# Will Deregulating Michigan's Toxic Air Emissions Put Residents at Risk? 

Backgrounder on the Proposed Administrative Rule Change

## ACKNOWLEDGEMENTS

## About the Michigan Environmental Council

Since 1980, the Michigan Environmental Council has been at the forefront of efforts to protect our Great Lakes, promote sustainable cities, safeguard public health and establish clean energy policies for a more vibrant economy. Representing over 65 member organizations throughout the state, MEC provides agenda-setting leadership at the State Capitol and with our Congressional delegation in Washington.

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## Recommendation to Deregulate Emissions of Over 500 Toxic Chemicals

Nine states including Michigan do not use a discrete list of toxic chemicals and therefore their programs regulate all toxic chemicals proposed to be emitted. Those states include Minnesota, Delaware, Georgia, Maryland, New Jersey, Oklahoma and Texas.

In Michigan, the regulated list currently includes about 1200 chemicals. When evaluating the potential impact that a chemical may have on the adjoining community, the state considers:

1. The toxicity of the chemical,
2. The quantity being emitted, and
3. How close the industrial facility or specific stack is to nearby homes.

However, the administration is currently assessing a proposal that would reduce the number of toxic chemicals regulated by more than 500 (the list of chemicals is included with the Air Toxics Workgroup report - Appendix B). Those chemicals fall into two categories:

1. Eliminating regulation of toxic chemicals that have not been tested for their impact on public health. Michigan's current regulation assumes any chemical that has not been tested for health impacts is very toxic and a default value is used for evaluating its impact on the local community. The permit application has the choice of using the default value or conducting basic health testing on the chemical to establish a specific value.
2. Eliminating regulation of the less toxic non-carcinogen chemicals (regardless of the quantity being emitted). The proposal arbitrarily draws the line based on toxicity and proposes deregulating the $25 \%$ least toxic chemicals currently regulated. Importantly, the chemicals would not be regulated regardless of the quantity being emitted.

## Concerns with the Proposed Changes

## 1. The changes fail to protect the public health of Michigan families

Under the proposal, industrial facilities will be allowed to emit chemicals that have not been tested for their impact on human health or natural resources. This makes Michigan families the equivalent of guinea pigs. It is our position that the company using the chemical (and reaping the financial benefits of its use) should bear the burden of demonstrating it is safe before emitting it into the air we breathe. Instead this proposal transfers to those living next to the factory the risk that the chemical can cause cancer or have other negative consequences.

Under current regulations, state regulators can credibly tell residents they have looked at the public health aspects of a new factory or proposed expansion and are basing their issuance of a permit on the demonstration that it will be safe for the community. If this proposal is adopted, regulators would have no basis to claim they have thoroughly examined its impact on public health. This has the potential to result in greater conflict between industrial facilities and
adjoining residents, and greater community resistance to new factories proposed in their communities.

## 2. The changes are not supported by sound science

There is no science behind the proposed change. The proposal to deregulate chemicals for which no safety data exists goes against our knowledge of toxic chemicals. The current program at least creates a presumption that an untested chemical is fairly toxic. If modeling shows that emitting a chemical is safe due to the quantity emitted, the state can issue a permit satisfied it has performed its duty to protect the health and safety of its residents.

The second category of chemicals being deregulated are those that have been found not to cause cancer and are less toxic than other chemicals, but which can still have impacts on public health. However, as explained above, the potential impact on human health is driven by both the toxicity of the chemical and the quantity of the chemical being emitted. The second category ignores this question of quantity and deregulates a chemical based solely on its toxicity. This change is also contrary to the science behind protecting people from the impacts of toxic chemicals.

## 3. Changes will have disproportionate impact on low-income areas and communities of color

Numerous studies have shown that residential neighborhoods next to industrial areas tend to have below-average income and have a greater likelihood to be communities of color ${ }^{1}$. By deregulating more than 500 chemicals the proposal will have the greatest impact in those communities with the highest concentration of industrial facilities and toxic air emissions. Residents in these areas of the state already are at greater risk because our program does not take into account the impacts of multiple pollutants from multiple sources when setting acceptable emissions limits. Further deregulating individual toxic chemicals will place these communities at even greater risk.

1 Race, Income, and Environmental Inequality in the United States, Liam Downey and Brian Hawkins, Sociol Perspect. Dec 1, 2008; 51(4): 759-781, doi: 10.1525/sop.2008.51.4.759

Racial and Socioeconomic Disparities in Residential Proximity to Polluting Industrial Facilities: Evidence From the Americans' Changing Lives Study, Paul Mohai, PhD, Paula M. Lantz, PhD, Jeffrey Morenoff, PhD, James S. House, PhD, and Richard P. Mero, MS, Am J Public Health. 2009 November; 99(Suppl 3): S649-S656., doi: 10.2105/AJPH.2007.131383

## 4. Department failed to consider the health and safety benefits of the current rule

The Executive Order that prompted the review of the Air Toxics rule enumerated seven factors that were supposed to be considered when reviewing an existing administrative rule. The first factor to be evaluated was the "health or safety benefits of the rules." In this case the department conducted no assessment of the potential health and safety risks that could be presented by deregulating over 500 toxic chemicals in Michigan. This failure to evaluate the potential impacts will place Michigan residents at risk.

## Where Did These Recommendations Come From?

The Environmental Advisory Rules Committee (ARC) was created by the Office of Regulatory Reinvention (ORR) in accordance with Executive Order 2011-5. The mission of the ORR is to ensure that Michigan's regulatory environment is simple, fair, efficient, and conducive to business growth and job creation. The purpose of the Environmental ARC was to produce recommendations to the ORR for changes to Michigan's existing environmental regulations.

Evaluations and recommendations were based on the application of the seven factors described in Executive Order 2011-5. Those seven factors are as follows:

1. Health or safety benefits of the rules;
2. Whether the rules are mandated by any applicable constitutional or statutory provision;
3. The cost of compliance with the rules, taking into account their complexity, reporting requirements and other factors;
4. The extent to which the rules conflict with or duplicate similar rules or regulations adopted by the state or federal government;
5. Extent to which the regulations exceed national or regional compliance requirements or other standards;
6. Date of last evaluation of the rules and the degree, if any, to which technology, economic conditions or other factors have changed regulatory activity covered by the rules since the last evaluation; and
7. Other changes or developments since implementation that demonstrate there is no continued need for the rules.

In December of 2011 the ARC issued its final recommendations to the administration. The recommendations within each category were listed in order of priority, based on the importance of the recommendation to the state's future.

Under the air program, the first recommendation addressed was the issue of the regulation of air toxics emitted from industrial facilities (Appendix A).

The ORR review of environmental rules led to the formation of a number of workgroups. One of those workgroups, the Air Toxics Workgroup (ATW) of the Air Quality Division, reviewed

Michigan’s Air Toxic Program and made a number of specific recommendations. The report of the workgroup is attached as Appendix B.

A summary of the history of Michigan's Air Toxic Program is included in workgroup report. In general the program has been operating since the early 1980s and has been reviewed by been the subject of three separate stakeholder groups, the last one in the late 1990s which resulted in our current program.

In Michigan for several decades, a source emitting any toxic air contaminant, not specifically exempted, had both the duty to identify and characterize such an emission in a quantitative manner, and then perform a community health risk assessment by evaluating the ambient consequences of those emissions against a system of screening values. By reducing the number of chemicals covered by the rules this change significantly weakens the community health risk assessment performed by the applicant.

## Air Quality Recommendations

## No. A-1

Subject: Air Toxics Rules
Regulation: R 336.1224-R 336.1232
Remedy: $\square$ Process $\boxtimes$ Rules $\square$ Statute

Background/Issue: In 1992, Michigan approved state-only air toxics regulations and, for the most part, they remain in effect today. The rule development and approval process required several years of prior discussions with industrial and environmental groups, well before passage of the federal 1990 Clean Air Act (CAA) amendments. Prior to passage of the CAA amendments, the federal regulations concerning potential air toxics were limited to approximately 5 compounds.

Title III of the 1990 Clean Air Act Amendments established a national regulatory program to minimize the emissions of the most significant air toxics. Since passage of the 1990 CAA, the federal government has developed numerous Maximum Achievable Control Technology (MACT) determinations for a wide variety of processes that typically emit hazardous air pollutants (HAPS). Any MACT-subject source is required to comply with (1) the federal emission limitations, (2) specified emissions control technologies, and (3) specific monitoring, testing, and reporting requirements. Therefore, Michigan's outdated air toxics regulations are in need of significant reform.

In developing each MACT standard, USEPA focused on the most significant HAPS emitted from a specific process type, then developed emission control strategies that are as stringent as the control systems required to meet the intent of Michigan Rule 224. Not all potential air pollutants were evaluated by USEPA during development of the MACT standards, nor should they have been. The USEPA regulations focus on specific organic and non-organic HAPS, while the Michigan program focuses on each individual pollutant. The emission control systems required by MACT are as stringent as those required by Rule 224, although MACT uses emission surrogates such as CO and PM to certify reductions in potential HAPS, whereas Michigan's air toxics regulations review each potential pollutant individually. The end results for determining the adequacy of a proposed emissions control system is essentially the same.

Recommendations: The Committee makes the following recommendations regarding Air Toxic rules:

- The parts of $R 336.1224$ dealing with compounds that are considered volatile organic compounds (VOCs) should be rescinded. Portions of R 336.1224 are redundant because R 336.1702 requires a control technology review for VOCs. VOC-based emission control is more effective under R 336.1702 and this entire regulation exceeds federal standards.
- Rule 336.1225 should be amended to specifically include the following:
- Limit permit modification reviews to those increases in a Hazard Index exceeding $10 \%$ above the previously permitted baseline.
- Exempt sources that are identified in a MACT source category.
- Exempt clean fuels such as natural gas, low sulfur \#2 Fuel Oil, and nonchemically treated biofuels.
- Exempt pollution control projects for existing sources from the air toxic regulations.
- Limit the number of air toxics to the federal HAPS list.
- Make the acceptable exposure limits consistent with other nearby states.
- Stop requiring permit holders to conduct elaborate and costly stack tests to provide emissions research data, since the DEQ does not use this information for subsequent permit reviews.
- R 336.1228 should be rescinded. This rule allows the Air Quality Division to go beyond the requirements of the rule for any reason.

Rationale/Comments: These proposed revisions to Michigan's air toxics regulations would serve to level the playing field with other states vying for additional industrial growth, and would not result in a back-sliding of the environmental programs (see supplemental document Issue A-1, Attachment 1). The current system of reviewing the impacts of every feasible, potential air toxic results in nothing more than a "numbers game" that only serves to heighten public anxiety, delay permit issuance and waste several hundreds of thousands of dollars for stack testing that could be put to better use for industrial expansions, process improvements and other more beneficial programs. Please review supplemental document Issue A-1, Attachments 2 and 2a related to the Public Participation Document for Frontier Renewable Resources PTI 166-09. You can see that only 6 of the 37 identified air toxics (based on the firing of natural gas) have a predicted air quality impact at least $50 \%$ of their individual ambient limitations. Furthermore, the Permit to Install document for the Mancelona Renewable Resources project contained in Issue A-1, Attachment 3 requires the company to spend tens of thousands of additional dollars to perform stack testing for 28 potential air toxics (the majority of which are predicted to be emitted in trace levels) generated from the burning of nonchemically treated biomass (wood).

This has also been a consistent issue for Michigan's Asphalt Plant industry for the past 20+ years (see supplemental document Issue A-1, Attachment 4). The AQD has required stack testing for compounds that have proven to have ambient impacts well below any state or federal standard. Had the AQD invested the time to compile the results of these historical stack tests, which have been consistently submitted to AQD for the past 20+ years, it would be readily apparent that the vast majority of these air toxics of concern would no longer be an issue. This would result in a more expedited air permitting process, reduce stack testing requirements for new Asphalt Plants and save companies several thousands of dollars in unnecessary stack testing.

Supplemental document Issue A-1, Attachment 5, provides a summary of issues related to a $2+$ year delay in issuance of an air permit for use of Biodiesel in a large utility boiler when the same fuel is widely used in this company's fleet vehicles.

Supporting document Issue A-1, Attachments 6 and 6a relate to air toxics issues for raw material and fuel substitutions that can be expanded to several other types of manufacturing and combustion processes. As noted in our final recommendations, we believe a Hazard Indexing
methodology be established by rule to allow for raw material / fuel substitutions that do not have a detrimental ambient impact. This is consistent with our recommendation in A-1, Rule 336.1225.

James Clift Comments: The environmental community believes that a company or person who wants to emit a toxic chemical into the environment should have the duty to demonstrate that emission of that toxic chemical will not adversely impact natural resources or public health. By requiring testing before chemical are released, we are providing companies with the incentive to develop safer alternatives through green chemistry.

If the federal government has developed a MACT, BACT or LEAR standard for a chemical, we support that standard being applied as the Michigan standard. However, if such a standard has not been developed, we believe the source should be subject to T-BACT. Exempting all chemicals not on the list federal list of hazardous air pollutants could provide an unwise incentive for companies to use chemicals for which there has been less testing and analysis versus the use of less toxic alternatives.

We support Rule 228 which is designed to protect Michigan citizens from persistent bioaccumative toxic chemicals.

# Air Toxics Workgroup <br> "TAC List" Discussion Paper - DRAFT <br> August 20, 2013 UPDATE 

## ORR (2011) Report Recommendation A-1(6):

R 336.1225 should be amended and specifically include the following: Limit the number of air toxics to the federal HAPs list.

## ATW Initial Discussion

Discussion of the "TAC list" issue at the 3/5/13 ATW meeting indicated that the "status quo" is characterized by some ATW members as burdensome and more extensive than other Region 5 state's programs. However, there are also reservations about the sufficiency of the HAPs list. And if the DEQ were to adopt a defined list of TACs for R225 applicability, then staff asked about a mechanism to ensure public health protection if health concerns are posed by the proposed emission of an unlisted compound. ATW members voted, using the "gradient of agreement" tool, on three options: 1. HAPs only; 2. HAPs plus, including a caveat to add other compounds; and, 3. maintaining the status quo. Although there were varied levels of acceptability for each option, the voting was relatively polarized for options 1 and 3, and option 2 was relatively closer to consensus. While the discussion and the voting at that point should not be mistaken for a final recommendation or decision, the feedback was sufficient to prompt DEQ to explore further the potential ways that a regulatory system based on a defined TAC list could be developed.

## Goal Statement and Guiding Concepts

The following goal statement was proposed, for purposes of consideration and discussion, and was accepted by the ATW:

The TAC list includes the federal HAPs list and other air toxics that may be reasonably anticipated to occur in NSR permitted air emissions, and which warrant the evaluation of ambient air impacts in PTI applications in order to help ensure public health and environmental protection while promoting regulatory certainty and efficiency.

The following set of "guiding concepts" for developing an "option 2" approach was provided for discussion purposes:

1. The TAC list should include the HAPs list, and should additionally include the air toxics that may be reasonably anticipated to occur in emissions from facilities requiring a Permit to Install ( PTI ), minus those substances that have relatively low toxicity. The regulated community would prefer an approach that is focused on the more relevant substances, that is less burdensome and provides greater certainty.
2. The DEQ would have the authority to add to the list or remove substances from the list through the rulemaking process.
3. Rule 203(1)(c) should continue to require PTI applicants to describe the "quantity of all air contaminants that are reasonably anticipated due to the operation of the proposed process equipment." However, for unlisted air toxics (i.e., non-TACs), the current language in Rule 203(1)(h) would not be interpreted to be applicable; i.e., the applicant would not be required to provide in the PTI application, "Data demonstrating that the emissions from the process will not have an unacceptable air quality impact in relation to all federal, state, and local air quality standards." So, for non-TACs, the permit applicant would need to identify the emission rates but would not be required to model the ambient air impacts or compare the impacts to screening levels or other health protective benchmarks.
4. The DEQ rules should provide the DEQ authority to evaluate the ambient air impacts and potential health concerns of non-TACs in a PTI application, and to impose restrictions on their emissions as necessary to ensure public health protection. Section 324.5512 of NREPA authorizes the department to promulgate rules for controlling or prohibiting air pollution, and to deny or revoke a permit to operate a source, process, or process equipment that would adversely affect human health or other conditions important to the life of the community. [The Natural Resources and Environmental Protection Act (NREPA) Act 451 of the Public Acts of 1994, Part 55 Air Pollution Control].
5. For non-TACs, a modeled maximum ambient air impact exceeding a health-protective benchmark, such as a screening level (SL) as currently derived by the DEQ, may or may not in itself provide sufficient weight of evidence to support DEQ action to ensure public health protection under \#4 above. The DEQ may additionally consider relevant scientific and case-by-case information (as done currently under Rule 226(d) and Rule 228).

## Potential Approaches to List Development

In 2010, AQD conducted a survey of State's air toxics programs to gather basic information on the scope of their programs, including the list of air toxics regulated. The survey found that 29 of the 50 states regulate air toxics in permit reviews, based on ambient air impact estimates and public health protective benchmarks. Of the 21 states that do not routinely perform air toxics risk assessment in NSR, many (if not all) have a "backstop" or "safety net" provision for case-specific risk assessment. Of the six states in EPA Region 5 , four states routinely evaluate air toxics ambient air impacts for public health acceptability. Illinois generally does not (but could in exceptional cases). Indiana performs such evaluations only in a limited number of cases, not "routinely." Complete information was not collected on what list of air toxics are included for all states, but the gathered information did indicate that program scope varied widely. The state's approach for establishing the regulated air toxics may be generally grouped into five categories, as listed in Table 1 below.

