydro-1,3,a,B 57647 4 P188 100 (45.4 Carbanic acid, 1-horzinitazor-2-yi, methyl caster 17004352 4 U277 10 (45.4 Carbanic acid, 2-hydroxy- compd. with (3aS-cis)-1,2,3,3,a,8,8a-haxahydro-1,3a,B 57647 4 P188 100 (45.4 Carbanic acid, 2-hydroxy-, compd. with (3aS-cis)-1,2,3,3,a,8,8a-haxahydro-1,3a,B 101 (45.4 2271 10 (45.4 Carbanic acid, 1-horzin-baxiny, 1-hornethydethydro, 1-horzin-baxiny distr 10280 10 (45.4 10127 10 (45.4 Carbanic acid, 1-horzin-baxiny, 1-horzin-bydethydro, 2-dimethydro, 2-dimethy	Aldicarb sulfone	1646884	4 *	P203	100 (45.4)
Bendicator phenol 22661826 4 U364 1000 (454 1.3-Benzodioxol-4-ol, 2.2-dimethyl- 2261826 4 U278 100 (454 1.3-Benzodioxol-4-ol, 2.2-dimethyl- 2261826 4 U278 100 (454 7-Benzolizardi - 2-ydimethyl-, methyl carbamate 22781233 4 U278 100 (454 7-Benzolizardi - 2-ydimethyl-, ocmpd, with (335-cis)-1.2.3.38.88-hexahydrol-1.3.8- trimethylgymolo[2,3-bi]indol-5-yl methyl ester 10605217 4 U372 10 (454 Carbamic acid, (1-Hobrajamidazol-2-yl, methyl ester 17804352 4 U271 10 (454 Carbamic acid, (1-Hobrajamidazol-2-yl, methyl ester 10805217 4 U320 10 (454 Carbamic acid, (1-Hobrajamidazol-2-yl, methyl ester 10279 4 U280 10 (454 Carbamic acid, (1-Hobrajamicazol-5-yl-terryl) ester 113904 P F189 1000 (454 Carbamic acid, (1-gherphyl-nebi(mino)carbonyl)-1-methyl-terryl-1-methyldinyl-2-2-dimethyl-1-methyldinyl-2-2-dimethyl-1-methyldinyldinyl-2-2-dimethyl-1-methyldinyldinyl 10 (454 Carbamic acid, (1, 2-pheryl-nebi(mino)carbonyl)-1-methyl-1-methyldinyldinyl-2-2-dimethyl-1-methyldinyldinyl-2-2-dimethyl 10 (454 C	Barban	101279	4	U280	10 (4.54)
Bendicator phenol 22661826 4 U364 1000 (454 1.3-Benzodioxol-4-ol, 2.2-dimethyl- 2261826 4 U278 100 (454 1.3-Benzodioxol-4-ol, 2.2-dimethyl- 2261826 4 U278 100 (454 7-Benzolizardi - 2-ydimethyl-, methyl carbamate 22781233 4 U278 100 (454 7-Benzolizardi - 2-ydimethyl-, ocmpd, with (335-cis)-1.2.3.38.88-hexahydrol-1.3.8- trimethylgymolo[2,3-bi]indol-5-yl methyl ester 10605217 4 U372 10 (454 Carbamic acid, (1-Hobrajamidazol-2-yl, methyl ester 17804352 4 U271 10 (454 Carbamic acid, (1-Hobrajamidazol-2-yl, methyl ester 10805217 4 U320 10 (454 Carbamic acid, (1-Hobrajamidazol-2-yl, methyl ester 10279 4 U280 10 (454 Carbamic acid, (1-Hobrajamicazol-5-yl-terryl) ester 113904 P F189 1000 (454 Carbamic acid, (1-gherphyl-nebi(mino)carbonyl)-1-methyl-terryl-1-methyldinyl-2-2-dimethyl-1-methyldinyl-2-2-dimethyl-1-methyldinyldinyl-2-2-dimethyl-1-methyldinyldinyl 10 (454 Carbamic acid, (1, 2-pheryl-nebi(mino)carbonyl)-1-methyl-1-methyldinyldinyl-2-2-dimethyl-1-methyldinyldinyl-2-2-dimethyl 10 (454 C		00704000	^	, 11070	400 (45 4)
Benomyl 17804352 4 U271 10 (454 1.3-Benzodizool-4-ol, 2.2-dimethyl- 2261826 U284 U000 (454 1.3-Benzodizool-4-ol, 2.2-dimethyl- 1563388 U367 100 (454 7-Benzofuranol, 2.3-dihydro-2.2-dimethyl- 1563388 U367 10 (454 Carbanic acid, 2-hydroxy-, compd, with (335-cic)-1.2,3,3,8,8-hexahydro-1,3,8,8 57647 P188 100 (45.4 Carbanic acid, 1-Hobenzinidazol-2-yl, methy ester 10605217 4 U372 10 (45.4 Carbanic acid, (abmethyl-1-4)-benzinidazol-2-yl, methy ester 101274 4 U280 10 (45.4 Carbanic acid, (abmethyl-1-4)-enzinidazol-2-yl, Emthyl ester 17804352 4 U271 10 (45.4 Carbanic acid, dimethyl-1-3-dimuthyl-2-benzinidazol-2-yl, methyl ester 17804352 4 U271 10 (45.4 Carbanic acid, dimethyl-1-3-methylpenyl ester 110274 4 U280 10 (4.54 Carbanic acid, dimethyl-3-methylphenyl ester 113980 P192 100 (45.4 Carbanic acid, dimethyl-3-methylphenyl ester 2306175 4 U499 10 (4.54 Ca			-		()
1.3-Benzodioxol-4-ol, 2.2-dimethyl- 22961826 4 U364 1000 (454 1.3-Benzodioxol-4-ol, 2.2-dimethyl-, methyl carbamate 22781233 4 U278 100 (454 7-Benzoduranol, 2.3-dihydro-2.2-dimethyl- 1563388 4 U367 10 (454 Benzoic acid, 2-hydroxy-, compd, with (385-cis)-1.2.3.3a,8.a-hexahydro-1.3a,8-// imethyl-groupd)-H-benzoimateach 57647 P188 100 (45.4 Carbamic acid, [H-benzimidacol-2-yl, methyl ester 10805217 4 U372 10 (45.4 Carbamic acid, [Gibuhylminy)-hydro-2-butynyl ester 10805217 4 P188 100 (45.4 Carbamic acid, [Gibuhylminy)-hydro-2-butynyl ester 101276 4 P189 1000 (45.4 Carbamic acid, [Gibuhylminy)-hydro-2-butynyl ester 1123415 P190 1000 (45.4 Carbamic acid, methyl-, 3-methyl-1-1-methylethyl ester 1123415 P190 1000 (45.4 Carbamic acid, [henryh-, 1-methylethyl) seter 2230175 4 U393 1000 (45.4 Carbamic acid, methyl-, 3-methyl-10-11-methylethyl seter 22303175 4 U393 1000 (45.4 Carbamic acid, [henryh-, 1-methylethyl] seter 23388409 U397 100 (45.4 <td< td=""><td>I Contraction of the second second</td><td></td><td>-</td><td></td><td>()</td></td<>	I Contraction of the second		-		()
1,3-Benzodioxol-4-ol, 2,2-dimethyl-, methyl carbamate 22781233 4 U278 100 (45.4 7-Benzoluranol, 2,3-dinydro-2,2-dimethyl- 1563388 4 U367 10 (45.4 Benzolic acid, 2-hydroxy- compd, with (3aS-cis)-1,2,3,3a,8,8-hexahydro-1,3a,5 57647 4 P188 100 (45.4 Carbamic acid, 11-benzimidazol-2-yl, methyl ester 10605217 4 U377 10 (45.4 Carbamic acid, [1-[butylamino]carboryl]-11-benzimidazol-2-yl], methyl ester 17804522 4 U278 100 (45.4 Carbamic acid, (Inbutyl-11-benzimidazol-2-yl], methyl ester 17804522 4 U271 10 (45.4 Carbamic acid, (Inbutyl-11-benzimidazol-2-yl], methyl ester 17804522 4 U278 100 (45.4 Carbamic acid, (Inbutyl-1-1-inbutylethyl)-11-benzimidazol-yl ester 119380 4 P192 100 (45.4 Carbamic acid, (12-phenylenebis(minocarbonothicyl)bis-, dimethyl ester 2364058 4 U409 10 (45.4 Carbamic acid, (12-phenylenebis(minocarbonothicyl)bis-, dimethyl ester 2388809 4 U337 1000 (45.4 Carbamic acid, (12-phenylenebis(minocarbonothicyl)bis-, dimethyl ester 238809 4 U337 1000 (45.4 Carbami		17004332	*		10 (4.54)