Table 1. State's approaches to the development of lists of regulated air toxics.

| Air toxics included in NSR <br> health risk assessment | Example states | \# states |
| :--- | :--- | :--- |
| HAPs only | CT; HA; VA | 3 |
| HAPs plus additional air <br> toxics of concern | KY (HAPs+112r list); LA; NM <br> (HAPs+OELs); NY (HAPs+112r list); <br> NC; ND; RI; VT; WV (HAPs+OELs) | 10 |
| All air toxics with OELs | AL | 1 |
| State-specific list | OH; WI; CA; ID; MA; NH; SC | 7 |
| No discrete list; virtually any <br> may be included | MI; MN; DE; GA; MD; NJ; OK; TX | 9 |

Conceptually, there are several potential approaches to constructing a R225 TAC list, including the following:

1. Adopt a list developed by another state I states.
2. Develop a "list of lists."
3. List those chemicals meeting listing criteria based on health hazards, potency, persistence and bioaccumulation.
4. Develop a list based on the HAPs and the current list of TACs with SLs, with exclusion criteria.

The tendency for air toxics to pose a public health concern is generally a function of the potency, the exposure potential (which depends on the quantity and duration of the emission, the dispersion, and background exposures), and the presence and susceptibility of the public to the exposure. A list of regulated air toxics that is unlimited may be a relatively more reliable approach to address all potential concerns; any approach to developing a defined list of regulated air toxics may potentially be less reliable. For example, a substance with relatively low toxicity may be unlisted, however, a combination of high emissions, poor dispersion, and the presence of an exposed public, can pose public health concerns even if the toxicity or potency is relatively low. A "backstop" plan for detecting and addressing such cases is important, and is discussed elsewhere in this paper. Having noted this general limitation of any defined list, the following is a brief description of the apparent strengths and weaknesses/limitations of the four general approaches listed above, for discussion purposes.

## 1. Adopt a list developed by another state I states.

The positives of this option include convenience, and consistency (with the chosen State(s), but not with others). The concern is that the available lists in Region 5 may not be regarded by the DEQ, ATW, and/or the public, as fully appropriate for Michigan. The Ohio EPA list (303 compounds or classes) is based on the HAPs list plus substances passing several inclusion and exclusion criteria. Their rationale for applying exclusion criteria contains a considerable number of professional judgments. Some of these criteria may be regarded by some as having a questionable basis; environmental groups have strongly objected and have brought a lawsuit against Ohio EPA over the list and the criteria used to develop the list. The Minnesota MPCA has an unlimited list of regulated air toxics. The Wisconsin DNR's list was derived in 2004 based on certain inclusion and exclusion criteria,
and consists of 535 substances ( 26 HAPs are not included). Of course, lists from states outside of EPA R5 may also be considered. There is no consistency in the state's lists or in the approaches used to derive the lists. It would be arguable to debate whose list is more appropriate for Michigan.

## 2. Develop a "list of lists."

This approach was recommended by the Michigan Air Toxics Policy Committee (1989) as a way to focus the required environmental acceptability assessments (with case-by-case assessment of other air toxics of concern at a specific site). They recommended a list of approximately 1200 substances, consisting of the substances with ACGIH or NIOSH OELs, the Michigan Critical Materials Register, the NTP and IARC lists of carcinogens, and the chemicals listed in the IJC's Great Lakes Water Quality Board 1987 Report on Great Lakes Water Quality. As noted in Table 1 above, some states have used the EPA's 112(r) chemical list for emergency preparedness (which consists of 77 acutely toxic chemicals, and 63 flammable gases and volatile flammable liquids). Another relevant list available today is the EPA's Toxics Release Inventory (TRI) list.

The strengths of this approach are the relative ease of compiling a list of lists, and, the contributing lists would presumably have some environmental relevance. The limitations of this approach are that many listed substances may be irrelevant to PTI air emissions in Michigan, and, many of the substances on lists such as the TRI may have inadequate data for SL development. Also, this approach can result in a very long list, which may be undesirable to the regulated community (guiding concept \#1 above).

## 3. List those chemicals meeting listing criteria based on health hazards, potency, persistence and bioaccumulation.

The strength of this approach is that the scientific defensibility may be relatively strong. The limitations of this approach are that it is a relatively labor intensive and time consuming initiative, the appropriate criteria may be difficult to establish, and the resulting list may not be the most relevant to the PTI program. Also, this approach (a version of which was implemented by Ohio EPA) may rely on multiple judgments for inclusion or exclusion that may be contested. A key element would be to establish well-reasoned, nonarbitrary inclusion and exclusion criteria, preferably derived by a consensus approach among multiple stakeholders.

## 4. Develop a TAC list based on the HAPs and the current MDEQ list of TACs with SLs, with exclusion criteria.

The strengths of this approach are relative efficiency of list development, the focus on air toxics that are relevant to PTI applications in Michigan, and the inclusion of those substances that have already been found to have sufficient toxicity data for SL development. As with \#3 above, a key element would be to derive well-reasoned, nonarbitrary criteria, but in this case, those would be more limited since they would only be exclusion criteria (i.e., criteria for not including certain substances that currently have SLs). The limitation of this approach is that the selection of the exclusion criteria may be debatable.

Further rationale for approach \#4: The initial universe of substances for assessment is the current SL list of 1202 substances (as of May, 2013). This list represents MDEQ's 21+ years of experience in evaluating air toxics in the New Source Review permitting program, under an open-ended TAC definition (excluding only a short list of exempted substances; currently 41). Over the last 21 years (since 1992), screening levels have been derived for TACs (under the open-ended definition) if they appeared in proposed emission characterizations for all categories of facilities (thermal, chemical, or general manufacturing). Data-poor chemicals were addressed relatively inclusively in the MDEQ program, i.e., SL derivation methods include the use of minimal data such as subchronic animal studies, LD50s, and LC50s. This list also includes 289 substances with inadequate toxicity data for SL derivation, which were assigned the default ITSL of $0.1 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual AT). Rather than propose the inclusion of all 1200+ substances on the future TAC list, some exclusion criteria may be reasonable in the interest of developing a shorter list that is more focused on the more relevant substances and is less burdensome on the regulated community (guiding concept \#1).

## Proposal for the TAC List

It was proposed that the MDEQ follow approach \#4 above, to develop a defined TAC list including the following:

1. Most EPA HAPs should be included, including all individual chemicals that EPA includes as members of HAP listed groups (e.g., metal compounds). For clarity, the individual chemical members of the HAP groups of polycyclic organic matter (POM) and glycol ethers should be listed individually and only if they meet the other qualifying criteria (based on the ITSL or carcinogenicity). The HAPs list includes many air toxics with well documented toxicity and with the potential for public exposure, based on air emissions data and/or ambient air monitoring data. The HAPs list is the focus of EPA's air toxics data collection and regulatory actions under the Clean Air Act. Ohio EPA adopted all HAPs into their Toxic Air Pollutant list. However, it may be noted that some of the HAPs have relatively limited toxicity datasets, and some of the HAPs have not been identified and addressed in Permit to Install applications. For some HAPs, it may not be reasonable to anticipate that they would appear in future PTI applications. Reasons to include all HAPs in a TAC list are: for simplicity; for consistency with EPA; and, for better clarity in communicating the basis for the list with the regulated community and other groups. Reasons to not include some HAPs in the TAC list are: to better focus on the air toxics most relevant to PTI applications; and, many HAPs do not have SLs and therefore may never have been identified in a PTI application. In some cases, DEQ has evaluated air toxics in PTI applications and not established a SL, but rather notified permits staff that the predicted ambient air impact is acceptable, in cases where the impact was very low and the toxicologist did not feel it was appropriate to establish a data-derived or default SL. Therefore, for the Table 2 list of HAPs without SLs, the Toxics Unit files were reviewed to determine if the substance had been evaluated for a PTI application (Table 2 has a column for "File Review Comments"). It is tentatively proposed that the potential TAC list exclude HAPs that do not have a SL and have not been encountered in a PTI application.
2. All carcinogens would be included (i.e., all compounds with a current IRSL, or, meeting the current rules' definition of a carcinogen (e.g., asphalt fumes)). See also the discussion of the carcinogenic PAHs in Table 5.
3. All substances with ITSLs at or below a cutoff value would be included; substances with only ITSLs that are above the cutoff values would be excluded (see discussion below).
4. It may be considered to exclude all substances with an ITSL of $0.1 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual averaging time) based on the default value and a lack of chemical-specific data sufficient for SL development. That would include 287 chemicals currently on the SL list. This approach is consistent with Guiding Concepts \#1 described earlier. This approach would also be consistent with the other EPA R5 State air toxics programs. It may be noted that Texas TCEQ utilizes a default effect screening level (ESL) of $2 \mu \mathrm{~g} / \mathrm{m}^{3}$ ( 1 hour averaging time) when data are lacking for ESL derivation. That default ESL is similar to the AQD default ITSL, using the EPA's Screen3 averaging time (AT) conversion factor of 0.08 for converting from 1 hour AT to annual AT $\left(2 \mu \mathrm{~g} / \mathrm{m}^{3}(1 \mathrm{hr}\right.$ AT $) \times 0.08=0.16 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual AT)).
5. Consistent with the Guiding Concepts described earlier, substances not on the TAC list would be identified in PTI applications, including information on the quantity of emissions (R203(1)(c)), but the applicant would not be required to include further information demonstrating the acceptability of the air quality impacts. MDEQ may still address those substances, with justification, by way of emission limits to protect the public health and/or adding substances to the TAC list via rulemaking.

## ITSL Cutoff Values

Criterion \#3 above mentions ITSL cutoff values. While initially proposed cutoff values for consideration may be largely arbitrary (e.g., proposing a $\mu \mathrm{g} / \mathrm{m}^{3}$ value or a percentile of an ITSL distribution), the final selection of an appropriate and reasonable cutoff is not arbitrarily selected. Careful consideration by staff and the ATW Members of the reasonableness of the approach, the magnitude of the resulting ITSL cutoff values, the resulting chemicals that meet or fail to meet the cutoff values, and the overall adequacy of the TAC list to meet the goal and the guiding concepts, followed by an ATW recommendation, make the approach more reasoned and deliberate.

The selection of a cutoff may take into consideration available and appropriate criteria utilized in other air quality protection activities. For example, for substances that may be anticipated to exist as particulates in air emissions and in ambient air, consider the primary NAAQS for particulate matter ( $150 \mu \mathrm{~g} / \mathrm{m}^{3}$ ( 24 hour) for $\mathrm{PM}_{10}$, and $12 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual) and $35 \mu \mathrm{~g} / \mathrm{m}^{3}$ (24 hour) for $\mathrm{PM}_{2.5}$ ); also consider that the ACGIH (2012 handbook; Appendix B) recommends TLVs of $3 \mathrm{mg} / \mathrm{m}^{3}$ (respirable particles) and $10 \mathrm{mg} / \mathrm{m}^{3}$ (inhalable particles) for Particles Not Otherwise Specified (PNOS).

The Wisconsin air toxics regulatory list is based on several qualifying criteria, including exclusion criteria of having an OEL (TLV) of greater than or equal to 100 ppm or $10 \mathrm{mg} / \mathrm{m}^{3}$.

A TLV of $10 \mathrm{mg} / \mathrm{m}^{3}$ would be associated with an AQD ITSL of $100 \mu \mathrm{~g} / \mathrm{m}^{3}(8 \mathrm{hr} \mathrm{AT})$ (utilizing an uncertainty factor of 100, as per the air toxics rules).

It may be considered that the EPA has de-listed some HAPs based upon a finding that there are adequate data on the health and environmental effects of these substances to determine that emissions may not reasonably be anticipated to cause adverse human health or environmental effects (Table 3).

The establishment of a cutoff may also consider the range of ITSL values thus far derived by DEQ. An assessment of the current SL values, and the selection of a reasonable percentile of the distribution of the current ITSLs, may help distinguish the relatively more toxic substances (in the majority of the distribution) from the relatively lower toxicity substances (in the minority of the distribution). Setting that cutoff may be guided by consideration of the range of current ITSL values. Rather than setting an a priori percentile of the distribution as the cutoff point, it was considered informative to describe the distribution (e.g., the $50^{\text {th }}, 75^{\text {th }}, 90^{\text {th }}, 95^{\text {th }}$ and $99^{\text {th }}$ percentiles). The distributions were determined after excluding from the dataset those substances with an ITSL of $0.1 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual AT) based on the default value. These percentiles were first determined for all current ITSLs, without distinction as to HAP or non-HAP status, and without regard to the various averaging times (ATs) associated with the screening levels. For substances with two ITSLs (acute and chronic), only the chronic (lower) ITSL was included in the assessment. The ITSL distributions were also determined for the following subsets: HAPs only; non-HAPs only; annual AT only; 24 AT only; 8 hr AT only; and, 1 hr AT only. The resulting summary statistics for the ITSL group datasets, as of May 2013, that were initially considered by the ATW are presented in Table 4a. It should be noted that an August update of the $75^{\text {th }}$ percentile values is presented in Table 4b.

For discussion purposes, staff initially pursued the potential content of a TAC list that includes the current ITSLs except for those exceeding the $75^{\text {th }}$ percentile cutoff point for each specific averaging time, in addition to the other listing criteria previously mentioned (in bold in Table 4a). This approach and proposed cutoff points were regarded by staff as reasonably inclusive, while providing a significant reduction in the current SL list (guiding concept \#1). Following ATW consideration and discussion at several meetings through the $8^{\text {th }}$ meeting on August $1^{\text {st }}, 2013$, the utilization of the $75^{\text {th }}$ percentile of the distribution for each ITSL averaging time appeared to gain acceptance by many Members, pending a final Workgroup recommendation. It should be noted that the updated ITSL cutoff values appear in Table 4b and Table $\mathbf{7}$ and in the document on the ATW website, "Proposed TAC List, August Update".

## Authority to Address Unlisted Air Toxics in PTI Applications

If the current TAC definition were to be changed to some defined list, then a key issue would be the DEQ's authority to address air toxics concerns that may arise for unlisted air toxics that are proposed for emission in a PTI application. A review of the authority of other state's air agencies, and of other MDEQ divisions, to address unlisted substances, is summarized in Table 6. It was proposed for discussion purposes that AQD adopt rule
language similar to that of MDEQ-WRD in Table 6. Following Workgroup discussion of issue A-1(9) regarding Rule 228, the Workgroup drafted a recommendation to retain Rule 228 with the addition of clarifying language, and a Member proposed that non-TACs could also be addressed by the AQD as appropriate under this authority.

## ITSLs With 1 Hour Averaging Times

Upon review of the proposed TAC list and ITSL cutoffs (Table 4a), it was noted that the $75^{\text {th }} \%$ ile cutoff value for the 1 hr AT ITSLs $\left(300 \mu \mathrm{~g} / \mathrm{m}^{3}\right)$ was not as high as for the 8 hr or 24 hr ITSLs. Staff responded that this group presumably has a relatively lower ITSL distribution because it includes a relatively more acutely toxic subset of the substances that have TLV occupational exposure levels. A Member asked staff to evaluate the chemicals with 1 hr AT ITSLs that do not meet the criteria for TAC listing; if they raise concerns, then it may be an option to include them in the TAC list. Staff evaluated this list of 33 chemicals; eight have 1 hr AT ITSLs above the $75^{\text {th }} \%$ ile value of $300 \mu \mathrm{~g} / \mathrm{m}^{3}$. Of these eight, one (methylene chloride) is a carcinogen and therefore will be on the TAC list. Another (hydrogen chloride) will be on the TAC list because it also has an annual AT ITSL $\left(20 \mu \mathrm{~g} / \mathrm{m}^{3}\right)$ that is below the $75^{\text {th }} \%$ ile cutoff for the annual AT. Staff do not feel that the remaining six raise particular concerns for being unlisted, therefore, it is proposed to not make an exception to the $75^{\text {th }}$ percentile cutoff for these chemicals:

| Chemical | CAS \# | $\mathbf{1} \mathbf{~ h r ~ A T ~ I T S L ~}$ <br> $\left(\mu \mathrm{g} / \mathbf{m}^{\mathbf{3}}\right)$ | Other ITSL |
| :--- | :--- | :--- | :--- |
| Ethylene glycol | $107-21-1$ | 1000 |  |
| Hexylene glycol | $107-41-5$ | 1210 |  |
| Methanol | $67-56-1$ | 3250 | $2700 \mu \mathrm{~g} / \mathrm{m}^{3}(8 \mathrm{hr}$ AT $) ;$ <br> this is above the $75^{\text {th }}$ <br> $\%$ \%ile. |
| Isoamyl acetate | $123-92-2$ | 5300 | 130000 (annual AT); <br> this is above the $75^{\text {th }}$ <br> $\%$ \%ile cutoff. |
| Trichlorofluoromethane | $75-69-4$ | 56200 | 5560000 |
| Hfc-227ea | $431-89-0$ |  |  |

## Listing of Chemical Groups

The Workgroup discussed how the EPA HAPs list contains chemical groups for metals, and also for glycol ethers, cyanide compounds, POM (polycyclic organic matter), etc. The listing of chemical groups gives the impression of a smaller list size. There are 187 HAPs including the chemical groups, but the actual size of the list of specific HAP chemicals is much larger. The inclusion of chemical groups in a regulatory list can enable a regulatory agency to add chemicals to the list (as new members of a listed group) very efficiently, but
this diminishes the goal of a list to be clear and as specific as possible. The Workgroup favored the clarity of specific chemical listings rather than the use of some of the groups as in EPA's HAPs list, although it is recognized that this contributes to a longer list than if groups were listed. Therefore, the proposed list includes specific PAHs and glycol ether compounds, etc., if they meet the criteria for listing. Regarding metal compounds, staff feels that in some cases these compounds should be listed separately, because toxicity (and the magnitude of the health protective screening level) is dependent on the specific metal compound. However, in other cases, different compounds of the same metal have toxicity that is primarily determined by the metal alone. In these cases, it seems inappropriate to list the metal forms individually, and then apply a footnote directing that their emissions and impacts should be evaluated additively (with adjustment of the MW to the atomic weight of the metal) for comparison to the screening level. Therefore, staff anticipates that some metals may be appropriately listed as a TAC group. The current SL list, and draft proposed TAC list, include some specific metal compounds that may be grouped together in the future, pending further review. For example, an initial review has tentatively identified the following cases where further assessment is warranted:
"Antimony and antimony compounds" may consolidate 5 current listings.
"Cobalt and cobalt compounds" may consolidate 3 current listings.
"Copper and copper compounds" may consolidate 4 current listings.
"Magnesium and magnesium compounds" may consolidate 7 current listings.
"Manganese and manganese compounds" may consolidate 4 current listings.
"Molybdenum water soluble compounds" may consolidate 3 current listings.
"Molybdenum water insoluble compounds" may consolidate 3 current listings.

## Merging of the Current Annual AT ITSLs With the Current RfC- and RfD-Based 24Hour AT ITSLs That Are Anticipated To Change to Annual AT

Concurrent with addressing the TAC list issue, the Workgroup explored the ORR Report's Recommendation A-1(7): Make acceptable exposure limits consistent with other nearby states. As a result of that discussion, the Workgroup is recommending that AQD utilize a default annual averaging time (AT) rather than a 24 hour AT for ITSLs that are based on the EPA RfC and RfD methodologies. AQD is agreeable to making that change. Therefore, for those chemicals, the change in AT from 24 hours to annual may be regarded as "impending". However, this issue crosses over to the "TAC List" issue, because the proposed TAC list criteria include ITSL cutoff values set at the $75^{\text {th }}$ percentile level for each AT. Those $75^{\text {th }}$ percentile values are statistically determined based on the distribution of all of the non-default ITSLs for each AT. Previous estimates (e.g., the April 2013 statistics in Table 4A) of the $75^{\text {th }}$ percentiles, TAC list size, and the TAC list of chemicals were based on the current ATs and $75^{\text {th }}$ percentiles, and did not account for this
impending change in ATs. Further, the proposed draft rule language for the TAC list issue will include specific ITSL cutoff values. Therefore, it seems appropriate and necessary to address this impending change in the ATs so that the specific ITSL cutoff values in the draft proposed rules will reflect the AT change. In other words, the ITSL cutoff values for both annual and 24 hour ATs in the proposed draft rules should reflect that impending change. Also, there was a concern that making that change could significantly change the $75^{\text {th }}$ percentile cutoff values, and potentially cause a significant change in the number of chemicals proposed for the TAC list.

Staff recognized this issue and completed the evaluation of this AT change after the August 1, 2013 ATW meeting. After all ITSLs with a current 24 hour AT based on the EPA RfC or RfD methodologies are changed to annual AT, only eight chemicals will still have a 24 hour AT ITSL. The characteristics of that group are described in Table 7. The previous $75^{\text {th }}$ percentile cutoff values and the number of chemicals in the proposed draft TAC list are also presented for comparison in Table 7. Although the AT conversion results in a relatively small set of chemicals ( $n=8$ ) that will have 24 hr AT ITSLs, the ITSLs in that group are well distributed (ranging from $2 \mu \mathrm{~g} / \mathrm{m}^{3}$ to $10000 \mu \mathrm{~g} / \mathrm{m}^{3}$ ), and the $75^{\text {th }} \%$ ile cutoff did not change greatly (an increase from 420 to $522 \mu \mathrm{~g} / \mathrm{m}^{3}$ ).

Based on these findings, it is proposed that the draft TAC list rules utilize the cutoffs that result from the conversion of the ITSL ATs as described above. The effect of merging the two groups (those with current annual AT ITSLs, and those with an impending AT change from 24 hours to annual AT) is an increase in the cutoff from $43 \mu \mathrm{~g} / \mathrm{m}^{3}$ to $100 \mu \mathrm{~g} / \mathrm{m}^{3}$ for the annual AT. The effect of this change is the inclusion of chemicals that currently have annual AT ITSLs that are above the prior annual AT cutoff of $43 \mu \mathrm{~g} / \mathrm{m}^{3}$, but which are at or below the new cutoff of $100 \mu \mathrm{~g} / \mathrm{m}^{3}$. Another effect of this change is the exclusion of chemicals that have current 24 hr AT ITSLs below the prior 24 hr AT cutoff of $420 \mu \mathrm{~g} / \mathrm{m}^{3}$, but above the new annual AT cutoff of $100 \mu \mathrm{~g} / \mathrm{m}^{3}$. The overall net effect of these changes is a small increase in the total list of TACs (a change from 750 to 756 chemicals). This is further described in Table 7. The graph below helps to visualize the distribution of the merged annual AT ITSLs, and the $75^{\text {th }}$ percentile cutoff value.


## Other Chemical Listing Discussions

In addition to the above criteria, procedures, and discussions, the ATW discussed the listing of two perfluorinated compounds (PFOS, CAS\# 1763-23-1; and PFOA, CAS\# 335-67-1), crystalline silica (from sources not meeting the current TAC list exemption; CAS\# 14808-60-7), carcinogenic PAHs, and asphalt fumes (CAS\# 8052-42-4) (see Table 5). Also, a Member requested that methyl isocyanate (CAS\# 624-83-9) be added to the proposed TAC list due to high toxicity and the potential that it could occur in a future permit application. Although mercury (CAS\# 7439-97-6) does not have a SL, the SL list has a footnote indicating that a benchmark for inhalation of elemental mercury $\left(0.3 \mu \mathrm{~g} / \mathrm{m}^{3}\right)$ would meet the cutoff criterion; mercury is included in the future TAC list.

## Proposed TAC List and Procedure

The proposed TAC list, based on the above criteria, procedure, and discussions, is 756 chemicals. This may be anticipated to change somewhat due to the routine updating of chemical risk assessments, the evaluation of "new" air toxics in permit applications, the potential consolidation of some metal compounds, etc. Further statistical information and a spreadsheet showing all current TACs, and the basis for chemicals meeting or not meeting the criteria for the proposed future TAC list, are available on the ATW website in an August 13, 2013 document, "Proposed TAC List, August 2013". The spreadsheet
includes a notation for the chemicals that currently have 24 hr AT ITSLs but with an impending change to an annual AT. The spreadsheet reflects the updated $75^{\text {th }} \%$ ile cutoff values as listed in Table 4b and Table 7.

Table 2. HAPs without SLs.

| Chemical and CAS \# | Toxics Unit File Review Comments |
| :---: | :---: |
| Acetamide 60-35-5 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 2-acetylaminofluorene 53-96-3 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 4-aminobiphenyl 92-67-1 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| o-anisidine $90-04-0$ | O-anisidine hydrochloride (134-29-2) has an IRSL. Therefore, include it in the TAC list. |
| Benzotrichloride (trichlorotoluene) 98-07-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Calcium cyanamide 156-62-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Captan 133-06-2 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Carbaryl $63-25-2$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Catechol $120-80-9$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Chloramben 133-90-4 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Chlordane 57-74-9 | Chlordane (technical) (12789-03-6) has an ITSL and IRSL. Therefore, it is proposed to include it in the TAC list. |
| Chloroacetic acid 79-11-8 | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| Chlorobenzilate 510-15-6 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Chloromethyl methyl ether 107-30-2 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 2,4-D, salts and esters 94-75-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| $\begin{aligned} & \text { DDE } \\ & 3547-04-4 \end{aligned}$ | DDD(TDE; 72-54-8), DDE(p,p'; 72-55-9) and DDT(50-29-3) have IRSLs. Therefore, it is proposed to include it in the TAC list. |
| Diazomethane 334-88-3 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 3,3-dimethoxybenzidine 119-90-4 | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| Dimethyl aminoazobenzene 60-11-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 3,3'-dimethyl benzidine 119-93-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Dimethyl carbamoyl chloride | There is no indication of a review for NSR permitting, |


| 79-44-7 | therefore, it is proposed to not include it in the TAC list. |
| :---: | :---: |
| 1,1-dimethyl hydrazine 57-14-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 1,2-diphenylhydrazine 122-66-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Ethyl carbamate (Urethane) 51-79-6 | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| $\begin{array}{\|l\|} \hline \text { Ethylene imine (Aziridine) } \\ 151-56-4 \\ \hline \end{array}$ | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| Hexamethylphosphoramide 680-31-9 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Hydroquinone 123-31-9 | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| Lindane (all isomers) 58-89-9 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Methoxychlor 72-43-5 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Methyl iodide (Iodomethane) $74-88-4$ | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| Methyl isocyanate 624-83-9 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 4,4-methylene bis(2chloroaniline) 101-14-4 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 4,4'-methylenedianiline 101-77-9 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| 4-nitrobiphenyl 92-93-3 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| N-Nitrosomorpholine 59-89-2 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| $\begin{aligned} & \text { Parathion } \\ & 56-38-2 \\ & \hline \end{aligned}$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| p-Phenylenediamine 106-50-3 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Phthalic anhydride 85-44-9 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| beta-Propiolactone $57-57-8$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| $\begin{aligned} & \text { Propoxur (Baygon) } \\ & \text { 114-26-1 } \end{aligned}$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Quinone (p-benzoquinone) 106-51-4 | This was evaluated for at least one NSR permit. Therefore, it is proposed to list it as a TAC. |
| Styrene oxide 96-09-3 | Styrene (also a HAP) has an IRSL. Styrene is metabolized to styrene oxide. Both are reasonably anticipated to be human carcinogens (NTP Report on Carcinogens, $12^{\text {th }}$ Ed.). Therefore, RETAIN on TAC list. |


| $\begin{array}{\|l} \hline \text { Titanium tetrachloride } \\ 7550-45-0 \end{array}$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| :---: | :---: |
| 2,4-toluene diamine 95-80-7 | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| $\begin{aligned} & \text { Trifluralin } \\ & 1582-09-8 \end{aligned}$ | There is no indication of a review for NSR permitting, therefore, it is proposed to not include it in the TAC list. |
| Lead compounds | Lead is a criteria pollutant; exempted from TAC defn. |
| Radionuclides (including radon) | A 1994 DEQ policy determination was that there were sufficient regulations by NRC, EPA, and MDCH, such that additional AQD permitting requirements would be unnecessary and duplicative. |
| Polycyclic organic matter (POM) | The TAC list should include specific compounds, for clarity, if they meet criteria (ITSLs or carcinogenicity). |
| Glycol ethers | The TAC list should include specific compounds, for clarity, if they meet criteria (ITSLs). |

Table 3. De-listed EPA HAPs.

| Delisted HAP | Date of delisting | AQD ITSL ( $\mu \mathrm{g} / \mathrm{m}^{3}$; AT) or RfC | comments |
| :---: | :---: | :---: | :---: |
| Caprolactam | 6/18/96 | $10 \mu \mathrm{~g} / \mathrm{m}^{3}$ ( 8 hr AT ) |  |
| Surfactant alcohol ethoxylates and their derivatives (SAED) (in glycol ethers HAP category) | 8/2/2000 | Ethylene glycol ether <br> 2-methoxy-1propanol (a nonSAED) used as a conservative surrogate to derive an RfC-like benchmark of 200 to $2000 \mu \mathrm{~g} / \mathrm{m}^{3}$ for SAEDs. | A hypothetical facility emission rate of 105 lbs total SAEDs/year was used in the petition for de-listing, and was relied upon in EPA's review. |
| Ethylene glycol monobutyl ether (2butoxyethanol) (in glycol ethers HAP category) | 11/29/04 | $\begin{aligned} & 1600 \mu \mathrm{~g} / \mathrm{m}^{3}(24 \mathrm{hr} \\ & \text { AT) } \end{aligned}$ |  |
| Methyl ethyl ketone | 12/19/05 | $\begin{aligned} & 5000 \mu \mathrm{~g} / \mathrm{m}^{3}(24 \mathrm{hr} \\ & \text { AT) } \end{aligned}$ |  |

Table 4a. ITSL value distribution (as of April, 2013). All values are in units of $\mu \mathrm{g} / \mathrm{m}^{3}$. (These statistics are based on only the air toxics with data-derived final SLs, i.e., excluding chemicals with only default-based ITSLs).

| ITSL <br> group | Mean | $\mathbf{5 0}^{\text {th }}$ \%ile | $\mathbf{7 5}^{\text {th }}$ \%ile | $\mathbf{9 0}^{\text {th }}$ \%ile | $\mathbf{9 5}^{\text {th }}$ \%ile | $\mathbf{9 9}^{\text {th }}$ \%ile |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| All ITSLs | 1375 | 24 | 140 | 1956 | 5000 | 23800 |
| HAPs only | 626 | 14.5 | 100 | 1000 | 3088 | 13572 |
| Non-HAPs <br> only | 1547 | 28 | 140 | 2300 | 5450 | 42850 |
| Annual AT <br> only | 482 | 14 | $\mathbf{4 3}$ | 140 | 300 | 1363 |
| 24 hr AT <br> only | 1789 | 60 | $\mathbf{4 2 0}$ | 2600 | 6000 | 46600 |
| 8 hr AT <br> only | 2760 | 86 | $\mathbf{2 8 5 0}$ | 6020 | 16710 | 30482 |
| 1 hr AT <br> only | 2741 | 15 | $\mathbf{2 9 0}$ | 1168 | 3046 | 44551 |

Table 4b. Updated $75^{\text {th }}$ percentile values (as of August, 2013) reflecting the change in averaging time from 24 hours to annual for ITSLs based on the RfC or RfD methodologies. All values are in units of $\mu \mathrm{g} / \mathrm{m}^{3}$. (These statistics are based on only the air toxics with dataderived final SLs, i.e., excluding chemicals with only default-based ITSLs).

| Averaging Time | $\mathbf{7 5}^{\text {th }}$ Percentile of Distribution <br> $\left(\mu \mathrm{g} / \mathrm{m}^{3}\right)$ |
| :--- | :--- |
| 1 hr | $\mathbf{3 0 0}$ |
| 8 hr | $\mathbf{2 3 3 0}$ |
| 24 hr | $\mathbf{5 2 2}$ |
| Annual | $\mathbf{1 0 0}$ |

Table 5. Additional air toxics ( $\mathrm{n}=23$ ) that are not on the TAC SL list, which are proposed to be added to the future TAC list:

| Substance | Comments on why there is no SL, but that listing as a TAC would be appropriate |
| :---: | :---: |
| Crystalline silica (14808-60-7) | Not a HAP. Some sources of crystalline silica are exempt from TAC definition. (AQD has recently set an ITSL at $3 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual AT)). Proposed to place it on the TAC list. The current TAC list exemption for certain sources would remain. |
| Asphalt fumes (8052-42-4) | Not a HAP as a mixture. The fumes contain carcinogens, but there is no IRSL for the mixture due to lack of a key study on the mixture. Based on a 1995 Scientific Advisory Panel recommendation, AQD has regulated the mixture utilizing the EPA RPFs for carcinogenic PAHs (see also below). Proposed to list this mixture as a TAC with an explanatory footnote (only) that would help clarify the regulatory approach. |
| Carcinogenic PAHs ( $\mathrm{n}=19$, in addition to those meeting other listing criteria) | The PAHs are HAPs as "POM." The EPA's risk assessment of the carcinogenic PAH group is currently in transition. The 1993 EPA guidance for the group is currently still in use by MDEQ (there are 7 carcinogenic $P A H s$, including $B(a) P$ and 6 with Relative Potency Factors (RPFs) relative to $\mathrm{B}(\mathrm{a}) \mathrm{P}$ ). CalOEHHA regulates 21 carcinogenic PAHs with RPFs. EPA has drafted a new scheme, with 25 carcinogenic PAHs with nonzero RPFs (including B(a)P); they are currently addressing the SAB review comments on that draft <br> (http://yosemite.epa.gov/sab/sabproduct.nsf/0/E65D909C98520C1 D85257501005E46AE?OpenDocument). Currently, 16 do not have SLs. Three additional PAHs have evidence of carcinogenicity, have CalOEHHA RPFs, and are not on the current SL list. Therefore, 19 additional substances for the TAC list are proposed, for this group. (In the $5 / 13 / 13$ spreadsheet of potential TACs, the basis for listing = "Carc7" ( $n=7$ ), "EPA Carc" ( $n=16$ ), or "CAL Carc" ( $n=3$ ). |
| Perfluorinated compounds (PFCs): PFOS and PFOA ( $\mathrm{n}=2$ ) | Perfluorooctane sulfonate (PFOS) and perfluorooctanoic acid (PFOA) are persistent bioaccumulative toxics (PBTs) that have been identified by MDEQ as emerging contaminants of concern. (http://www.michigan.gov/deq/0,4561,7-135-3308-266777-,00.html). PFCs have recently been detected in Michigan groundwater and in several species of aquatic and terrestrial wildlife. Although the presence of PFCs in air emission sources subject to NSR permitting has not yet been characterized, it is proposed that these two PFCs be listed as TACs. (In the 8/13/13 spreadsheet of potential TACs, the basis for listing = "Emerging". |