1,3-Benzodioxol-4-ol, 2,2-dimethyl-, methyl carbamate 22781233 4 U278 100 (45.4 7-Benzoluranol, 2,3-dinydro-2,2-dimethyl- 1563388 4 U367 10 (45.4 Benzolic acid, 2-hydroxy- compd, with (3aS-cis)-1,2,3,3a,8,8-hexahydro-1,3a,5 57647 4 P188 100 (45.4 Carbamic acid, 11-benzimidazol-2-yl, methyl ester 10605217 4 U377 10 (45.4 Carbamic acid, [1-[butylamino]carboryl]-11-benzimidazol-2-yl], methyl ester 17804522 4 U278 100 (45.4 Carbamic acid, (Inbutyl-11-benzimidazol-2-yl], methyl ester 17804522 4 U271 10 (45.4 Carbamic acid, (Inbutyl-11-benzimidazol-2-yl], methyl ester 17804522 4 U278 100 (45.4 Carbamic acid, (Inbutyl-1-1-inbutylethyl)-11-benzimidazol-yl ester 119380 4 P192 100 (45.4 Carbamic acid, (12-phenylenebis(minocarbonothicyl)bis-, dimethyl ester 2364058 4 U409 10 (45.4 Carbamic acid, (12-phenylenebis(minocarbonothicyl)bis-, dimethyl ester 2388809 4 U337 1000 (45.4 Carbamic acid, (12-phenylenebis(minocarbonothicyl)bis-, dimethyl ester 238809 4 U337 1000 (45.4 Carbami	1.3-Benzodioxol-4-ol. 2.2-dimethyl-	22961826	4	U364	1000 (454)
Benzoic acid, 2-hydroxy-, compd. with (3aS-cis)-1,2,3,3a,8,8-hexahydro-1,3a,8- trimethylpyrrol2,3-8)Indol-5-yi methyl carbanate ester (1:1) 57647 4 P188 100 (45.4 Carbamic acid, 11-benzimidazol-2-yi, methyl ester 17804352 4 U372 10 (45.4 Carbamic acid, 11-benzimidazol-2-yi, methyl ester 17804352 4 U372 10 (45.4 Carbamic acid, 13-chlorophenyl-, 4-chloro-2-burynl ester 17804352 4 U271 10 (45.4 Carbamic acid, 13-chlorophenyl-, 4-chloro-2-burynl ester 17804352 4 P188 100 (45.4 Carbamic acid, 14-phenyl-neithyl-11-(methyl-thylp-ixc)-5-yi ester 118980 4 P192 100 (45.4 Carbamic acid, 112-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23564058 4 U373 1000 (45.4 Carbamic acid, 12-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23564058 4 U383 1000 (45.4 Carbamic acid, 12-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23564058 4 U383 1000 (45.4 Carbamic acid, 12-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23654058 4 U383 1000 (45.4 Carbamic acid, 12-chem			4		100 (45.4)
Benzoic acid, 2-hydroxy-, compd. with (3aS-cis)-1,2,3,3a,8,8-hexahydro-1,3a,8- trimethylpyrrol2,3-8)Indol-5-yi methyl carbanate ester (1:1) 57647 4 P188 100 (45.4 Carbamic acid, 11-benzimidazol-2-yi, methyl ester 17804352 4 U372 10 (45.4 Carbamic acid, 11-benzimidazol-2-yi, methyl ester 17804352 4 U372 10 (45.4 Carbamic acid, 13-chlorophenyl-, 4-chloro-2-burynl ester 17804352 4 U271 10 (45.4 Carbamic acid, 13-chlorophenyl-, 4-chloro-2-burynl ester 17804352 4 P188 100 (45.4 Carbamic acid, 14-phenyl-neithyl-11-(methyl-thylp-ixc)-5-yi ester 118980 4 P192 100 (45.4 Carbamic acid, 112-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23564058 4 U373 1000 (45.4 Carbamic acid, 12-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23564058 4 U383 1000 (45.4 Carbamic acid, 12-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23564058 4 U383 1000 (45.4 Carbamic acid, 12-phenylenebis(minocarbonothioy/l)bis-, dimethyl ester 23654058 4 U383 1000 (45.4 Carbamic acid, 12-chem	* * * * *		*	*	
trimethylpyrold[2.3-b]indol-5-yl methyl ester 57647 4 P188 100 (45.4 Carbamic add, 11-berzinidazol-2-yl, methyl ester 10605217 4 U372 10 (45.4 Carbamic add, (11-bulylaminocarbonyl-1-thenzinidazol-2-yll, methyl ester 10605217 4 U271 10 (45.4 Carbamic add, (12-bulyne) (10-biorehyl-2-bulynyl ester 101294 1028 100 (45.4 Carbamic add, (12-bulyne) (10-biorehyl-2-bulynyl ester 644644 P191 100 (45.4 Carbamic add, (12-bylenyl-1-thenkylethyl)-1-Hyyrazol-5-yl ester 644644 P191 100 (45.4 Carbamic add, (12-phenylenebis(minocarbonothioyl))bis-, dimethyl ester 23664058 U409 10 (45.4 Carbamic add, (12-phenylenebis(minocarbonothioyl))bis-, dimethyl ester 2303175 U389 1000 (45.4 Carbamic add, phenyl-, 1-methylethyl ester 2303175 U389 1000 (45.4 Carbamic add, phenyl-, 5-(phenylenebis(minocarbonothioyl)) ester 5288184 P189 1000 (45.4 Carbamic add, (12-methylethyl)- S-(2,3-trichioro-2-propenyl) ester 5288809 U387 5000 (2270 Carbamic add, phenyl-, S-(phenylenebis(minocarbonothyloyl) ester 5288148 P189 1000 (45.4 Carbourd phenol<	7-Benzofuranol, 2,3-dihydro-2,2-dimethyl-	1563388 *	4 *	U367 *	10 (4.54)