Table 6. Authority to address unlisted substances.

| Agency | Description of authority |
| :--- | :--- |
| MDEQ-Water Resources <br> Division (WRD) | NREPA Part 8 rules regulate surface water discharges of "toxic <br> substances," which are defined as those included in three lists of <br> substances (several hundred) and, "Any other toxic substances <br> that the department determines are of concern at a specific site." |
| MDEQ-Remediation and <br> Redevelopment Division <br> (RRD) | NREPA Part 201 rules define "hazardous substance" as three lists of <br> substances (several hundred), and, "Any substance that the <br> department demonstrates, on a case by case basis, poses an <br> unacceptable risk to the public health, safety, or welfare, or the <br> environment, considering the fate of the material, dose- <br> response, toxicity, or adverse impact on natural resources." |
| Ohio EPA - Air | Ohio EPA has a list of 303 chemicals/classes of regulated air toxics. <br> Language in administrative code and in rules gives authority for their <br> Director to evaluate unlisted air toxics (personal communication with <br> Paul Koval, 2/21/13). |
| Wisconsin DNR - Air | There are 535 listed "hazardous air contaminants" <br> substances/groups; this was established in 2004, based on criteria <br> specified in their code. Authority to address unlisted substances: <br> "Code: NR 445.03 General limitations. No person may cause, <br> allow or permit emissions into the ambient air of any hazardous <br> substance in a quantity or concentration or for a duration that is <br> injurious to human health, plant or animal life unless the <br> purpose of that emission is for the control of plant or animal life. <br> Hazardous substances include but are not limited to the <br> hazardous air contaminants listed in Tables A to C of s. NR <br> 445.07." |
| Minnesota PCA - Air | MN does not have a defined list of regulated air toxics. Statute: "The <br> Pollution Control Agency may issue, continue in effect or deny <br> permits, under such conditions as it may prescribe for the <br> prevention of pollution, for the emission of air contaminants..." |

Table 7. The Effects of Converting the 24 Hour AT ITSLs Based on the RfD or RfC Methodologies to Annual AT ITSLs.
\(\left.$$
\begin{array}{|l|l|l|l|}\hline & \begin{array}{l}\text { May 13, 2013 } \\
\text { Draft } \\
\text { Discussion } \\
\text { Paper }\end{array} & \begin{array}{l}\text { Current } \\
\text { Discussion } \\
\text { Paper }\end{array} & \text { Comments } \\
\hline \begin{array}{l}\text { Number of } \\
\text { chemicals with } \\
\text { an annual AT } \\
\text { ITSL }\end{array} & 389 & 620 & \begin{array}{l}\text { The current number reflects the } \\
\text { conversion from 24 hr AT to } \\
\text { annual AT for all RfC- and RfD- } \\
\text { based ITSLs. }\end{array} \\
\hline \begin{array}{l}\text { Number of } \\
\text { chemicals with } \\
\text { a } 24 \text { hr AT ITSL }\end{array} & 239 & 8 & \begin{array}{l}\text { Same as above. }\end{array} \\
\hline \begin{array}{l}75^{\text {th } \% \text { ile cutoff }} \\
\text { for annual AT } \\
\left(\mu \mathrm{g} / \mathrm{m}^{3}\right)\end{array} & 43 & 100 & \begin{array}{l}\text { The current cutoff is significantly } \\
\text { higher than previous, due to the } \\
\text { new, larger group of chemicals } \\
\text { in the annual AT group. }\end{array} \\
\hline \begin{array}{l}75^{\text {th } \% \text { \%ile cutoff }} \\
\text { for } 24 \text { hr AT } \\
\left(\mu \mathrm{g} / \mathrm{m}^{3}\right)\end{array} & 420 & 522 & \begin{array}{l}\text { The AT conversion will result in } \\
8 \text { remaining chemicals with a } 24 \\
\text { hr AT. Only two of these } 8 \\
\text { chemicals (TCE and }\end{array}
$$ <br>

tetrachloroethylene) have 24 hr\end{array}\right\}\)| AT ITSLs that are above the |
| :--- |
| cutoff of 522 $\mu g / m^{3}$; they would |
| be listed as TACs based on |
| carcinogenicity. |

## APPENDIX J:

## Potential Defined TAC List

## APPENDIX J: Proposed TAC List - August Update

| Table 1. Percentiles of Initial Threshold Screening Levels |  |  |  |
| :---: | :---: | :---: | :---: |
| Initial Threshold Screening Levels (ITSLs) Grouped by Averaging Time | 75th <br> Percentile of ITSL Group | $\begin{gathered} \text { Count of 1st ITSL } \\ <75 \text { th\% } \\ \hline \end{gathered}$ | $\begin{gathered} \text { Count of 2nd ITSL } \\ <75 \text { th\% } \\ \hline \end{gathered}$ |
| 1 hr | $300 \mu \mathrm{~g} / \mathrm{m}^{3}$ | 18 | 7 |
| 8 hr | 2330 g/ $\mathrm{m}^{3}$ | 135 | 10 |
| 24 hr | $522 \mu \mathrm{~g} / \mathrm{m}^{3}$ | 2 | 4 |
| Annual | $100 \mu \mathrm{~g} / \mathrm{m}^{3}$ | 479 | 0 |
| Note: Some Toxic Air Contaminant (TACs) have 2 ITSLs each with different averaging times. One or both ITSLs may be less than 75th percentile cutoff. |  |  |  |

Table 2. Basis for TACs on the Future Screening Level List

| Count only if "1st ITSL" | 578 |
| ---: | :---: |
| Count only if "1st ITSL, 2nd ITSL" | 18 |
| Count only if "1st ITSL, 2nd ITSL, Carc*" | 3 |
| Count only if "1st ITSL, Carc" | 40 |
| Count only if "Carc" | 81 |
| $* *$ Count only if "Added" | 36 |
| Total Number of Future TACs | 756 |

* "Carc" = carcinogenic compounds. All carcinogenic TACs have Initial Risk Screening Levels (IRSLs), except Asphalt Fumes.
**Note: Asphalt fumes and crystalline silica were listed as "carc" and "1st ITSL", respectively, therefore, are counted in those groups above, despite being technically "Added".

| Table 3. Other Information |  |
| :---: | :---: |
| Total number of compounds evaluated | 1231 |
| 1st ITSL < 75\% | 639 |
| 2nd ITSL < $75 \%$ | 21 |
| *TACs removed from List | 475 |
| Default ITSLs | 289 |
| Added TACs | 38 |
| Number of TACs that had 24 hr averaging time, but were converted to annual averaging time | 231 |
| Number of TACs with annual averaging time that were previously 24 hr averaging time AND had values < 75th percentile, therefore are on the new TAC list | 139 |
| **IRSLs | 123 |
| IRSLs with no ITSLs < 75 th\% | 80 |

[^0]Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants
*NEW AvgT "annual*" Some ITSLs with 24-hr avg. time were converted to annual avg. because ITSL was derived to protect for chronic effects.

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| -- | 100 sxl | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | 2-(1-ethoxyethoxy)-6-(trifluoromethyl)benzenethiol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | 2-mercapto-3-(trifluoromethyl)-phenol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | 4-chloro-2-ethyoxy-6-fluoropyrimidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | atlox 848 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | cyclic (phme)2(me)2, d4 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | cyclopentyldichlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | dicyclopentyldichlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | disiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | ethomeen t/30 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | heptamethyl-1-vinyl-1,7dichlorotetrasilazane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | n-chloro-2,6-difluorobenzamide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | sponto 11 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | sponto 723 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | t-det c-40 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | witconol al 69-66 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| -- | o-(1-ethoxyethyl)-2-(propylthio)-3(trifluoromethyl)phenol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| --1 | biosam tp-1.5 | YES | 1st ITSL |  | 0.02 | 1 hr | YES |  |  |  |  |
| --2 | purafect 4000g | YES | 1st ITSL |  | 0.02 | 1 hr | YES |  |  |  |  |
| --3 | fyre-zyme | YES | 1st ITSL |  | 0.15 | annual | YES |  |  |  |  |
| --4 | 1,1,2,4-tetramethyl-1-1-1-sila-2-azacyclopentane | YES | 1st ITSL |  | 0.7 | annual | YES |  |  |  |  |
| --5 | epoxy resin solution | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| --6 | n-butylglucamine | YES | 1st ITSL |  | 6.4 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th } \% \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \end{aligned}$ AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 t h \% \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| --7 | polyglycol 26-3 | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| --8 | ad acid | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| --9 | triethylammonium suleptanate | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| -1-0 | amyl acetate (mixture) | NO |  | >75th\% | 1100 | annual* |  |  |  |  |  |
| 50-00-0 | formaldehyde | YES | 1st ITSL, Carc |  | 9 | 8 hr | YES |  |  |  | 0.08 |
| 50-03-3 | hydrocortisone acetate | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 50-21-5 | lactic acid | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 50-28-2 | estradiol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 50-29-3 | ddt | YES | Carc |  |  |  |  |  |  |  | 0.01 |
| 50-32-8 | benzo(a)pyrene | YES | Carc |  |  |  |  |  |  |  | 0.0005 |
| 51-28-5 | 2,4-dinitrophenol | YES | 1st ITSL |  | 7 | annual* | YES |  |  |  |  |
| 51-79-6 | Ethyl carbamate (Urethane) | YES | $\begin{array}{\|c} \hline \text { HAP Table } \\ 2 \\ \hline \end{array}$ |  |  |  |  |  |  |  |  |
| 53-36-1 | methyl predisolone acetate | YES | 1st ITSL |  | 43 | annual | YES |  |  |  |  |
| 53-70-3 | Dibenz(a,h)anthracene | YES | Carc7 |  |  |  |  |  |  |  |  |
| 56-23-5 | carbon tetrachloride | YES | 1st ITSL, Carc |  | 100 | annual* | YES |  |  |  | 0.17 |
| 56-49-5 | 3-methyl cholanthrene | YES | Cal Carc |  |  |  |  |  |  |  |  |
| 56-55-3 | benz(a)anthracene | YES | Carc7 |  |  |  |  |  |  |  |  |
| 56-81-5 | glycerol | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 57-11-4 | stearic acid | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 57-12-5 | cyanide | YES | 1st ITSL |  | 50 | 1 hr | YES |  |  |  |  |
| 57-15-8 | chlorobutanol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 57-41-0 | phenytoin | YES | Carc |  |  |  |  |  |  |  | 0.07 |
| 57-55-6 | propylene glycol | NO |  | >75th\% | 6000 | annual |  |  |  |  |  |
| 57-83-0 | progesterone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 57-97-9 | 7,12-dimethyl benzanthracene | YES | Cal Carc |  |  |  |  |  |  |  |  |
| 58-36-6 | 10,10'-oxybisphenoxarsine oxide | YES | 1st ITSL |  | 0.2 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{gathered} \text { 1st ITSL } \\ <=75 \text { th\% } \end{gathered}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 60-29-7 | ethyl ether | NO |  | >75th\% | 12000 | 8 hr |  |  |  |  |  |
| 60-34-4 | methyl hydrazine | YES | 1st ITSL, Carc |  | 0.03 | annual* | YES |  |  |  | 0.0087 |
| 60-57-1 | dieldrin | YES | Carc |  |  |  |  |  |  |  | 0.0002 |
| 62-53-3 | aniline | YES | 1st ITSL, 2nd ITSL, Carc |  | 1 | annual | YES | 76 | 8 hr | YES | 0.6 |
| 62-73-7 | dichlorvos | YES | 1st ITSL |  | 0.5 | annual* | YES |  |  |  |  |
| 62-75-9 | n-nitrosodimethylamine | YES | Carc |  |  |  |  |  |  |  | 7E-05 |
| 63-05-8 | androstenedione | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 64-02-8 | ethylenediamine tetra-acetic acid, tetrasodium salt | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 64-04-0 | beta phenylethylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 64-17-5 | ethyl alcohol | NO |  | >75th\% | 19000 | 8 hr |  |  |  |  |  |
| 64-18-6 | formic acid | YES | 1st ITSL |  | 2 | annual* | YES |  |  |  |  |
| 64-19-7 | acetic acid | YES | 1st ITSL |  | 250 | 8 hr | YES |  |  |  |  |
| 64-67-5 | diethyl sulfate | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 66-25-1 | hexanaldehyde | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 67-56-1 | methanol | NO |  | >75th\% | 3250 | 1 hr |  |  |  |  |  |
| 67-63-0 | isopropyl alcohol | NO |  | >75th\% | 220 | annual* |  |  |  |  |  |
| 67-64-1 | acetone | NO |  | >75th\% | 5900 | 8 hr |  |  |  |  |  |
| 67-66-3 | chloroform | YES | Carc |  |  |  |  |  |  |  | 0.4 |
| 67-68-5 | dimethylsulfoxide | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 67-72-1 | hexachloroethane | YES | $\begin{aligned} & \text { 1st ITSL, } \\ & \text { 2nd ITSL, } \\ & \text { Carc } \end{aligned}$ |  | 30 | annual | YES | 1600 | 8 hr | YES | 0.1 |
| 68-12-2 | N,N-dimethylformamide | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 71-23-8 | n-propyl alcohol | NO |  | >75th\% | 730 | annual |  |  |  |  |  |
| 71-36-3 | n-butanol | NO |  | >75th\% | 350 | annual* |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd <br> ITSL <br> AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 71-41-0 | amyl alcohol | NO |  | >75th\% | 120 | annual |  |  |  |  |  |
| 71-43-2 | benzene | YES | 1st ITSL, 2nd ITSL, Carc |  | 30 | annual | YES | 30 | 24 hr | YES | 0.1 |
| 71-55-6 | methyl chloroform | NO |  | >75th\% | 6000 | annual* |  |  |  |  |  |
| 72-54-8 | DDD (TDE) | YES | Carc |  |  |  |  |  |  |  | 0.01 |
| 72-55-9 | DDE, p, p'- | YES | Carc |  |  |  |  |  |  |  | 0.01 |
| 74-83-9 | methyl bromide | YES | 1st ITSL |  | 5 | annual* | YES |  |  |  |  |
| 74-85-1 | ethylene | NO |  | >75th\% | 6240 | annual* |  |  |  |  |  |
| 74-87-3 | methyl chloride | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 90 | annual* | YES |  |  |  | 1.6 |
| 74-88-4 | Methyl iodide (lodomethane) | YES | $\begin{gathered} \text { HAP Table } \\ 2 \end{gathered}$ |  |  |  |  |  |  |  |  |
| 74-89-5 | methylamine | YES | 1st ITSL |  | 64 | 8 hr | YES |  |  |  |  |
| 74-90-8 | hydrogen cyanide | YES | 1st ITSL, 2nd ITSL |  | 0.8 | annual | YES | 50 | 1 hr | YES |  |
| 74-93-1 | methyl mercaptan | YES | 1st ITSL |  | 10 | 1 hr | YES |  |  |  |  |
| 74-97-5 | chlorobromomethane | NO |  | >75th\% | 10600 | 8 hr |  |  |  |  |  |
| 74-99-7 | methyl acetylene | NO |  | >75th\% | 16500 | 8 hr |  |  |  |  |  |
| 75-00-3 | ethyl chloride | NO |  | >75th\% | 10000 | annual* |  |  |  |  |  |
| 75-01-4 | vinyl chloride | YES | $\begin{gathered} \hline \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 100 | annual* | YES |  |  |  | 0.11 |
| 75-04-7 | ethylamine | YES | 1st ITSL |  | 92 | 8 hr | YES |  |  |  |  |
| 75-05-8 | acetonitrile | YES | 1st ITSL |  | 60 | annual* | YES |  |  |  |  |
| 75-07-0 | acetaldehyde | YES | $\begin{gathered} \hline \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 9 | annual* | YES |  |  |  | 0.5 |
| 75-09-2 | methylene chloride | YES | Carc |  | 2000 | annual |  | 14000 | 1 hr |  | 60 |
| 75-12-7 | formamide | YES | Carc |  | 600 | annual* |  |  |  |  | 0.2 |
| 75-15-0 | carbon disulfide | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 75-18-3 | dimethylsulfide | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 t h \% \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 75-21-8 | ethylene oxide | YES | Carc |  |  |  |  |  |  |  | 0.03 |
| 75-25-2 | bromoform | YES | Carc |  |  |  |  |  |  |  | 0.9 |
| 75-27-4 | bromodichloromethane | YES | Carc |  |  |  |  |  |  |  | 0.06 |
| 75-28-5 | isobutane | NO |  | >75th\% | 23800 | 8 hr |  |  |  |  |  |
| 75-29-6 | 2-chloropropane | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 75-31-0 | isopropylamine | YES | 1st ITSL |  | 120 | 8 hr | YES |  |  |  |  |
| 75-34-3 | 1,1-dichloroethane | NO |  | >75th\% | 500 | annual* |  |  |  |  |  |
| 75-35-4 | vinylidene chloride (1,1dichloroethylene) | NO |  | >75th\% | 200 | annua** |  |  |  |  |  |
| 75-36-5 | acetyl chloride | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 75-37-6 | 1,1-difluoroethane | NO |  | >75th\% | 40000 | annual* |  |  |  |  |  |
| 75-38-7 | vinylidene fluoride | YES | 1st ITSL |  | 30 | annual ${ }^{*}$ | YES |  |  |  |  |
| 75-44-5 | phosgene | YES | 1st ITSL |  | 0.3 | annual ${ }^{*}$ | YES |  |  |  |  |
| 75-45-6 | chlorodifluoromethane | NO |  | >75th\% | 50000 | annual* |  |  |  |  |  |
| 75-50-3 | trimethylamine | YES | 1st ITSL |  | 120 | 8 hr | YES |  |  |  |  |
| 75-52-5 | nitromethane | YES | $\begin{aligned} & \text { 1st ITSL, } \\ & \text { Carc } \end{aligned}$ |  | 70 | annua** | YES |  |  |  | 0.1 |
| 75-54-7 | methyldichlorosilane | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 75-55-8 | 1,2-propylenimine | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 75-56-9 | propylene oxide | YES | $\begin{aligned} & \text { 1st ITSL, } \\ & \text { Carc } \end{aligned}$ |  | 30 | annual* | YES |  |  |  | 0.3 |
| 75-64-9 | t-butylamine | YES | 1st ITSL |  | 60 | annual | YES |  |  |  |  |
| 75-65-0 | t-butanol | NO |  | >75th\% | 1890 | annual* |  |  |  |  |  |
| 75-68-3 | 1-chloro-1,1-difluoroethane | NO |  | >75th\% | 50000 | annual* |  |  |  |  |  |
| 75-69-4 | trichlorofluoromethane | NO |  | >75th\% | 56200 | 1 hr |  |  |  |  |  |
| 75-71-8 | dichlorodifluoromethane | NO |  | >75th\% | 49500 | 8 hr |  |  |  |  |  |
| 75-75-2 | methane sulfonic acid | YES | 1st ITSL |  | 1.4 | annual | YES |  |  |  |  |
| 75-76-3 | tetramethylsilane | NO |  | >75th\% | 1300 | annual |  |  |  |  |  |
| 75-77-4 | trimethylchlorosilane | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants
*NEW AvgT "annual*" Some ITSLs with 24-hr avg. time were converted to annual avg. because ITSL was derived to protect for chronic effects.