Carbamic acid, 1H-benzimidazol-2-yl, methyl ester 10605217 4 U372 10 (4 54 Carbamic acid, 1-(Lotylamino)carbonyl)-1H-benzimidazol-2-yl], methyl ester 17804352 4 U271 10 (4 54 Carbamic acid, (Gaboryhamino)-thiojmethyl-2,3-dihydro-2,2-dimethyl-7-benzofuranyl ester 108279 4 U280 10 (4 54 Carbamic acid, (Gaburyhamino)-thiojmethyl-2,3-dihydro-2,2-dimethyl-7-benzofuranyl ester 55285148 P189 1000 (454 Carbamic acid, imethyl-, 3-methyl-1-(1-methylethyl)-1H-pyrazol-5-yl ester 1129415 P190 1000 (454 Carbamic acid, fl-2-phenylenebis(iminocarbonothioyi))bis, dimethyl ester 23664058 U409 10 (4 54 Carbamotic acid, bis(1-methylethyl) S-(2,3-strichloro-2-propenyl) ester 230564058 U409 100 (454 Carbamotic acid, dipropyl-, S-(phenylmethyl) ester 2303175 U389 1000 (454 Carbanototic acid, bis(1-methylethyl)- S-(2,3-strichloro-2-propenyl) ester 230564058 U367 10 (4 54 Carbonatinic acid, dipropyl-, S-(phenylmethyl) ester 5288148 P189 1000 (454 Carbonatinic acid, dipropyl-, S-(phenylmethyl) ester 5952261 U337 5000 (227 <tr< td=""><td></td><td></td><td></td><td>_</td><td></td></tr<>				_	
Carbamic acid, 11-{[butylamino]carboryl]-11-benzmidazol-2-yl]-methyl ester 17804352 4 U271 10 (4.54 Carbamic acid, (3c-thorophenyl), 4. c-thioro-2-butyny ester 55285148 4 P189 1000 (454 Carbamic acid, (imethyl-1:(interthy-amino)-acbornyl)-7-benzohranyl ester 55285148 4 P191 10 (4.54 Carbamic acid, imethyl-1:(interthy-amino)-acbornyl-7-benzohranyl ester 1129415 4 P192 1000 (454 Carbamic acid, imethyl-1:(interthy-amino)-acbornyl-1-feyrazol-5-yl ester 1129415 4 P190 1000 (454 Carbamic acid, 11,2-prenylenebis/(minocarbonthioyl)bis-, dimethyl ester 2364058 4 U409 10 (4.54 Carbamic acid, phenyl-, 1-methylethyl, S-(2.3.3-trichloro-2-propenyl) ester 2303175 4 U389 100 (454 Carbamothioic acid, bi(-n-thylethyl), S-(2.3.3-trichloro-2-propenyl) ester 2303175 4 U387 5000 (227 Carbonthioic acid, bi(-n-thylethyl) S-(2.3.3-trichloro-2-propenyl) ester 52885148 4 P189 1000 (454 Carbonthioic acid, 2-(dimethyl-thyl-0-([(methylamino)-carbonyl]0xime 26419738 4 U395 5000 (227 <t< td=""><td>trimethylpyrrolo[2,3-b]indol-5-yl methylcarbamate ester (1:1)</td><td>57647</td><td>4 *</td><td>P188 *</td><td>100 (45.4)</td></t<>	trimethylpyrrolo[2,3-b]indol-5-yl methylcarbamate ester (1:1)	57647	4 *	P188 *	100 (45.4)
Carbamic acid, 11-{[butylamino]carboryl]-11-benzmidazol-2-yl]-methyl ester 17804352 4 U271 10 (4.54 Carbamic acid, (3c-thorophenyl), 4. c-thioro-2-butyny ester 55285148 4 P189 1000 (454 Carbamic acid, (imethyl-1:(interthy-amino)-acbornyl)-7-benzohranyl ester 55285148 4 P191 10 (4.54 Carbamic acid, imethyl-1:(interthy-amino)-acbornyl-7-benzohranyl ester 1129415 4 P192 1000 (454 Carbamic acid, imethyl-1:(interthy-amino)-acbornyl-1-feyrazol-5-yl ester 1129415 4 P190 1000 (454 Carbamic acid, 11,2-prenylenebis/(minocarbonthioyl)bis-, dimethyl ester 2364058 4 U409 10 (4.54 Carbamic acid, phenyl-, 1-methylethyl, S-(2.3.3-trichloro-2-propenyl) ester 2303175 4 U389 100 (454 Carbamothioic acid, bi(-n-thylethyl), S-(2.3.3-trichloro-2-propenyl) ester 2303175 4 U387 5000 (227 Carbonthioic acid, bi(-n-thylethyl) S-(2.3.3-trichloro-2-propenyl) ester 52885148 4 P189 1000 (454 Carbonthioic acid, 2-(dimethyl-thyl-0-([(methylamino)-carbonyl]0xime 26419738 4 U395 5000 (227 <t< td=""><td>Carbamic acid. 1H-benzimidazol-2-vl. methyl ester</td><td>10605217</td><td>4</td><td>U372</td><td>10 (4.54)</td></t<>	Carbamic acid. 1H-benzimidazol-2-vl. methyl ester	10605217	4	U372	10 (4.54)
Carbamic acid, (3-chlorophenyl), 4-chloro-2-butyny ester 10129 4 U280 10 (4.54 Carbamic acid, (dibutylamino)-thlormethyl-2-adimethyl-7-benzohranyl ester 55285148 4 P181 10 (4.54 Carbamic acid, (dimethyl-1-1[(dimethyl-amino)carbonyl)E-methyl-1+pyrazol-5-yl ester 644644 4 P191 10 (4.54 Carbamic acid, (imethyl-1, 3-methylphenyl ester 1129415 4 P180 1000 (454 Carbamic acid, (j 2-pherylenebis(minocarbonthoyl))bis-, dimethyl ester 23664058 4 U409 10 (4.54 Carbamic acid, (j 1-pheryl-henyl ester 23664058 4 U393 1000 (454 Carbamothicic acid, bis(1-methylethyl)- S-(2.3.3-richloro-2-propenyl) ester 2307175 4 U389 5000 (227 Carbanothicic acid, dipropyl-, S-(phenylmethyl) ester 5288148 4 P189 1000 (454 Carbanotalian 55285148 4 U395 5000 (227 Carbanotalian 55285148 4 P189 1000 (454 Carbanothicic acid, al, clamamate 5952261 U395 5000 (227 Carbosulfan 5952261 U395 </td <td></td> <td></td> <td>-</td> <td></td> <td>10 (4.54)</td>			-		10 (4.54)
Carbamic acid, (idiburytamino)-thio[methyl-12,3-dihydro-2,2-dimethyl-7-benzofuranyl ester 55285149 4 P189 1000 (454 Carbamic acid, dimethyl-1,3-methyl-1-1(-methylethyl)-1H-pyrazol-5-yl ester 1129415 4 P190 1000 (454 Carbamic acid, dimethyl-1,3-methyl-1-(1-methylethyl)-1H-pyrazol-5-yl ester 1129415 4 P190 1000 (454 Carbamic acid, phenyl-, 1-methylethyl ester 23664058 4 U409 10 (454 Carbamic acid, phenyl-, 1-methylethyl ester 2303175 4 U389 1000 (454 Carbamici acid, dincopt/, S-(phenylmethyl) ester 2303175 4 U389 100 (454 Carbanotinici acid, dipropt/, S-(phenylmethyl) ester 2303175 4 U387 100 (454 Carbanotinici acid, dipropt/, S-(phenylmethyl) ester 2303175 4 U387 100 (454 Carbonotin phenol 1563388 U387 100 (454 Carbonotin phenol 55226148 4 P189 1000 (454 Diethylene glycol, dicarbamate 64006 P202 10 (454 Libolane-2-carboxaldehyde, 2,4-dimethyl-1-0-[(methylamino)-carbonylloxime 26419738			4		10 (4.54)
Carbamic acid, dimethyl-, 3-methyl-1-(1-methylethyl)-1H-pyrazol-5-yl ester 119380 4 P192 100 (45.4 Carbamic acid, methyl-, 3-methylphenyl ester 1129415 4 P190 1000 (45.4 Carbamic acid, II, 2-phenylenebis(minocarbonothioyl)]bis-, dimethyl ester 23564058 4 U409 10 (45.4 Carbamchioic acid, bis(1-methylethyl) ester 2303175 4 U389 100 (45.4 Carbamchioic acid, dipropl., S-(2,3,3-trichloro-2-propenyl) ester 5288809 4 U372 100 (45.4 Carbamchioic acid, dipropl., S-(phenylmethyl) ester 2303175 4 U389 100 (45.4 Carbanothioic acid, dipropl., S-(phenylmethyl) ester 5288809 4 U372 10 (45.4 Carbourin phenol 1045517 4 U372 10 (45.4 Carbourin phenol 55285148 P189 1000 (45.4 m-Curmenyl methylcarbamate 64006 P202 10 (4.54 Diethylene glycol, dicarbamate 5952261 U395 5000 (2270 Dimetilan 644644 P191 1 (0.454 100 (45.4 1,3-Dithiolane-2-		55285148	4	P189	1000 (454)
Carbamic acid, dimethyl-, 3-methyl-1-(1-methylethyl)-1H-pyrazol-5-yl ester 119380 4 P192 100 (45.4 Carbamic acid, methyl-, 3-methylphenyl ester 1129415 4 P190 1000 (45.4 Carbamic acid, lif 2-phenylenebis(minocarbonothioyl)]bis-, dimethyl ester 23564058 4 U409 10 (45.4 Carbamchioic acid, bis(1-methylethyl) ester 2303175 4 U389 100 (45.4 Carbamchioic acid, dipropl., S-(2,3,3-trichloro-2-propenyl) ester 5288809 4 U372 10 (45.4 Carbamchioic acid, dipropl., S-(phenylmethyl) ester 106065217 4 U372 10 (45.4 Carbosulfan 106065217 4 U372 10 (45.4 Carbosulfan 55285148 P189 1000 (45.4 m-Cumenyl methylcarbamate 64006 P202 10 (4.54 Diethylene glycol, dicarbamate 5952261 U395 5000 (2270 Dimetilan 644644 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-0-((methylamino)-carbonyl]oxime 26419738 P185 100 (45.4 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl		644644	4	P191	1 (0.454)
Carbamic acid, [1,2-phenylenebis(iminocarbonothioyl)]bis-, dimethyl ester 23564058 4 U409 10 (4.54 Carbamic acid, phenyl-, 1-methylethyl ester 2303175 4 U389 100 (454 Carbamothioic acid, bis(1-methylethyl)-, S-(2,3,3-trichloro-2-propenyl) ester 2303175 4 U389 100 (454 Carbamothioic acid, dipropyl-, S-(phenylmethyl) ester 5288809 4 10372 10 (4.54 Carbonatzim 10605217 4 U387 5000 (2270 Carbonatzim 10665217 4 U372 10 (4.54 Carbosulfan 55285148 4 P189 1000 (454 m-Curmenyl methylcarbamate 5952261 4 U395 5000 (2270 Diethylene glycol, dicarbamate 5952261 4 U395 5000 (2270 Dimetilan 644644 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2.4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-H(methylamino)-carbonyl]oxyl-2-oxo-, methyl 2313520 P194 100 (45.4 Ethanimid			4		100 (45.4)
Carbamic acid, phenyl-, 1-methylethyl ester 122429 4 U373 1000 (454 Carbamothioic acid, dipropyl-, S-(2,3,3-trichloro-2-propenyl) ester 2303175 4 U389 100 (454 Carbamothioic acid, dipropyl-, S-(phenylmethyl) ester 10605217 4 U372 10 (454 Carbendzim 10605217 4 U372 10 (454 Carbondzim 1553388 4 U367 10 (454 Carbondzim 1553388 4 U377 10 (454 Carbondzim 55285148 4 P189 1000 (454 Carbondzim 64006 4 P202 10 (4.54 Diethylene glycol, dicarbamate 64006 4 P202 10 (4.54 J.a-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Libanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, NN-[thiobis[(methylamino)-carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4 Ethanimidothioic acid, N,N-[thiobis[(methylamino)-carbonyloxy]]bis-, dimethyl ester 5952261 <td>Carbamic acid, methyl-, 3-methylphenyl ester</td> <td>1129415</td> <td>* 4</td> <td>* P190</td> <td>1000 (454)</td>	Carbamic acid, methyl-, 3-methylphenyl ester	1129415	* 4	* P190	1000 (454)
Carbamic acid, phenyl-, 1-methylethyl ester 122429 4 U373 1000 (454 Carbamothioic acid, dipropyl-, S-(2,3,3-trichloro-2-propenyl) ester 2303175 4 U389 100 (454 Carbamothioic acid, dipropyl-, S-(phenylmethyl) ester 10605217 4 U372 10 (454 Carbendzim 10605217 4 U372 10 (454 Carbondzim 1553388 4 U367 10 (454 Carbondzim 1553388 4 U377 10 (454 Carbondzim 55285148 4 P189 1000 (454 Carbondzim 64006 4 P202 10 (4.54 Diethylene glycol, dicarbamate 64006 4 P202 10 (4.54 J.a-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Libanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, NN-[thiobis[(methylamino)-carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4 Ethanimidothioic acid, N,N-[thiobis[(methylamino)-carbonyloxy]]bis-, dimethyl ester 5952261 <td>* * * * *</td> <td></td> <td>*</td> <td>*</td> <td></td>	* * * * *		*	*	
Carbamothioic acid, bis(1-methylethyl)-, S-(2,3,3-trichloro-2-propenyl) ester 2303175 4 U389 100 (45.4 Carbamothioic acid, dipropyl-, S-(phenylmethyl) ester 10605217 4 U372 10 (4.54 Carboduran phenol 10605217 4 U372 10 (4.54 Carboduran phenol 55285148 4 P189 1000 (454 Carbosulfan 55285148 4 P189 1000 (454 m-Cumenyl methylcarbamate 64006 4 P202 10 (4.54 Diethylene glycol, dicarbamate 5952261 U395 5000 (2270 Dimetilan 644644 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 U394 5000 (2270 Ethanimidothioic acid, 2-(dimethylamino)-Carbonyloxyl]oxine 23135220 P194 100 (45.4 Ethanimidothioic acid, N,N'-[thiobis](methylimino)carbonyloxyl]bis-, dimethyl ester 59669260 U410 100 (45.4 Ethanimidothioic acid, N,N'-[thiobis](methylimino)carbonyloxyl]bis-, dimethyl ester 5952261 U395 5			-		10 (4.54)