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{gathered} \text { 1st ITSL } \\ <=75 \text { th\% } \end{gathered}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 75-78-5 | dimethyldichlorosilane | YES | 1st ITSL |  | 6.2 | annual | YES |  |  |  |  |
| 75-79-6 | methyltrichlorosilane | NO |  | >75th\% | 109 | annual |  |  |  |  |  |
| 75-94-5 | vinyltrichlorosilane | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 76-05-1 | trifluoroacetic acid | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 76-13-1 | 1,1,2-trichloro-1,2,2-trifluoroethane | NO |  | >75th\% | 19140 | annual ${ }^{\text {* }}$ |  |  |  |  |  |
| 76-14-2 | dichlorotetrafluoroethan | NO |  | >75th\% | 69000 | 8 hr |  |  |  |  |  |
| 76-44-8 | heptachlor | YES | Carc |  |  |  |  |  |  |  | 0.0008 |
| 76-83-5 | triphenyl methyl chloride | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 77-47-4 | hexachlorocyclopentadiene | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |
| 77-48-5 | 1,3-dibromo-5,5-dimethylhydantoin | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 77-58-7 | dibutyl tin dilaurate | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 77-73-6 | dicyclopentadiene | YES | 1st ITSL |  | 1 | annual* | YES |  |  |  |  |
| 77-76-9 | 2,2-dimethoxypropane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 77-78-1 | dimethyl sulfate | YES | 1st ITSL |  | 0.5 | 8 hr | YES |  |  |  |  |
| 77-93-0 | triethyl citrate | NO |  | >75th\% | 290 | annual |  |  |  |  |  |
| 78-07-9 | ethyltriethoxysilane | YES | 1st ITSL |  | 44 | annual | YES |  |  |  |  |
| 78-10-4 | ethyl silicate | YES | 1st ITSL |  | 850 | 8 hr | YES |  |  |  |  |
| 78-59-1 | isophorone | YES | $\begin{aligned} & \text { 1st ITSL, } \\ & \text { Carc } \end{aligned}$ |  | 280 | 1 hr | YES |  |  |  | 3.7 |
| 78-78-4 | 2-methyl butane | NO |  | >75th\% | 17700 | 8 hr |  |  |  |  |  |
| 78-79-5 | isoprene | YES | Carc |  |  |  |  |  |  |  | 0.02 |
| 78-83-1 | isobutyl alcohol | YES | 1st ITSL |  | 1500 | 8 hr | YES |  |  |  |  |
| 78-84-2 | isobutyraldehyde | NO |  | >75th\% | 160 | annual* |  |  |  |  |  |
| 78-87-5 | propylene dichloride | YES | 1st ITSL |  | 4 | annual* | YES |  |  |  |  |
| 78-92-2 | sec-butanol | NO |  | >75th\% | 3000 | 8 hr |  |  |  |  |  |
| 78-93-3 | methyl ethyl ketone | NO |  | >75th\% | 5000 | annual* |  |  |  |  |  |
| 78-96-6 | monoisopropanolamine | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 79-00-5 | 1,1,2-trichloroethane | YES | Carc |  |  |  |  |  |  |  | 0.06 |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | $\begin{aligned} & \text { Future } \\ & \text { TAC? } \end{aligned}$ | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { Avg } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \mathrm{th} \% \\ \hline \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 79-01-6 | trichloroethylene | YES | 1st ITSL, Carc |  | 2 | annual | YES | 10000 | 24 hr |  | 0.2 |
| 79-06-1 | acrylamide | YES | 1st ITSL, Carc |  | 6 | annual* | YES |  |  |  | 0.005 |
| 79-09-4 | propionic acid | YES | 1st ITSL |  | 300 | 8 hr | YES |  |  |  |  |
| 79-10-7 | acrylic acid | YES | 1st ITSL |  | 1 | annual* | YES |  |  |  |  |
| 79-11-8 | Chloroacetic acid | YES | $\begin{gathered} \text { HAP Table } \\ 2 \\ \hline \end{gathered}$ |  |  |  |  |  |  |  |  |
| 79-14-1 | hydroxyacetic acid/ glycolic acid | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 79-20-9 | methyl acetate | NO |  | >75th\% | 6100 | 8 hr |  |  |  |  |  |
| 79-24-3 | nitroethane | YES | 1st ITSL |  | 60 | annual* | YES |  |  |  |  |
| 79-29-8 | 2,3-dimethylbutane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 79-31-2 | isobutyric acid | YES | 1st ITSL |  | 0.9 | annual | YES |  |  |  |  |
| 79-34-5 | 1,1,2,2-tetrachloroethane | YES | Carc |  |  |  |  |  |  |  | 0.02 |
| 79-41-4 | methacrylic acid | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 79-46-9 | 2-nitropropane | YES | 1st ITSL, Carc |  | 20 | annual* | YES |  |  |  | 0.0004 |
| 79-92-5 | camphene | YES | 1st ITSL |  | 80 | annual | YES |  |  |  |  |
| 80-15-9 | cumene hydroperoxide | YES | 1st ITSL |  | 6 | annual* | YES |  |  |  |  |
| 80-43-3 | dicumyl peroxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 80-56-8 | pinene, alpha | YES | 1st ITSL |  | 1120 | 8 hr | YES |  |  |  |  |
| 80-62-6 | methyl methacrylate | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 80-73-9 | $\mathrm{n}, \mathrm{n}$ '-dimethylethyleneurea | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 82-68-8 | pentachloronitrobenzene | YES | 1st ITSL |  | 11 | annual* | YES |  |  |  |  |
| 83-32-9 | acenaphthene | NO |  | >75th\% | 210 | annual* |  |  |  |  |  |
| 84-66-2 | diethyl phthalate | YES | 1st ITSL |  | 50 | 8 hr | YES | 2800 | 24 hr |  |  |
| 85-01-8 | phenanthrene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 85-68-7 | butyl benzyl phthalate | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 86-73-7 | fluorene | NO |  | >75th\% | 140 | annual* |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | 2nd ITSL $\left(\mu \mathrm{g} / \mathrm{m}^{3}\right)$ | 2nd <br> ITSL <br> AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 86-74-8 | carbazole | YES | Carc |  |  |  |  |  |  |  | 0.4 |
| 87-61-6 | 1,2,3-trichlorobenzene | YES | 1st ITSL |  | 27 | annual* | YES |  |  |  |  |
| 87-62-7 | 2,6-xylidine | YES | Carc |  |  |  |  |  |  |  | 0.78 |
| 87-68-3 | hexachlorobutadiene | YES | Carc |  |  |  |  |  |  |  | 0.05 |
| 87-86-5 | pentachlorophenol | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 20 | annual* | YES |  |  |  | 0.009 |
| 87-90-1 | 1,3,5-trichloroisocyanuric acid | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 88-06-2 | 2,4,6-trichlorophenol | YES | Carc |  |  |  |  |  |  |  | 0.3 |
| 88-12-0 | n-vinylpyrrolidinone | YES | Carc |  |  |  |  |  |  |  | 0.04 |
| 88-65-3 | o-bromobenzoic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 88-73-3 | 1-chloro-2-nitrobenzene | YES | Carc |  |  |  |  |  |  |  | 0.21 |
| 88-85-7 | dinoseb | YES | 1st ITSL |  | 4 | annual* | YES |  |  |  |  |
| 90-02-8 | salicylaldehyde | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 90-12-0 | 1-methyl naphthalene | YES | Carc |  | 250 | annual* |  |  |  |  | 0.14 |
| 90-43-7 | o-phenylphenol | YES | Carc |  |  |  |  |  |  |  | 1.1 |
| 90-72-2 | 2,4,6-tri(dimethylaminomethyl)phenol | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 91-01-0 | benzhydrol | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 91-17-8 | decahydronaphthalene | YES | Carc |  |  |  |  |  |  |  | 0.03 |
| 91-20-3 | naphthalene | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 3 | annual* | YES |  |  |  | 0.08 |
| 91-22-5 | quinoline | YES | Carc |  |  |  |  |  |  |  | 0.001 |
| 91-44-1 | 7-diethylamino-4-methyl coumarin | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 91-57-6 | 2-methylnaphthalene | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 91-59-8 | 2-naphthylamine | YES | Carc |  |  |  |  |  |  |  | 0.0001 |
| 91-94-1 | dichlorobenzidine | YES | Carc |  |  |  |  |  |  |  | 0.002 |
| 92-52-4 | biphenyl | YES | 1st ITSL |  | 13 | 8 hr | YES |  |  |  |  |
| 92-87-5 | benzidine | YES | Carc |  |  |  |  |  |  |  | 2E-05 |
| 93-14-1 | guaifenesin | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | 1st ITSL $\left(\mu \mathrm{g} / \mathrm{m}^{3}\right)$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 93-58-3 | methyl benzoate | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 93-59-4 | peroxybenzoic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 93-83-4 | oleoyl diethanolamine | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 94-96-2 | 2-ethyl-1,3-hexanediol | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 95-16-9 | benzothiazole | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 95-38-5 | oyel hydroxyethylimidazoline | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 95-47-6 | o-xylene | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 95-48-7 | o-cresol | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 95-49-8 | monochlorotoluene | YES | 1st ITSL |  | 70 | annual* | YES |  |  |  |  |
| 95-50-1 | 1,2-dichlorobenzene | NO |  | >75th\% | 300 | annual* |  |  |  |  |  |
| 95-51-2 | 2-chloroaniline | YES | 1st ITSL |  | 10 | annual* | YES |  |  |  |  |
| 95-53-4 | o-toluidine | YES | Carc |  |  |  |  |  |  |  | 0.07 |
| 95-57-8 | 2-chlorophenol | YES | 1st ITSL |  | 18 | annual* | YES |  |  |  |  |
| 95-63-6 | 1,2,4-trimethylbenzene | YES | 1st ITSL, 2nd ITSL |  | 50 | annual | YES | 1200 | 8 hr | YES |  |
| 95-65-8 | 3,4-dimethyl phenol | YES | 1st ITSL |  | 3.5 | annual* | YES |  |  |  |  |
| 95-74-9 | 3-chloro-p-toluidine | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 95-87-4 | 2,5-dimethylphenol | YES | 1st ITSL |  | 0.7 | annual | YES |  |  |  |  |
| 95-93-2 | 1,2,4,5-tetramethyl benzene | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 95-94-3 | 1,2,4,5-tetrachlorobenzene | YES | 1st ITSL |  | 1 | annual* | YES |  |  |  |  |
| 95-95-4 | 2,4,5-trichlorophenol | NO |  | >75th\% | 350 | annual* |  |  |  |  |  |
| 96-12-8 | dibromochloropropane | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 0.2 | annual* | YES |  |  |  | 0.0001 |
| 96-14-0 | 3-methylpentane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 96-18-4 | 1,2,3-trichloropropane | YES | 1st ITSL |  | 0.3 | annual* | YES |  |  |  |  |
| 96-23-1 | 1,3-dichloro-2-propanol | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 3 | annual* | YES |  |  |  | 0.07 |
| 96-29-7 | methylethylketoxime | YES | Carc |  |  |  |  |  |  |  | 2.5 |
| 96-33-3 | methyl acrylate | YES | 1st ITSL |  | 70 | annual* | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{array}{\|c\|} \hline \text { 1st ITSL } \\ <=75 \text { th\% } \end{array}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 96-37-7 | methylcyclopentane | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 96-45-7 | ethylene thiourea | YES | 1st ITSL, Carc |  | 0.28 | annual* | YES |  |  |  | 0.01 |
| 96-48-0 | gamma-butyrolactone | NO |  | >75th\% | 280 | annual* |  |  |  |  |  |
| 96-49-1 | ethylene carbonate | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 96-80-0 | diisopropylaminoethanol | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 97-64-3 | ethyl lactate | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 97-85-8 | isobutyl isobutyrate | NO |  | >75th\% | 300 | annual |  |  |  |  |  |
| 97-86-9 | isobutyl methacrylate | NO |  | >75th\% | 600 | annual |  |  |  |  |  |
| 97-88-1 | n-butyl methacrylate | NO |  | >75th\% | 569 | annual |  |  |  |  |  |
| 97-95-0 | 2-ethyl butanol | YES | 1st ITSL |  | 40 | annual | YES |  |  |  |  |
| 97-99-4 | tetrahydrofuryl methanol | YES | 1st ITSL |  | 52 | annual | YES |  |  |  |  |
| 98-00-0 | furfuryl alcohol | YES | 1st ITSL, Carc |  | 1 | annual* | YES |  |  |  | 0.03 |
| 98-01-1 | furfural | YES | Carc |  |  |  |  |  |  |  | 0.06 |
| 98-06-6 | tert-butylbenzene | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 98-13-5 | phenyltrichlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 98-17-9 | m-trifluoromethylphenol | YES | 1st ITSL |  | 0.08 | annual | YES |  |  |  |  |
| 98-29-3 | t-butylcatechol | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 98-56-6 | p-chlorobenzotrifluoride | YES | 1st ITSL |  | 70 | annual ${ }^{\text {* }}$ | YES |  |  |  |  |
| 98-82-8 | cumene | YES | Carc |  | 400 | annual* |  |  |  |  | 0.1 |
| 98-83-9 | alpha-methyl styrene | NO |  | >75th\% | 230 | annual* |  |  |  |  |  |
| 98-84-0 | dl-alpha phenylethylamine | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 98-86-2 | acetophenone | YES | 1st ITSL |  | 490 | 8 hr | YES |  |  |  |  |
| 98-95-3 | nitrobenzene | YES | 1st ITSL, Carc |  | 9 | annual* | YES |  |  |  | 0.025 |
| 99-87-6 | p-isopropyltoluene | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 99-97-8 | $n, \mathrm{n}$-dimethyl-p-toluidine | YES | 1st ITSL |  | 28 | annual | YES |  |  |  |  |
| 100-02-7 | 4-nitrophenol | YES | 1st ITSL |  | 0.7 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th } \% \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \end{aligned}$ | $\begin{array}{\|c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 100-06-1 | 4-methoxyacetophenone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 100-36-7 | 2-diethylaminoethylamine | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 100-37-8 | 2-diethylaminoethanol (deae) | YES | 1st ITSL |  | 4 | annual* | YES |  |  |  |  |
| 100-40-3 | 4-vinylcyclohexene | YES | 1st ITSL |  | 4 | 8 hr | YES |  |  |  |  |
| 100-41-4 | ethylbenzene | YES | Carc |  | 1000 | annual* |  |  |  |  | 3 |
| 100-42-5 | styrene | YES | Carc |  | 1000 | annual* |  |  |  |  | 1.7 |
| 100-44-7 | benzyl chloride | YES | Carc |  |  |  |  |  |  |  | 0.02 |
| 100-46-9 | benzylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 100-51-6 | benzyl alcohol | NO |  | >75th\% | 5000 | annual* |  |  |  |  |  |
| 100-52-7 | benzaldehyde | YES | Carc |  |  |  |  |  |  |  | 0.4 |
| 100-85-6 | benzyltrimethylammonium hydroxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 100-97-0 | hexamethylenetetramine | YES | 1st ITSL |  | 100 | annual | YES |  |  |  |  |
| 101-68-8 | methylene diphenyl diisocyanate | YES | 1st ITSL |  | 0.6 | annual* | YES |  |  |  |  |
| 101-84-8 | diphenyloxide | YES | 1st ITSL |  | 70 | 8 hr | YES |  |  |  |  |
| 102-69-2 | tripropylamine | YES | 1st ITSL |  | 0.2 | annual | YES |  |  |  |  |
| 102-71-6 | triethanolamine | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 102-76-1 | triacetin | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 102-79-4 | butyldiethanolamine | YES | 1st ITSL |  | 14 | annual | YES |  |  |  |  |
| 102-81-8 | 2-n-dibutylaminoethanol | YES | 1st ITSL |  | 28 | annual* | YES |  |  |  |  |
| 102-82-9 | tributylamine | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 103-09-3 | 2-ethylhexyl acetate | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 103-11-7 | 2-ethylhexyl acrylate | YES | 1st ITSL |  | 18 | annual | YES |  |  |  |  |
| 103-23-1 | di (2-ethylhexyl) adipate | YES | Carc |  |  |  |  |  |  |  | 3 |
| 103-33-3 | azobenzene | YES | Carc |  |  |  |  |  |  |  | 0.03 |
| 103-63-9 | 2-bromoethyl benzene | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 103-65-1 | propylbenzene | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 103-83-3 | benzyl dimethylamine | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 103-99-1 | N -stearoyl-4-aminophenol | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 104-15-4 | p-toluenesulfonic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 104-51-8 | n-butylbenzene | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 104-68-7 | diethylene glycol monophenyl ether | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 104-75-6 | 2-ethylhexylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 104-76-7 | 2-ethylhexanol | YES | 1st ITSL |  | 70 | annual | YES |  |  |  |  |
| 104-78-9 | n,n-diethyl-1,3-propanediamine | NO |  | >75th\% | 140 | annual |  |  |  |  |  |
| 104-87-0 | p-tolualdehyde | NO |  | >75th\% | 440 | annual* |  |  |  |  |  |
| 105-39-5 | ethyl chloroacetate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 105-53-3 | diethylmalonate | YES | 1st ITSL |  | 50 | annual | YES |  |  |  |  |
| 105-56-6 | ethyl cyanoacetate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 105-58-8 | diethyl carbonate | NO |  | >75th\% | 5000 | annual* |  |  |  |  |  |
| 105-59-9 | methyldiethanolamine | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 105-60-2 | caprolactam | YES | 1st ITSL |  | 10 | 8 hr | YES |  |  |  |  |
| 105-67-9 | 2,4-dimethylphenol | YES | 1st ITSL |  | 70 | annual* | YES |  |  |  |  |
| 106-36-5 | propyl propionate | YES | 1st ITSL |  | 84 | annual | YES |  |  |  |  |
| 106-42-3 | p-xylene | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 106-46-7 | 1,4-dichlorobenzene | YES | Carc |  | 800 | annual* |  |  |  |  | 0.14 |
| 106-49-0 | p-toluidine | YES | Carc |  |  |  |  |  |  |  | 0.03 |
| 106-51-4 | Quinone (p-benzoquinone) | YES | $\begin{gathered} \text { HAP Table } \\ 2 \end{gathered}$ |  |  |  |  |  |  |  |  |
| 106-79-6 | dimethyl decanedioate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 106-88-7 | 1,2-butylene oxide | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 20 | annual* | YES |  |  |  | 1.2 |
| 106-89-8 | epichlorohydrin | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 1 | annual* | YES |  |  |  | 0.8 |
| 106-91-2 | glycidyl methacrylate | YES | 1st ITSL, 2nd ITSL |  | 0.8 | annual | YES | 16 | 24 hr | YES |  |
| 106-92-3 | allyl glycidyl ether | YES | Carc |  |  |  |  |  |  |  | 0.1 |
| 106-93-4 | ethylene dibromide | YES | 1st ITSL, |  | 9 | annual* | YES |  |  |  | 0.002 |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ |  | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  |  |  | Carc |  |  |  |  |  |  |  |  |
| 106-94-5 | propyl bromide | YES | 1st ITSL |  | 49 | annual | YES |  |  |  |  |
| 106-97-8 | butane | NO |  | >75th\% | 23800 | 8 hr |  |  |  |  |  |
| 106-99-0 | 1,3-butadiene | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 2 | annual* | YES |  |  |  | 0.03 |
| 107-00-6 | ethylacetylene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 107-02-8 | acrolein | YES | 1st ITSL, 2nd ITSL |  | 0.16 | annual | YES | 5 | 1 hr | YES |  |
| 107-03-9 | 1-propanethiol | YES | 1st ITSL |  | 16 | 1 hr | YES |  |  |  |  |
| 107-05-1 | allyl chloride | YES | 1st ITSL, 2nd ITSL |  | 1 | annual | YES | 31 | 8 hr | YES |  |
| 107-06-2 | 1,2-dichloroethane | YES | Carc |  |  |  |  |  |  |  | 0.04 |
| 107-10-8 | propylamine | NO |  | >75th\% | 112 | annual |  |  |  |  |  |
| 107-13-1 | acrylonitrile | YES | $\begin{gathered} \hline \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 2 | annual* | YES |  |  |  | 0.01 |
| 107-15-3 | ethylene diamine | YES | 1st ITSL |  | 0.03 | annual | YES |  |  |  |  |
| 107-18-6 | allyl alcohol | YES | 1st ITSL |  | 18 | annual* | YES |  |  |  |  |
| 107-21-1 | ethylene glycol | NO |  | >75th\% | 1000 | 1 hr |  |  |  |  |  |
| 107-31-3 | methyl formate | YES | 1st ITSL |  | 1250 | 8 hr | YES |  |  |  |  |
| 107-39-1 | diisobutylene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 107-41-5 | hexylene glycol | NO |  | >75th\% | 1210 | 1 hr |  |  |  |  |  |
| 107-46-0 | hexamethyldisiloxane | NO |  | >75th\% | 240 | annual |  |  |  |  |  |
| 107-51-7 | octamethyltrisiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 107-54-0 | 3,5-dimethyl-1-hexyn-3-0 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 107-66-4 | dibutyl phosphate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 107-68-6 | n-methyl taurine | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 107-71-1 | t-butyl peroxyacetate | YES | 1st ITSL |  | 0.06 | annual | YES |  |  |  |  |
| 107-83-5 | 2-methylpentane | NO |  | >75th\% | 17600 | 8 hr |  |  |  |  |  |
| 107-87-9 | methyl propyl ketone | NO |  | $>75$ th\% | 5300 | 8 hr |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 107-92-6 | butyric acid | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 107-98-2 | propylene glycol monomethyl ether | NO |  | >75th\% | 2000 | annual* |  |  |  |  |  |
| 108-01-0 | dimethylethanolamine | YES | 1st ITSL, 2nd ITSL |  | 5.2 | annual | YES | 220 | 8 hr | YES |  |
| 108-03-2 | 1-nitropropane | YES | 1st ITSL |  | 900 | 8 hr | YES |  |  |  |  |
| 108-05-4 | vinyl acetate | NO |  | >75th\% | 200 | annual* |  |  |  |  |  |
| 108-08-7 | 2,4-dimethylpentane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 108-10-1 | methyl isobutyl ketone | NO |  | >75th\% | 3000 | annual* |  |  |  |  |  |
| 108-11-2 | methyl amyl alcohol | YES | 1st ITSL |  | 1000 | 8 hr | YES |  |  |  |  |
| 108-16-7 | dimethylamino-2-propanol | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 108-18-9 | diisopropylamine | YES | 1st ITSL |  | 200 | 8 hr | YES |  |  |  |  |
| 108-20-3 | diisopropyl ether | NO |  | >75th\% | 358 | annual* |  |  |  |  |  |
| 108-21-4 | isopropyl acetate | NO |  | >75th\% | 4200 | 8 hr |  |  |  |  |  |
| 108-31-6 | maleic anhydride | YES | 1st ITSL |  | 0.1 | 8 hr | YES |  |  |  |  |
| 108-32-7 | propylene carbonate | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 108-38-3 | m-xylene | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 108-46-3 | resorcinol | YES | 1st ITSL |  | 27 | annual* | YES |  |  |  |  |
| 108-60-1 | bis(2-chloroisopropyl)ether | NO |  | >75th\% | 140 | annual* |  |  |  |  |  |
| 108-65-6 | propylene glycol monomethyl ether acetate | NO |  | >75th\% | 3000 | annual* |  |  |  |  |  |
| 108-67-8 | 1,3,5-trimethyl benzene | YES | 1st ITSL, 2nd ITSL |  | 50 | annual | YES | 1200 | 8 hr | YES |  |
| 108-68-9 | 3,5-dimethylphenol | YES | 1st ITSL |  | 0.8 | annual | YES |  |  |  |  |
| 108-78-1 | melamine | YES | Carc |  |  |  |  |  |  |  | 1.5 |
| 108-82-7 | 2,6-dimethyl-4-heptanol | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 108-83-8 | diisobutyl ketone | YES | 1st ITSL |  | 1500 | 8 hr | YES |  |  |  |  |
| 108-86-1 | bromobenzene | YES | 1st ITSL |  | 60 | annual* | YES |  |  |  |  |
| 108-87-2 | methylcyclohexane | NO |  | >75th\% | 16000 | 8 hr |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | $\begin{aligned} & \text { Basis } \\ & \text { for Yes } \end{aligned}$ | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \\ \hline \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 108-88-3 | toluene | NO |  | >75th\% | 5000 | annual* |  |  |  |  |  |
| 108-90-7 | chlorobenzene | YES | 1st ITSL |  | 50 | annual | YES | 4400 | 8 hr |  |  |
| 108-94-1 | cyclohexanone | YES | 1st ITSL |  | 800 | 8 hr | YES |  |  |  |  |
| 108-95-2 | phenol | YES | 1st ITSL |  | 190 | 8 hr | YES |  |  |  |  |
| 108-99-6 | 3-picoline | YES | 1st ITSL |  | 80 | annual | YES |  |  |  |  |
| 109-06-8 | alpha-picoline | YES | 1st ITSL |  | 24 | annual* | YES |  |  |  |  |
| 109-56-8 | isopropylethanolamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 109-60-4 | n-propyl acetate | NO |  | >75th\% | 8350 | 8 hr |  |  |  |  |  |
| 109-65-9 | 1-bromobutane | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 109-66-0 | pentane | NO |  | >75th\% | 17700 | 8 hr |  |  |  |  |  |
| 109-69-3 | n-butyl chloride | NO |  | >75th\% | 1500 | annual* |  |  |  |  |  |
| 109-70-6 | 1-bromo-3-chloropropane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 109-83-1 | 2-methylaminoethanol | YES | 1st ITSL |  | 38 | annual | YES |  |  |  |  |
| 109-86-4 | 2-methoxyethanol | YES | 1st ITSL |  | 20 | annual* | YES |  |  |  |  |
| 109-89-7 | diethylamine | YES | 1st ITSL |  | 150 | 8 hr | YES |  |  |  |  |
| 109-92-2 | ethyl vinyl ether | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 109-94-4 | ethyl formate | NO |  | >75th\% | 3000 | 8 hr |  |  |  |  |  |
| 109-99-9 | tetrahydrofuran | NO |  | >75th\% | 8000 | annual |  |  |  |  |  |
| 110-00-9 | furan | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 4 | annual* | YES |  |  |  | 0.0002 |
| 110-12-3 | methyl isoamy ketone | YES | 1st ITSL |  | 2300 | 8 hr | YES |  |  |  |  |
| 110-16-7 | maleic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 110-19-0 | isobutyl acetate | YES | 1st ITSL |  | 480 | 8 hr | YES |  |  |  |  |
| 110-30-5 | $\mathrm{n}, \mathrm{n}$ '-ethylene bis-octadecanamide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 110-43-0 | methyl n-amyl ketone | YES | 1st ITSL |  | 2330 | 8 hr | YES |  |  |  |  |
| 110-49-6 | ethylene glycol monomethyl ether acetate | YES | 1st ITSL |  | 31 | annual ${ }^{*}$ | YES |  |  |  |  |
| 110-54-3 | n-hexane | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th } \% \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \end{aligned}$ | $\begin{array}{\|c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 110-58-7 | amylamine | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 110-61-2 | succinonitrile | YES | 1st ITSL |  | 0.8 | annual | YES |  |  |  |  |
| 110-62-3 | valeraldehyde | YES | 1st ITSL |  | 1760 | 8 hr | YES |  |  |  |  |
| 110-63-4 | 1,4 butanediol | YES | 1st ITSL |  | 79 | annual | YES |  |  |  |  |
| 110-71-4 | ethylene glycol dimethyl ether | YES | 1st ITSL |  | 24 | annual* | YES |  |  |  |  |
| 110-73-6 | 2-ethylaminoethanol | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 110-80-5 | 2-ethoxyethanol | NO |  | >75th\% | 200 | annual* |  |  |  |  |  |
| 110-82-7 | cyclohexane | NO |  | >75th\% | 6000 | annual* |  |  |  |  |  |
| 110-83-8 | cyclohexene | NO |  | >75th\% | 10000 | 8 hr |  |  |  |  |  |
| 110-86-1 | pyridine | YES | 1st ITSL |  | 3.5 | annual* | YES |  |  |  |  |
| 110-89-4 | piperidine | NO |  | >75th\% | 140 | annual |  |  |  |  |  |
| 110-97-4 | diisopropanolamine | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 111-13-7 | 2-octanone | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 111-15-9 | ethylene glycol monoethyl ether acetate | NO |  | >75th\% | 293 | annual* |  |  |  |  |  |
| 111-30-8 | glutaraldehyde | YES | 1st ITSL, 2nd ITSL |  | 0.08 | annual | YES | 0.2 | 1 hr | YES |  |
| 111-42-2 | diethanolamine | YES | 1st ITSL |  | 5 | annual* | YES |  |  |  |  |
| 111-44-4 | bis-2-chloroethylether | YES | Carc |  |  |  |  |  |  |  | 0.003 |
| 111-46-6 | diethylene glycol | NO |  | >75th\% | 21000 | annual* |  |  |  |  |  |
| 111-75-1 | 2-butylaminoethanol | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 111-76-2 | 2-butoxyethanol | NO |  | >75th\% | 1600 | annual* |  |  |  |  |  |
| 111-77-3 | diethylene glycol monomethyl ether | NO |  | >75th\% | 190 | annual* |  |  |  |  |  |
| 111-84-2 | n-nonane | NO |  | >75th\% | 550 | annual* |  |  |  |  |  |
| 111-90-0 | diethylene glycol monoethyl ether | NO |  | >75th\% | 1750 | annual* |  |  |  |  |  |
| 111-92-2 | dibutylamine | YES | 1st ITSL |  | 23 | annual | YES |  |  |  |  |
| 112-06-1 | n-heptyl acetate | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 112-07-2 | ethylene glycol monobutyl ether acetate | NO |  | >75th\% | 17600 | annual* |  |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | 2nd <br> ITSL <br> $<=75 t h \%$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 112-15-2 | diethylene glycol monoethyl ether acetate | YES | 1st ITSL |  | 18 | annual | YES |  |  |  |  |
| 112-24-3 | triethylene tetramine | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 112-25-4 | ethylene glycol monohexyl ether | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 112-34-5 | butyl carbitol | YES | 1st ITSL |  | 20 | annual* | YES |  |  |  |  |
| 112-48-1 | ethylene glycol dibutyl ether | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 112-50-5 | triethylene glycol monoethyl ether | YES | 1st ITSL |  | 100 | annual | YES |  |  |  |  |
| 112-55-0 | n-dodecyl mercaptan | YES | 1st ITSL |  | 8 | 8 hr | YES |  |  |  |  |
| 112-80-1 | oleic acid | NO |  | >75th\% | 242 | annual |  |  |  |  |  |
| 115-07-1 | propylene | NO |  | >75th\% | 1500 | annual* |  |  |  |  |  |
| 115-10-6 | dimethyl ether | YES | 1st ITSL |  | 66 | annual | YES |  |  |  |  |
| 115-11-7 | isobutylene | YES | 1st ITSL |  | 21 | annual | YES |  |  |  |  |
| 115-19-5 | methyl butynol | YES | 1st ITSL |  | 6.5 | annual | YES |  |  |  |  |
| 116-11-0 | 2-methoxy-1-propene | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 116-14-3 | tetrafluoroethylene | YES | Carc |  |  |  |  |  |  |  | 0.4 |
| 117-81-7 | diethyl hexyl phthalate | YES | Carc |  |  |  |  |  |  |  | 0.2 |
| 117-84-0 | di-n-octyl phthalate | NO |  | >75th\% | 470 | annual* |  |  |  |  |  |
| 118-52-5 | 1,3-dichloro-5,5-dimethylhydantoin | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 118-74-1 | hexachlorobenzene | YES | Carc |  |  |  |  |  |  |  | 0.002 |
| 118-91-2 | o-chlorobenzoic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 119-53-9 | benzoin | YES | 1st ITSL |  | 32 | annual* | YES |  |  |  |  |
| 119-90-4 | 3,3-dimethoxybenzidine | YES | $\begin{gathered} \hline \text { HAP Table } \\ 2 \end{gathered}$ |  |  |  |  |  |  |  |  |
| 120-07-0 | phenyldiethanolamine | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 120-12-7 | anthracene | NO |  | >75th\% | 1000 | annual* |  |  |  |  |  |
| 120-82-1 | 1,2,4-trichlorobenzene | YES | 1st ITSL |  | 4 | annual* | YES |  |  |  |  |
| 120-83-2 | 2,4-dichlorophenol | YES | 1st ITSL |  | 77 | annual | YES |  |  |  |  |
| 121-14-2 | 2,4-dinitrotoluene | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 2 | 8 hr | YES |  |  |  | 0.009 |