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Carbamothioic acid, dipropyl-, Ś-(phenylmethyl) ester 52888809 4 U387 5000 (227C Carboduran phenol 1563384 4 U367 10 (4,54 Carboduran phenol 1563384 4 U367 10 (4,54 Carbonuran phenol 1563384 4 U367 10 (4,54 Carbonuran phenol 55285148 4 P189 1000 (454 Carbonuran phenol 5952261 4 U395 5000 (227C Diethylene glycol, dicarbamate 5952261 4 U395 5000 (227C Dimetilan 644644 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 U394 5000 (227C ester 30558431 U395 5000 (227C 2315220 P194 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-l((methylamino)carbonyl]oxyl-2-oxo-, methyl ester 30558431 U395 5000 (227C Formparanate 59669260 U410 100 (45.4 100 (45.4 Eth	* * * * * *		*	*	
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Carbofuran phenol 1563388 4 U367 10 (4.54 Carbosulfan 55285148 4 P189 1000 (454 m-Cumenyl methylcarbamate 64006 4 P202 10 (4.54 Diethylene glycol, dicarbamate 5952261 4 U395 5000 (2270 Dimetilan 644644 4 P191 1 (0.454 1.3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-[(methylamino)-carbonyl]oxy]-2-oxo-, methyl ester 30558431 U394 5000 (2270 Ethanimidothioic acid, 2-(dimethylamino)-N-[[(methylamino)carbonyl]oxy]-2-oxo-, methyl ester 30558431 U394 5000 (2270 Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester 59669260 U410 100 (45.4 Ethanol, 2,2'-oxybis-, dicarbamate 5952261 U395 5000 (2270 Formetanate hydrochloride 23422539 P198 100 (45.4 Isolan 1179380 P192 100 (45.4 Isolan 119380 P192 100 (45.4 Alaganese, bis(dimethylcarbamate 64006			-		
Carbosulfan 55285148 4 P189 1000 (454 m-Cumenyl methylcarbamate 64006 4 P202 10 (4.54 Diethylene glycol, dicarbamate 5952261 4 U395 5000 (2270 Dimetilan 644644 4 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, 2-(dimethylamino)-N-h[((methylamino)carbonyl]oxy]-2-oxo-, methyl 23135220 4 P194 100 (45.4 Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyl0xy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4 Ethanol, 2,2'-oxybis-, dicarbamate 5952261 U395 5000 (2270 Formparanate 17702577 4 P197 100 (45.4 Isolan 119380 4 P192 100 (45.4 Alsopopylphenyl N-methylcarbamate 64006 4 P202 10 (45.4 Manganese, bis(dimethylcarbamate 15339363 4 P196 10			-		
m-Cumenyl methylcarbamate 64006 4 P202 10 (4.54 Diethylene glycol, dicarbamate 5952261 4 U395 5000 (2270 Dimetilan 644644 4 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, N,N'-[thiobis[(methylamino)carbonylloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4 Ethanol, 2,2'-oxybis-, dicarbamate 5952261 U395 5000 (2270 Formparanate 17702577 P198 100 (45.4 Isolan 117380 P192 100 (45.4 Alsolan 119380 P192 100 (45.4 Manganese, bis(dimethylcarbamate 64006 P202 10 (4.54 Manganes	Carbofuran phenol	1563388	* 4	U367	10 (4.54)
Diethylene glycol, dicarbamate 5952261 4 U395 5000 (2270 Dimetilan 644644 4 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 Ethanimidothioic acid, 2-(dimethylamino)-N-lig(methylamino)carbonyl]oxy]-2-oxo-, methyl ester 30558431 4 U394 5000 (2270 23135220 4 P194 100 (45.4 Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4 Ethanol, 2,2'-oxybis-, dicarbamate 5952261 4 U395 5000 (2270 Formparanate 17702577 4 P198 100 (45.4 Isolan 119380 4 P192 100 (45.4 Alsopropylphenyl N-methylcarbamate 64006 4 P202 10 (45.4 Manganese, bis(dimethylcarbamate 15339363 4 P196 10 (4.54 Manganese dimethyldithiocarbamate 0.0 (45.4 10.0 (45.4 10.0	Carbosulfan	55285148	4	P189	1000 (454)
Dimetilan 644644 4 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270) Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270) ester 23135220 4 P194 100 (45.4) Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonylloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4) Ethanol, 2,2'-oxybis-, dicarbamate 5952261 4 U395 5000 (2270) Formetanate hydrochloride 23422539 4 P194 100 (45.4) Isolan 119380 4 P192 100 (45.4) Isolan 119380 4 P192 100 (45.4) Manganese, bis(dimethylcarbamate 64006 4 P202 10 (4.54) Manganese dimethyldithiocarbamate 15339363 4 P196 10 (4.54) Manganese dimethyldithiocarbamate 15339363 4 P196 <t< td=""><td>m-Cumenyl methylcarbamate</td><td>64006</td><td>* 4</td><td>* P202</td><td>10 (4.54)</td></t<>	m-Cumenyl methylcarbamate	64006	* 4	* P202	10 (4.54)
Dimetilan 644644 4 P191 1 (0.454 1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4 Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270) Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270) ester 23135220 4 P194 100 (45.4) Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonylloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4) Ethanol, 2,2'-oxybis-, dicarbamate 5952261 4 U395 5000 (2270) Formetanate hydrochloride 23422539 4 P194 100 (45.4) Isolan 119380 4 P192 100 (45.4) Isolan 119380 4 P192 100 (45.4) Manganese, bis(dimethylcarbamate 64006 4 P202 10 (4.54) Manganese dimethyldithiocarbamate 15339363 4 P196 10 (4.54) Manganese dimethyldithiocarbamate 15339363 4 P196 <t< td=""><td>Niethylene glycel diaethamate</td><td>5052261</td><td>*</td><td>*</td><td>. ,</td></t<>	Niethylene glycel diaethamate	5052261	*	*	. ,
1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[((methylamino)-carbonyl]oxime 26419738 4 P185 100 (45.4) Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270) Ethanimidothioic acid, 2-(dimethylamino)-N-[[(methylamino)carbonyl]oxy]-2-oxo-, methyl 23135220 4 P194 100 (45.4) Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4) Ethanol, 2,2'-oxybis-, dicarbamate 5952261 4 U395 5000 (2270) Formparanate 5952261 4 U395 5000 (2270) Formparanate 17702577 4 P198 100 (45.4) Isolan 119380 4 P192 100 (45.4) Alsopropylphenyl N-methylcarbamate 64006 4 P202 10 (4.54) Manganese dimethyldithiocarbamate 15339363 4 P196 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, monohydrochloride 23422539 4 P196 10 (4.54) Monganese dimethyldithiocarbamate 10 (4.54) 15339363 4 P196 </td <td></td> <td>5952261</td> <td>*</td> <td>U395 *</td> <td>5000 (2270)</td>		5952261	*	U395 *	5000 (2270)
Ethanimidothioic acid, 2-(dimethylamino)-N-hydroxy-2-oxo-, methyl ester 30558431 4 U394 5000 (2270) Ethanimidothioic acid, 2-(dimethylamino)-N-[[(methylamino)carbonyl]oxy]-2-oxo-, methyl ester 23135220 4 P194 100 (45.4) Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4) Ethanol, 2,2'-oxybis-, dicarbamate 5952261 U395 5000 (2270) Formetanate hydrochloride 23422539 4 P198 100 (45.4) Isolan 119380 P192 100 (45.4) Isolan 119380 P192 100 (45.4) Manganese, bis(dimethylcarbamate 64006 P202 10 (45.4) Manganese dimethylcarbamate 15339363 P196 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, 23422539 4 P196 10 (4.54)	Dimetilan	644644	4 *	P191 *	1 (0.454)
Ethanimidothioic acid, 2-(dimethylamino)-N-[[(methylamino)carbonyl]oxy]-2-oxo-, methyl 23135220 4 P194 100 (45.4) Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4) Ethanol, 2,2'-oxybis-, dicarbamate 5952261 4 U395 5000 (2270) Formetanate hydrochloride 23422539 4 P198 100 (45.4) Formparanate 117702577 4 P197 100 (45.4) Isolan 119380 4 P192 100 (45.4) Manganese, bis(dimethylcarbamate 64006 4 P202 10 (45.4) Manganese dimethyldithiocarbamate 15339363 4 P196 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, 23422539 4 P198 100 (45.4) 100 (45.4) 15339363 4 P196 10 (4.54) 10 (4.54) Manganese dimethyldithiocarbamate 23422539 4 P196 10 (4.54) Manganese dimethyldithiocarbamate 23422539 4 P196 10 (4.54) Monohydrochloride N,N-dimeth	1,3-Dithiolane-2-carboxaldehyde, 2,4-dimethyl-, O-[(methylamino)-carbonyl]oxime	26419738	4	P185	100 (45.4)
ester 23135220 4 P194 100 (45.4 Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester 59669260 4 U410 100 (45.4 Ethanol, 2,2'-oxybis-, dicarbamate 59552261 4 U395 5000 (2270) Formetanate hydrochloride 23422539 4 P198 100 (45.4) Formparanate 17702577 4 P197 100 (45.4) Isolan 119380 4 P192 100 (45.4) Alsolan 119380 4 P192 100 (45.4) Manganese, bis(dimethylcarbamate * 64006 4 P202 10 (4.54) Manganese dimethyldithiocarbamate * 15339363 4 P196 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, 23422539 4 P196 10 (4.54) Monohydrochloride * * * 15339363 4 P196 10 (4.54) Monohydrochloride * * * * 10 (4.54) * * * * * * *		30558431	* 4	* U394	5000 (2270)
************************************		23135220	4	P194	100 (45.4)
* * * * 5952261 4 U395 5000 (2270 Formetanate hydrochloride * * 23422539 4 P198 100 (45.4) Formparanate * * * * 100 (45.4) Formparanate * * * 100 (45.4) Isolan * * * * * * * * * * * * * * * * * * * * * *	* * * * *		*	*	,
Formetanate hydrochloride *<	Ethanimidothioic acid, N,N'-[thiobis[(methylimino)carbonyloxy]]bis-, dimethyl ester	59669260	4 *	U410 *	100 (45.4)
Formetanate hydrochloride *<	Ethanol, 2,2'-oxybis-, dicarbamate	5952261	4 *	U395	5000 (2270)
Isolan 119380 4 P192 100 (45.4) 3-Isopropylphenyl N-methylcarbamate * * 64006 4 P202 10 (4.54) Manganese, bis(dimethylcarbamodithioato-S,S')- * * 15339363 4 P196 10 (4.54) Manganese dimethyldithiocarbamate * * * 15339363 4 P196 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, 23422539 4 P198 100 (45.4)		23422539	4	P198	100 (45.4)
Isolan 119380 4 P192 100 (45.4) 3-Isopropylphenyl N-methylcarbamate * * 64006 4 P202 10 (4.54) Manganese, bis(dimethylcarbamodithioato-S,S')- * * 15339363 4 P196 10 (4.54) Manganese dimethyldithiocarbamate * * * 15339363 4 P196 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, 23422539 4 P198 100 (45.4)	Formparanate	17702577	* 4	* P197	100 (45.4)
* * * * * * * * * * * * * * * * * * *			*	*	. ,
Manganese, bis(dimethylcarbamodithioato-S,S')- * * * * * * * 10 (4.54) Manganese dimethyldithiocarbamate * * * 15339363 4 P196 10 (4.54) Manganese dimethyldithiocarbamate * * 10 (4.54) * 10 (4.54) Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, * 10 (4.54) Monohydrochloride * * * * 10 (4.54)	* * * * * *		*	F 192 *	100 (45.4)
* * * * * * * * * * * * * * * * * * *	3-Isopropylphenyl N-methylcarbamate	64006	4 *	P202	10 (4.54)
Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, monohydrochloride	Manganese, bis(dimethylcarbamodithioato-S,S')	15339363	* 4	P196	10 (4.54)
Methanimidamide, N,N-dimethyl-N'-[3-[[(methylamino)- carbonyl]oxy]phenyl]-, monohydrochloride	Manganese dimethyldithiocarbamate	15339363	4 *	P196	10 (4.54)
monohydrochloride				-	
		23422539	4	P198	100 (45.4)
	Methanimidamide, N,N-dimethyl-N'-[2-methyl-4-[[(methylamino)carbonyl]oxy]phenyl]	17702577	4	P197	100 (45.4)