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| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{array}{\|c} \text { 1st ITSL } \\ <=75 \text { th\% } \end{array}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \\ \hline \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 121-43-7 | trimethoxyborine | YES | 1st ITSL |  | 18 | annual | YES |  |  |  |  |
| 121-44-8 | triethylamine | YES | 1st ITSL |  | 7 | annual* | YES |  |  |  |  |
| 121-69-7 | dimethylaniline | YES | Carc |  |  |  |  |  |  |  | 0.085 |
| 121-93-7 | isopropyldiethanolamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 122-20-3 | triisopropanolamine (tipa) | YES | 1st ITSL |  | 19 | annual | YES |  |  |  |  |
| 122-60-1 | phenyl glycidyl ether | YES | Carc |  |  |  |  |  |  |  | 0.1 |
| 122-79-2 | phenyl acetate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 122-99-6 | ethylene glycol monophenyl ether | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 123-03-5 | cetylpyridinium chloride | YES | 1st ITSL |  | 1.8 | annual | YES |  |  |  |  |
| 123-05-7 | 2-ethylhexanal | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 123-19-3 | dipropyl ketone | NO |  | >75th\% | 250 | annual* |  |  |  |  |  |
| 123-31-9 | Hydroquinone | YES | $\begin{array}{\|c} \hline \text { HAP Table } \\ 2 \end{array}$ |  |  |  |  |  |  |  |  |
| 123-38-6 | propionaldehyde | YES | 1st ITSL |  | 8 | annual* | YES |  |  |  |  |
| 123-42-2 | diacetone alcohol | NO |  | >75th\% | 2375 | 8 hr |  |  |  |  |  |
| 123-51-3 | isoamyl alcohol | YES | 1st ITSL |  | 360 | 8 hr | YES |  |  |  |  |
| 123-54-6 | 2,4-pentanedione | YES | 1st ITSL |  | 25 | annual* | YES |  |  |  |  |
| 123-72-8 | butyraldehyde | YES | 1st ITSL |  | 7 | annual* | YES |  |  |  |  |
| 123-86-4 | n-butyl acetate | NO |  | >75th\% | 7100 | 8 hr |  |  |  |  |  |
| 123-91-1 | 1,4-dioxane | YES | 1st ITSL, Carc |  | 100 | annual* | YES |  |  |  | 0.04 |
| 123-92-2 | isoamyl acetate | NO |  | >75th\% | 2700 | 8 hr |  | 5300 | 1 hr |  |  |
| 124-04-9 | adipic acid | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 124-07-2 | octanoic acid | YES | 1st ITSL |  | 33 | annual | YES |  |  |  |  |
| 124-17-4 | diethylene glycol monobutyl ether acetate | YES | 1st ITSL |  | 25 | annual* | YES |  |  |  |  |
| 124-26-5 | octadecanamide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 124-28-7 | $\mathrm{N}, \mathrm{N}$-dimethyl octadecylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 124-41-4 | sodium methylate | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{gathered} \text { 1st ITSL } \\ <=75 \text { th\% } \\ \hline \end{gathered}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \\ \hline \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 124-48-1 | chlorodibromomethane | YES | Carc |  |  |  |  |  |  |  | 0.04 |
| 124-63-0 | methyl sulfonyl chloride | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 124-68-5 | 2-amino-2-methyl-1-propanol | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 124-70-9 | methylvinyldichlorosilane | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 126-06-7 | 3-bromo-1-chloro-5,5dimethylhydantoin | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 126-30-7 | neopentyl glycol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 126-72-7 | tris(2,3-dibromopropyl) phosphate | YES | Carc |  |  |  |  |  |  |  | 0.002 |
| 126-73-8 | tributyl phosphate | YES | 1st ITSL |  | 22 | 8 hr | YES |  |  |  |  |
| 126-86-3 | actylenic diol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 126-99-8 | beta-chloroprene | YES | 1st ITSL, Carc |  | 20 | annual* | YES |  |  |  | 0.002 |
| 127-18-4 | tetrachloroethylene | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 40 | annual | YES | 1400 | 24 hr |  | 4 |
| 127-91-3 | pinene, beta | YES | 1st ITSL |  | 1120 | 8 hr | YES |  |  |  |  |
| 128-04-1 | sodium dimethyl dithiocarbamate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 128-37-0 | 2,6-di-tert-butyl-p-cresol | YES | Carc |  |  |  |  |  |  |  | 1 |
| 129-00-0 | pyrene | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 131-11-3 | dimethylphthalate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 131-17-9 | diallyl phthalate | YES | Carc |  |  |  |  |  |  |  | 0.1 |
| 132-64-9 | dibenzofuran | YES | 1st ITSL |  | 4 | annual* | YES |  |  |  |  |
| 134-29-2 | o-ansidine hydrochloride | YES | Carc |  |  |  |  |  |  |  | 0.04 |
| 135-98-8 | sec-butylbenzene | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 136-47-0 | tetracaine hyrochloride | YES | 1st ITSL |  | 0.3 | annual | YES |  |  |  |  |
| 136-52-7 | cobalt 2-ethylhexanoate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 137-26-8 | thiram | YES | 1st ITSL |  | 17.5 | annual* | YES |  |  |  |  |
| 137-32-6 | 2-methyl-1-butanol | YES | 1st ITSL |  | 13 | annual | YES |  |  |  |  |
| 140-31-8 | aminoethylpiperazine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 140-88-5 | ethyl acrylate | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | $\begin{aligned} & \text { Future } \\ & \text { TAC? } \end{aligned}$ | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd <br> ITSL <br> AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 141-32-2 | butyl acrylate | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 141-43-5 | ethanolamine | YES | 1st ITSL |  | 80 | 8 hr | YES |  |  |  |  |
| 141-62-8 | decamethyltetrasiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 141-63-9 | linear dimethylsiloxanes,MD3M(\&higher) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 141-78-6 | ethyl acetate | NO |  | >75th\% | 3200 | annual* |  |  |  |  |  |
| 141-79-7 | mesityl oxide | YES | 1st ITSL |  | 400 | 8 hr | YES |  |  |  |  |
| 141-91-3 | 2,6-dimethyl morpholine | NO |  | >75th\% | 377 | annual |  |  |  |  |  |
| 141-97-9 | ethyl acetoacetate | YES | 1st ITSL |  | 46 | annual | YES |  |  |  |  |
| 142-29-0 | cyclopentene | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 142-59-6 | ethylene bisthiocarbamate disodium | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 142-71-2 | cupric acetate | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 142-82-5 | heptane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 142-84-7 | di-n-propylamine | YES | 1st ITSL |  | 1.5 | annual | YES |  |  |  |  |
| 142-96-1 | dibutyl ether | YES | 1st ITSL |  | 33 | annual | YES |  |  |  |  |
| 143-29-3 | butylcarbitol formal | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 144-62-7 | oxalic acid | YES | 1st ITSL |  | 10 | 8 hr | YES |  |  |  |  |
| 144-79-6 | diphenylmethylchlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 145-73-3 | endothall | YES | 1st ITSL |  | 35 | annual* | YES |  |  |  |  |
| 147-14-8 | copper phthalocyanine | YES | 1st ITSL |  | 21 | annual | YES |  |  |  |  |
| 147-24-0 | benadryl hcl | YES | 1st ITSL |  | 50 | annual | YES |  |  |  |  |
| 147-94-4 | cytarabine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 149-57-5 | 2-ethylhexanoic acid | YES | 1st ITSL |  | 64 | annual | YES |  |  |  |  |
| 149-73-5 | trimethylorthoformate | NO |  | >75th\% | 800 | annual |  |  |  |  |  |
| 151-56-4 | Ethylene imine (Aziridine) | YES | $\begin{gathered} \hline \text { HAP Table } \\ 2 \\ \hline \end{gathered}$ |  |  |  |  |  |  |  |  |
| 156-59-2 | cis-1-2,dichloroethylene | YES | 1st ITSL |  | 7 | annual* | YES |  |  |  |  |
| 156-60-5 | trans-1-2-dichloroethylene | YES | 1st ITSL |  | 70 | annual* | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th } \% \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \\ \hline \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 t h \% \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 189-55-9 | Dibenzo[a,i]pyrene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 189-64-0 | Dibenzo[a,h]pyrene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 191-24-2 | benzo(g,h,i)perylene | YES | 1st ITSL |  | 12 | annual* | YES |  |  |  |  |
| 191-26-4 | Anthanthrene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 191-30-0 | Dibenzo[a,l]pyrene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 192-65-4 | Dibenzo[a,e]pyrene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 193-09-9 | Naphtho[2,3e]pyrene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 193-39-5 | Indeno(1,2,3-cd)pyrene | YES | Carc7 |  |  |  |  |  |  |  |  |
| 199-54-2 | Benz[e]aceanthrylene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 202-33-5 | Benz[j]aceanthrylene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 202-94-8 | Benz[b,c]aceanthrylene, 11H | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 202-98-2 | Cyclopenta[d, e,f]chrysene, 4 H | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 205-12-9 | Benzo[c]fluorene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 205-82-3 | Benzo[j]fluoranthene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 205-99-2 | Benzo(b)fluoranthene | YES | Carc7 |  |  |  |  |  |  |  |  |
| 206-44-0 | fluoranthene | NO |  | >75th\% | 140 | annual* |  |  |  |  |  |
| 207-08-9 | Benzo(k)fluoranthene | YES | Carc7 |  |  |  |  |  |  |  |  |
| 208-96-8 | acenaphthylene | YES | 1st ITSL |  | 35 | annual* | YES |  |  |  |  |
| 211-91-6 | Benz[l]aceanthrylene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 215-58-7 | Dibenz[a,c]anthracene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 218-01-9 | Chrysene | YES | Carc7 |  |  |  |  |  |  |  |  |
| 280-57-9 | triethylenediamine | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 287-92-3 | cyclopentane | NO |  | >75th\% | 17200 | 8 hr |  |  |  |  |  |
| 300-57-2 | allyl benzene | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 302-01-2 | hydrazine | YES | Carc |  |  |  |  |  |  |  | 0.0002 |
| 302-22-7 | chlormadinone acetate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 303-81-1 | novobiocin | YES | 1st ITSL |  | 40 | annual* | YES |  |  |  |  |
| 309-00-2 | aldrin | YES | Carc |  |  |  |  |  |  |  | 0.0002 |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $1 s t$ ITSL $\left(\mu \mathrm{g} / \mathrm{m}^{3}\right)$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 313-06-4 | estradiol cypionate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 319-84-6 | alpha-hexachlorocyclohexane | YES | Carc |  |  |  |  |  |  |  | 0.0006 |
| 330-54-1 | diuron | YES | 1st ITSL |  | 7 | annual* | YES |  |  |  |  |
| 335-67-1 | PFOA | YES | Emerging |  |  |  |  |  |  |  |  |
| 338-98-7 | isoflupredone acetate | YES | 1st ITSL |  | 0.01 | annual | YES |  |  |  |  |
| 353-50-4 | carbonyl fluoride | YES | 1st ITSL |  | 54 | 8 hr | YES |  |  |  |  |
| 358-67-8 | trifluoropropylmethyl dimethoxysilane | YES | 1st ITSL |  | 100 | annual | YES |  |  |  |  |
| 359-07-9 | 2-bromo-1,1-difluoro ethane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 363-51-9 | 2-chloro-6-fluorobenzenamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 363-72-4 | pentafluorobenzene | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 366-18-7 | 2,2'-bipyridyl | YES | 1st ITSL |  | 0.8 | annual | YES |  |  |  |  |
| 382-21-8 | perfluoroisobutylene | YES | 1st ITSL |  | 0.8 | 1 hr | YES |  |  |  |  |
| 385-00-2 | 2,6-difluorobenzoic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 431-89-0 | hfc-227ea | NO |  | >75th\% | 130000 | annual |  | 5560000 | 1 hr |  |  |
| 460-73-1 | 1,1,1,3,3-pentafluoropropane | NO |  | >75th\% | 2000 | annual* |  |  |  |  |  |
| 461-58-5 | cyanoguanidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 463-58-1 | carbonyl sulfide | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 505-48-6 | suberic acid | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 509-14-8 | tetranitromethane | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 0.4 | 8 hr | YES |  |  |  | 7E-05 |
| 513-35-9 | amylene | NO |  | >75th\% | 106 | annual |  |  |  |  |  |
| 513-37-1 | dimethylvinyl chloride | YES | Carc |  |  |  |  |  |  |  | 0.008 |
| 513-85-9 | 2,3-butanediol | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 526-73-8 | 1,2,3-trimethylbenzene | YES | 1st ITSL, 2nd ITSL |  | 50 | annual | YES | 1200 | 8 hr | YES |  |
| 526-75-0 | 2,3-dimethyl phenol | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 532-27-4 | alpha chloroacetophenone | YES | 1st ITSL |  | 0.03 | annual* | YES |  |  |  |  |
| 534-52-1 | dinitro-o-cresol | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 540-49-8 | 1,2-dibromoethylene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 540-59-0 | 1,2-dichloroethylene | YES | 1st ITSL |  | 35 | annual* | YES |  |  |  |  |
| 540-84-1 | 2,2,4-Trimethyl Pentane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 540-88-5 | tert-butyl acetate | NO |  | >75th\% | 9500 | 8 hr |  |  |  |  |  |
| 540-97-6 | dodecamethylcyclohexasiloxane | NO |  | >75th\% | 400 | annual |  |  |  |  |  |
| 541-02-6 | decamethylcyclopentasiloxane | NO |  | >75th\% | 200 | annual* |  |  |  |  |  |
| 541-05-9 | hexamethylcyclotrisiloxane | YES | 1st ITSL |  | 50 | annual | YES |  |  |  |  |
| 541-73-1 | 1,3-dichlorobenzene | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 541-85-5 | ethyl amyl ketone | NO |  | >75th\% | 220 | annual* |  |  |  |  |  |
| 542-75-6 | 1,3-dichloropropene | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 20 | annual* | YES |  |  |  | 0.2 |
| 542-88-1 | bis(chloromethyl)ether | YES | Carc |  |  |  |  |  |  |  | 2E-05 |
| 546-93-0 | magnesium carbonate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 552-45-4 | alpha-chloro-ortho-xylene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 556-67-2 | octamethylcyclotetrasilo | YES | 1st ITSL |  | 75 | annual* | YES |  |  |  |  |
| 557-04-0 | magnesium stearate | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 557-05-1 | zinc stearate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 563-47-3 | 3-chloro-2-methylpropene | YES | Carc |  |  |  |  |  |  |  | 0.03 |
| 565-59-3 | 2,3-dimethylpentane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 576-26-1 | 2,6-dimethyl phenol | YES | 1st ITSL |  | 2 | annual* | YES |  |  |  |  |
| 584-84-9 | 2,4-toluene diisocyanate | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 0.07 | annual* | YES |  |  |  | 0.03 |
| 589-34-4 | 3-methylhexane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 590-01-2 | n-butyl propionate | NO |  | >75th\% | 102 | annual |  |  |  |  |  |
| 590-86-3 | isovaleraldehyde | NO |  | >75th\% | 800 | annual |  |  |  |  |  |
| 591-22-0 | 3,5-lutidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 591-27-5 | m-aminophenol | NO |  | >75th\% | 390 | annual |  |  |  |  |  |
| 591-76-4 | 2-methylhexane | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | 2nd <br> ITSL <br> $<=75 t h \%$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 591-78-6 | methyl n-butyl ketone | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 592-09-6 | trifluoropropyltrichlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 592-42-7 | 1,5-hexanediene | NO |  | >75th\% | 264 | annual |  |  |  |  |  |
| 592-76-7 | 1-heptene | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 592-84-7 | butyl formate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 593-51-1 | methylamine hydrochloride | YES | 1st ITSL |  | 64 | 8 hr | YES |  |  |  |  |
| 593-60-2 | vinyl bromide | YES | 1st ITSL |  | 3 | annual* | YES |  |  |  |  |
| 606-46-2 | n,n-diethyl-o-toluine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 608-31-1 | 2,6-dichlorobenzenamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 611-14-3 | 1-ethyl-2-methylbenzene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 612-00-0 | 1,1-diphenylethane | YES | 1st ITSL |  | 0.8 | annual | YES |  |  |  |  |
| 613-48-9 | n,n-diethyl-p-toludine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 616-38-6 | dimethyl carbonate | NO |  | >75th\% | 300 | annual |  |  |  |  |  |
| 617-94-7 | ```phenyl isopropanol (2-phenyl-2- propanol)``` | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 620-23-5 | m-tolualdehyde | NO |  | >75th\% | 440 | annual* |  |  |  |  |  |
| 621-64-7 | n-nitroso-di-n-propylamine | YES | Carc |  |  |  |  |  |  |  | 0.0005 |
| 621-77-2 | triamylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 622-96-8 | p-ethyl toluene | NO |  | >75th\% | 350 | annual* |  |  |  |  |  |