TABLE 302.4.—LIST OF HAZARDOUS SUBSTANCES AND REPORTABLE QUANTITIES—Continued [Note: All Comments/Notes Are Located at the End of This Table]

	Hazardous subst	ance		CASRN	Statutory code†	RCRA waste number	Final RQ pounds (Kg)
* Metolcarb	*	*	* *	1129415	* 4	* P190	1000 (454)
* Oxamyl	*	*	* *	23135220	* 4	* P194	100 (45.4)
* Phenol, 3-(1-methylethyl)-, meth Phenol, 3-methyl-5-(1-methyleth				64006 2631370	* 4 4	* P202 P201	10 (4.54) 1000 (454)
* Physostigmine Physostigmine salicylate				57476 57647	* 4 4	* P204 P188	100 (45.4) 100 (45.4)
Promecarb	*	*	* *	2631370	* 4	* P201	1000 (454)
Propanal, 2-methyl-2-(methyl-su	, O–[(methy	* lamino)carbonyl] ox	ime	1646884	* 4	P203	100 (45.4)
Propham	*	*	* *	122429	4	 U373	1000 (454)
Prosulfocarb	*	*	* *	52888809	4	U387	5000 (2270)
Pyrrolo[2,3-b]indol-5-ol, 1,2,3 (ester), (3aS-cis)				57476	4	P204	100 (45.4)
* Thiodicarb	*	*	* *	59669260	* 4	* U410	100 (45.4)
* Thiophanate-methyl	*	*	* *	23564058	* 4	* U409	10 (4.54)
Tirpate	*	*	* *	26419738	* 4	* P185	100 (45.4)
Triallate	*	*	* * *	2303175	* 4		100 (45.4)
Zinc, bis(dimethylcarbamodithio	م. ato-S,S′)	*	* *	137304	^ 4	P205	10 (4.54)
Ziram*	*	*	* *	137304	4 *	P205	10 (4.54)
K156 Organic waste (including heavy and decantates) from the pro ing does not apply to wastes butvlcarbamate.)	y ends, still bottom	s, light ends, spent ates and carbamoy	t solvents, filtrates, oximes. (This list-		4	K156	10 (4.54)
K157 Wastewaters (including scrubbe waters) from the production of not apply to wastes gene butylcarbamate.)	er waters, condense of carbamates and	er waters, washwate carbamoyl oximes	ers, and separation . (This listing does		4	K157	10 (4.54)
K158 Bag house dusts and filter/sepa bamoyl oximes. (This listing d of 3- iodo-2-propynyl n- butylo	aration solids from t does not apply to w	he production of ca	arbamates and car-		4	K158	10 (4.54)
K159 Organics from the treatment of	, ,				4	K159	10 (4.54)
Purification solids (including filtr dust and floor sweepings from (This listing does not include	ration, evaporation, n the production of	and centrifugation	solids), bag-house		4	K161	1 (0.454)
K178 Nonwastewaters from the produ [This listing does not apply to production exempt under sect	uction of titanium di o chloride process v tion 261.4(b)(7).].	oxide by the chlorid vaste solids from tit	e-ilmenite process.		* 4	* K178	1000 (454)

+ Indicates the statutory source as defined by 1, 2, 3, and 4 below.

4- Indicates that the statutory source for designation of this hazardous substance under CERCLA is RCRA section 3001. 1* Indicates that the 1-pound RQ is a CERCLA statutory RQ.

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3. Appendix A to § 302.4 is amended by revising the following entries, to read as follows:

APPENDIX A TO § 302.4—SEQUENTIAL CAS REGISTRY NUMBER LIST OF CERCLA HAZARDOUS SUBSTANCES

			Hazardous su	bstance		
*	*	*	*	*	*	*
57476 57647	Physostigmine. Pyrrolo[2,3-b]indol-5-ol, 1,2,3,3a,8,8a Benzoic acid, 2-hydroxy-, compd. v ester (1:1). Physostigmine salicylate.					5-yl methylcarbamat
* 64006	* m-Cumenyl methylcarbamate. 3-Isopropylphenyl N-methylcarbamate Phenol, 3-(1-methylethyl)-, methyl car		*	*	*	*
101279	Barban. Carbamic acid, (3-chlorophenyl)-, 4-c	hloro-2-butyn	yl ester.	-	-	-
* 119380	* Carbamic acid, dimethyl-, 3-methyl-1- Isolan.	(1-methyleth	* yl)-1H-pyrazol-5-yl و	* ester.	*	*
* 122429	* Carbamic acid, phenyl-, 1-methylethy Propham.	* I ester.	*	*	*	*
* 137304	* Zinc, bis(dimethylcarbamodithioato-S, Ziram.	* S′)	*	*	*	*
* 644644	* Carbamic acid, dimethyl-,1-[(dimethyl- Dimetilan.	* -amino)carbo	* nyl]-5-methyl-1H-py	* vrazol-3-yl ester.	*	*
* 1129415	* Carbamic acid, methyl-, 3-methylpher Metolcarb.	* nyl ester.	*	*	*	*
* 1563388	* 7-Benzofuranol, 2,3-dihydro-2,2-dime Carbofuran phenol.	* thyl	*	*	*	*
* 1646884	* Aldicarb sulfone. Propanal, 2-methyl-2-(methyl-sulfonyl	*)-, O-[(methy	* lamino)carbonyl] ox	* ime.	*	*
* 2303175	* Carbamothioic acid, bis(1-methylethyl Triallate.	*)-, S-(2,3,3-tr	* richloro-2-propenyl)	* ester.	*	*
* 2631370	* Phenol, 3-methyl-5-(1-methylethyl)-, r Promecarb.	* nethyl carbar	* nate.	*	*	*
* 5952261	* Ethanol, 2,2'-oxybis-, dicarbamate. Diethylene glycol, dicarbamate.	*	*	*	*	*
* 10605217	* Carbamic acid, 1H-benzimidazol-2-yl, Carbendazim.	* methyl ester	*	*	*	*
* 5339363	* Manganese, bis(dimethylcarbamodith Manganese dimethyldithiocarbamate.		*	*	*	*
* 7702577	*	*	* nethylamino)carbon	* vlloxvlphenvll-	*	*
7804352	Benomyl. Carbamic acid, [1-[(butylamino)carbon				*	*
2781233	Bendiocarb. 1,3-Benzodioxol-4-ol, 2,2-dimethyl-, n	nethyl carban	nate.			
22961826 23135220	Bendiocarb phenol. 1,3-Benzodioxol-4-ol, 2,2-dimethyl Ethanimidothioic acid, 2-(dimethylami	no)-N- [[(met	hylamino)carbonyl]c	oxy]-2-oxo-, methyl es	ster.	
23422539	Oxamyl. Methanimidamide, N,N-dimethyl-N'-[3 Formetanate hydrochloride.					
00004000	Carbamic acid, [1,2-phenylenebis(imit	nocarbonothi	oyl)]bis-, dimethyl e	ster.		
23564058	Thiophanate-methyl.	*	*	*	*	*

APPENDIX A TO § 302.4—SEQUENTIAL CAS REGISTRY NUMBER LIST OF CERCLA HAZARDOUS SUBSTANCES—Continued

CASRN	Hazardous substance						
*	*	*	*	*	*	*	
30558431	Ethanimidothioic acid, 2-(dime A2213.	ethylamino)-N-hydrox	ky-2-oxo-, methyl este	er.			
*	*	*	*	*	*	*	
52888809	Carbamothioic acid, dipropyl- Prosulfocarb.	, S-(phenylmethyl) e	ster.				
*	*	*	*	*	*	*	
55285148	Carbamic acid, [(dibutylaminc Carbosulfan.)-thio]methyl-, 2,3-di	hydro-2,2-dimethyl-7-	benzofuranyl ester.			
*	*	*	*	*	*	*	
59669260	Ethanimidothioic acid, N,N'-[the second seco	niobis[(methylimino)c	arbonyloxy]]bis-, dime	ethyl ester.			
*	*	*	*	*	*	*	

PART 355—EMERGENCY PLANNING AND NOTIFICATION

Authority: 42 U.S.C. 11002, 11004, and 11048.

entries, to read as follows (footnotes "*" and "h" have been republished without change):

1. The authority citation for part 355 continues to read as follows:

2. Appendices A and B in part 355 are amended by revising the following

APPENDIX A TO PART 355—THE LIST OF EXTREMELY HAZARDOUS SUBSTANCES AND THEIR THRESHOLD PLANNING QUANTITIES

[Alphabetical Order]

CAS No.	Chemical name	Notes	Reportable quantity* (pounds)	Threshold planning quan- tity (pounds)
*	* * *	*	*	
26419–73–8	Carbamic Acid, Methyl-, O-(((2,4–Dimethyl-1, 3-Dithiolan-2 yl)Methylene)Amino)	2	100	100/10,000
*	* * * * * *	*	*	
644–64–4*	Dimetilan	*	1 *	500/10,000
23422–53–9	Formetanate Hydrochloride	(^h)	100	500/10,000
17702–57–7*	* Formparanate* * *	*	100 *	100/10,000
119–38–0*	sopropylmethyl-pyrazolyl Dimethylcarbamate	 *	100 *	500
1129–41–5*	* Metolcarb * * *	 *	1,000	100/10,000
23135–22–0*	* Oxamyl * * *	 *	100 *	100/10,000
64–00–6*	* Phenol, 3-(1–Methylethyl)-, Methylcarbamate	 *	10 *	500/10,000
	* Physostigmine Physostigmine, Salicylate (1:1)		100 100	100/10,000 100/10,000
- 2631–37–0	* * * * Promecarb		* 1,000 *	500/10,000 *

*Only the statutory or final RQ is shown. For more information, see 40 CFR Table 302.4.

Notes:

^h Revised TPQ based on new or re-evaluated toxicity data.

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APPENDIX B TO PART 355—THE LIST OF EXTREMELY HAZARDOUS SUBSTANCES AND THEIR THRESHOLD PLANNING QUANTITIES [CAS Number Order]

CAS No.	Chemical name					Reportable quantity* (pounds)	Threshold planning quan- tity (pounds)**
*	*	*	*	*		*	*
57–47–6	Physostigmine	*	*	*		* 100	100/10,000
57–64–7 *	Physostigmine, Salicylate (1:1)	*	*	*		* 100	100/10,000 *
64-00-6	Phenol, 3-(1-Methylethyl)-, Methylo					* 10	500/10,000 *
119–38–0 *	Isopropylmethyl-pyrazolyl Dimethy	carbamate	*	*		* 100	* 500
644–64–4 *	Dimetilan*	*	*	*		* 1	500/10,000 *
1129–41–5 *	Metolcarb*	*	*	*		* 1,000	100/10,000
2631–37–0 *	Promecarb*	*	*	*	(^h)	* 1,000	500/10,000 *
17702–57–7 *	Formparanate	*	*	*		* 100	100/10,000 *
23135–22–0 23422–53–9 *	Oxamyl Formetanate Hydrochloride				(^h)	100 100	,
26419–73–8 *	Carbamic Acid, Methyl-, O-(((2,4-D)imethyl-1, 3- Di	thiolan-2- yl)N *	ethylene)Amino)		* 100	100/10,000 *
*Only the sta	tutory or final RQ is shown. For mo	re information, s	see 40 CFR Ta	able 302.4.	*		
^h Revised TP *	Q based on new or re-evaluated to	xicity data.		*	*		

[FR Doc. 03–30166 Filed 12–3–03; 8:45 am] BILLING CODE 6560–50–P