| 622-97-9 | 4-methylstyrene | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 624-41-9 | 2-methyl butyl acetate | NO |  | >75th\% | 1100 | annual* |  |  |  |  |  |
| 624-54-4 | n-pentyl proprionate | YES | 1st ITSL |  | 21 | annual | YES |  |  |  |  |
| 624-83-9 | Methyl isocyanate | YES | ATW |  |  |  |  |  |  |  |  |
| 624-92-0 | dimethyl disulfide | YES | 1st ITSL |  | 28 | annual | YES |  |  |  |  |
| 626-38-0 | 2-pentyl acetate | NO |  | >75th\% | 2600 | 8 hr |  |  |  |  |  |
| 626-67-5 | n-methylpiperidine | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 627-20-3 | cis-2-pentene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 627-30-5 | 3-chloro-1-propanol | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{array}{\|c} \text { 1st ITSL } \\ <=75 \text { th } \% \end{array}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 627-83-8 | octadecanoic acid, 1,2-ethanediyl ester | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 628-63-7 | n -amyl acetate | NO |  | >75th\% | 1100 | annual* |  |  |  |  |  |
| 629-11-8 | 1,6-hexanediol | YES | 1st ITSL |  | 14 | annual | YES |  |  |  |  |
| 629-73-2 | 1-hexadecene | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 630-20-6 | 1,1,1,2-tetrachloroethane | YES | Carc |  |  |  |  |  |  |  | 0.1 |
| 630-93-3 | sodium dilantin | YES | Carc |  |  |  |  |  |  |  | 0.04 |
| 632-22-4 | tetramethyl urea | YES | 1st ITSL, <br> 2nd ITSL |  | 0.8 | annual | YES | 230 | 24 hr | YES |  |
| 634-66-2 | 1,2,3,4-tetrachlorobenzene | NO |  | >75th\% | 120 | annual* |  |  |  |  |  |
| 634-90-2 | 1,2,3,5-tetrachlorobenzene | YES | 1st ITSL |  | 12 | annual* | YES |  |  |  |  |
| 637-92-3 | ethyl tertiary butyl ether | NO |  | >75th\% | 373 | annual* |  |  |  |  |  |
| 644-62-2 | meclofenamic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 646-06-0 | 1,3-dioxolane | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 668-45-1 | chlorofluorobenzonitrile | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 677-21-4 | 3,3,3-trifluoropropene | NO |  | >75th\% | 280 | annual |  |  |  |  |  |
| 684-93-5 | n-nitroso-n-methylurea | YES | Carc |  |  |  |  |  |  |  | 2E-06 |
| 694-87-1 | benzocyclobutene | NO |  | >75th\% | 220 | annual |  |  |  |  |  |
| 696-82-2 | 2,4,6-trifluoropyrimidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 701-64-4 | monophenyl phosphoric acid | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 756-79-6 | dimethyl methyl phosphonate | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 763-69-9 | ethyl-3-ethyloxypropionate | NO |  | >75th\% | 134 | annual* |  |  |  |  |  |
| 770-35-4 | propylene glycol phenyl ether | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 778-25-6 | diphenylmethylsilanol | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 807-28-3 | tetraphenyldimethyldisiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 811-97-2 | 1,1,1,2-tetrafluoroethane | NO |  | >75th\% | 80000 | annual* |  |  |  |  |  |
| 814-68-6 | acryloyl chloride | YES | 1st ITSL |  | 0.3 | annual | YES |  |  |  |  |
| 822-06-0 | hexamethylene diisocyanate | YES | 1st ITSL |  | 0.01 | annual* | YES |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 827-52-1 | cyclohexylbenzene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 836-30-6 | n-nitrodiphenylamine | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 838-85-7 | diphenyl phosphoric acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 859-18-7 | lincomycin hydrochloride | YES | 1st ITSL |  | 75 | annual* | YES |  |  |  |  |
| 868-77-9 | 2-hydroxyethyl methacrylate | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 872-36-6 | vinylene carbonate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 872-50-4 | N-methylpyrrolidone | NO |  | >75th\% | 700 | annual* |  |  |  |  |  |
| 947-19-3 | 1-hydroxcyclohexyl phenyl ketone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 981-34-0 | betamethasone 11 | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 992-94-9 | methylsilane | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 993-07-7 | trimethylsilane | NO |  | >75th\% | 340 | annual |  |  |  |  |  |
| 994-05-8 | tertiary amyl methyl ether | YES | 1st ITSL |  | 62 | annual* | YES |  |  |  |  |
| 996-35-0 | dimethylisopropylamine | NO |  | >75th\% | 200 | annual |  |  |  |  |  |
| 999-97-3 | hexamethyldisilazane | NO |  | >75th\% | 206 | annual |  |  |  |  |  |
| 1009-93-4 | hexamethylcyclotrisilazane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1047-16-1 | quinacridone pigment | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1066-35-9 | dimethylchlorosilane | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 1066-40-6 | trimethylsilanol | YES | 1st ITSL |  | 65 | annual | YES |  |  |  |  |
| 1067-25-0 | propyltrimethoxysilane | NO |  | >75th\% | 1000 | annual |  |  |  |  |  |
| 1070-10-6 | 2-ethylhexyltitanate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1072-53-3 | 1,3,2-dioxathiolane,2,2-dioxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1072-63-5 | 1-vinylimidazol | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 1074-40-4 | 4,6-dichloro-2-methoxypyrimidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1111-74-6 | dimethylsilane | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 1112-39-6 | dimethyldimethoxysilane | YES | 1st ITSL |  | 90 | annual | YES |  |  |  |  |
| 1116-54-7 | n-nitrosodiethanolamine | YES | Carc |  |  |  |  |  |  |  | 0.0012 |
| 1120-71-4 | 1,3-propane sultone | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 1122-82-3 | cyclohexyl isothiocyanate | NO |  | default | 0.1 | annual |  |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1156-19-0 | tolazamide | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 1163-19-5 | decabromodiphenyl oxide | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 25 | annual* | YES |  |  |  | 5 |
| 1184-85-6 | methyl methane sulfonamide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1185-55-3 | trimethoxymethylsilane | YES | 1st ITSL |  | 80 | annual | YES |  |  |  |  |
| 1194-02-1 | p-fluorobenzonitrile | YES | 1st ITSL |  | 0.5 | annual | YES |  |  |  |  |
| 1300-72-7 | sodium xylenesulfonate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1306-38-3 | cerium oxide | YES | 1st ITSL |  | 0.9 | annual* | YES |  |  |  |  |
| 1308-14-1 | chromium (+3) hydroxide | YES | 1st ITSL |  | 0.5 | annual* | YES |  |  |  |  |
| 1308-38-9 | chromium 3 oxide | YES | 1st ITSL |  | 0.5 | annual* | YES |  |  |  |  |
| 1309-42-8 | magnesium hydroxide | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 1309-48-4 | magnesium oxide | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 1309-64-4 | antimony trioxide | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |
| 1310-53-8 | germanium dioxide | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 1310-58-3 | potassium hydroxide | YES | 1st ITSL |  | 20 | 1 hr | YES |  |  |  |  |
| 1310-66-3 | lithium hydroxide | YES | 1st ITSL |  | 0.25 | 8 hr | YES |  |  |  |  |
| 1310-73-2 | sodium hydroxide | YES | 1st ITSL |  | 20 | 1 hr | YES |  |  |  |  |
| 1313-27-5 | molybdenum trioxide | YES | 1st ITSL, Carc |  | 5 | 8 hr | YES |  |  |  | 0.12 |
| 1313-96-8 | niobium oxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1314-13-2 | zinc oxide | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 1314-28-9 | rhenium oxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1314-32-5 | thallic oxide | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |
| 1314-62-1 | vanadium pentoxide | YES | 1st ITSL |  | 0.5 | 1 hr | YES |  |  |  |  |
| 1317-33-5 | molybdenum disulfide | YES | 1st ITSL |  | 30 | 8 hr | YES |  |  |  |  |
| 1319-77-3 | cresol (mixed isomers) | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 1320-67-8 | propylene glycol monomethyl ether | NO |  | >75th\% | 2000 | annual* |  |  |  |  |  |
| 1328-53-6 | phthalocyanine pigment green | NO |  | default | 0.1 | annual |  |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1330-20-7 | mixed xylenes | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 1330-86-5 | adipate plasticizer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1332-21-4 | asbestos | YES | Carc |  |  |  |  |  |  |  | 2E-05 |
| 1332-58-7 | kaolin | YES | 1st ITSL |  | 20 | 8 hr | YES |  |  |  |  |
| 1333-13-7 | tert-butyl-m-cresol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1333-86-4 | carbon black | YES | 1st ITSL |  | 30 | 8 hr | YES |  |  |  |  |
| 1336-21-6 | ammonium hydroxide | NO |  | >75th\% | 200 | annual* |  |  |  |  |  |
| 1336-36-3 | polychlorinated biphenyls | YES | Carc |  |  |  |  |  |  |  | 0.002 |
| 1338-23-4 | methyl ethyl ketone peroxide | YES | 1st ITSL |  | 15 | 1 hr | YES |  |  |  |  |
| 1345-04-6 | antimony trisulfide | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |
| 1345-05-7 | lithopone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1405-10-3 | neomycin sulfate | NO |  | >75th\% | 280 | annual |  |  |  |  |  |
| 1445-45-0 | trimethyl-o-acetate | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 1477-55-0 | 1,3-bis(aminomethyl)benzenen | YES | 1st ITSL |  | 1 | 1 hr | YES |  |  |  |  |
| 1559-35-9 | ethylene glycol mono-2-ethylhexyl ether | YES | 1st ITSL |  | 37 | annual | YES |  |  |  |  |
| 1559-36-0 | diethylene glycol mono-2-ethylhexyl ether | YES | 1st ITSL |  | 22 | annual | YES |  |  |  |  |
| 1559-37-1 | triethylene glycol mono-2-ethyhexyl ether | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1569-01-3 | 1-propoxy-2-propanol | YES | 1st ITSL |  | 86 | annual* | YES |  |  |  |  |
| 1569-02-4 | propylene glycol monoethyl ether (beta) | NO |  | >75th\% | 240 | annual |  |  |  |  |  |
| 1589-47-5 | 2-methoxy-1-propanol | NO |  | >75th\% | 660 | annual* |  |  |  |  |  |
| 1590-87-0 | disilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1623-15-0 | monobutyl phosphoric acid | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 1634-04-4 | methyl t-butyl ether | NO |  | >75th\% | 3000 | annual* |  |  |  |  |  |
| 1643-19-2 | t-n-butyl ammonium bromide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1702-17-6 | clopyralid | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1717-00-6 | 1,1-dichloro-1-fluoroethane | NO |  | >75th\% | 12800 | annual* |  |  |  |  |  |
| 1719-58-0 | dimethylvinylchlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1746-01-6 | 2,3,7,8-tetrachlorodibenzo(p)dioxin | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 2E-06 | annual | YES |  |  |  | 2E-08 |
| 1758-88-9 | 2-ethyl-1,4-dimethyl benzene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1760-24-3 | n-(3-(trimethoxysilyl)propyl)ethylenediamine | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 1761-71-3 | 4,4'-diaminodicyclohexylmethane | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 1763-23-1 | PFOS | YES | Emerging |  |  |  |  |  |  |  |  |
| 1873-88-7 | heptamethyltrisiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 1897-52-5 | 2,6-difluorobenzonitrile | YES | 1st ITSL |  | 0.2 | annual | YES |  |  |  |  |
| 1912-83-0 | stannous octoate | YES | 1st ITSL |  | 1 | 8 hr | YES |  |  |  |  |
| 2031-67-6 | methyltriethoxysilane | YES | 1st ITSL |  | 23 | annual | YES |  |  |  |  |
| 2050-92-2 | diamylamine | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 2157-45-1 | tetra-2-methoxyethoxy-silane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2160-93-2 | t-butyldiethanolamine | YES | 1st ITSL |  | 9 | annual | YES |  |  |  |  |
| 2238-07-5 | diglycidyl ether | YES | 1st ITSL |  | 0.5 | 8 hr | YES |  |  |  |  |
| 2370-88-9 | cyclic methylhydrogensiloxane, d4 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2374-14-3 | cyclic methyltrifluoropropylsiloxane, $\mathrm{d} 3$ | YES | 1st ITSL |  | 0.6 | annual | YES |  |  |  |  |
| 2403-89-6 | 1,2,2,6,6-pentamethyl-4-piperidinol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2426-08-6 | n-butyl glycidyl ether | YES | 1st ITSL, <br> 2nd ITSL |  | 300 | 1 hr | YES | 160 | 8 hr | YES |  |
| 2467-02-9 | bisphenolf | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2476-74-6 | flumethasone 6 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2530-85-0 | organofunctional silane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2627-86-3 | L-alpha-phenylethylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2627-95-4 | tetramethyldivinyldisiloxane | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 2627-97-6 | dimethyldiphenydivinylsiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 2682-20-4 | 2-methyl-4-isothiazolin-3-one | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 2687-91-4 | 1-ethyl-2-pyrrolidone | YES | 1st ITSL |  | 4.9 | annual | YES |  |  |  |  |
| 2768-02-7 | vinyltrimethoxysilane | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 2807-30-9 | ethylene glycol monopropyl ether | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 2837-89-0 | 2-chloro-1,1,1,2-tetrafluoroethane | NO |  | >75th\% | 5000 | annual* |  |  |  |  |  |
| 2919-66-6 | melengesterol acetate | YES | 1st ITSL |  | 2 | annual* | YES |  |  |  |  |
| 2981-10-4 | piperdinocyclohexene | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 2996-92-1 | phenyltrimethoxysilane | YES | 1st ITSL |  | 60 | annual | YES |  |  |  |  |
| 3006-82-4 | t-butylperoxy-2-ethylhexanoate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3006-86-8 | 1,1-di-(tert-buytlperoxy)cyclohexane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3020-12-0 | $\begin{aligned} & \text { o-(1-ethoxyethyl)-3- } \\ & \text { (trifluoromethyl)phenol } \end{aligned}$ | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3033-62-3 | bis (2-dimethylaminoethyl) ether | YES | 1st ITSL |  | 0.05 | annual* | YES |  |  |  |  |
| 3052-70-8 | (1-methylethylidene)bis(1,1dimethylpropyl)peroxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3081-01-4 | santoflex 14 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3144-09-0 | methanesulfonamide | YES | 1st ITSL |  | 44 | annual | YES |  |  |  |  |
| 3153-26-2 | vanadium oxide bis (2,4pentanedionate) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3236-53-1 | trimethyl hexamethylenediamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3277-26-7 | tetramethyldihydrogendisiloxane | NO |  | >75th\% | 120 | annual |  |  |  |  |  |
| 3290-92-4 | trimethylolpropane trimethacrylate | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 3390-61-2 | tetraphenyldimethyl-2phenylmethyltrisiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3399-73-3 | cyclohexenylethylamine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3697-24-3 | 5-methylchrysene | YES | Cal Carc |  |  |  |  |  |  |  |  |
| 3731-51-9 | 2-(aminomethyl)pyridine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3764-01-0 | 2,4,6-trichloropyrimidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3779-63-3 | aliphatic polyisocyanate-1 | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. |  | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th } \% \\ & \hline \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 3814-34-4 | 2-ethylbutyl bromide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3843-16-1 | distearyldimethylammonium methosulfate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3982-82-9 | tetraphenyldimethyl-2dimethyltrisiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 3986-89-8 | progesterone 4 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 4098-71-9 | isophorone diisocyanate | YES | 1st ITSL |  | 0.45 | 8 hr | YES |  |  |  |  |
| 4109-96-0 | dichlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 4170-30-3 | crotonaldehyde | YES | 1st ITSL |  | 9 | 1 hr | YES |  |  |  |  |
| 4221-98-1 | p-mentha-1,5-diene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 4253-34-3 | methyltriacetoxysilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 4420-74-0 | 3-mercaptopropyltrimethoxysilane | YES | 1st ITSL |  | 2.4 | annual | YES |  |  |  |  |
| 4435-53-4 | butoxyl | YES | 1st ITSL |  | 14 | annual | YES |  |  |  |  |
| 4444-67-1 | di-sec-butylamine | NO |  | >75th\% | 417 | annual |  |  |  |  |  |
| 4620-70-6 | t-butylaminoethanol | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 4652-27-1 | 4-methoxy-3-buten-2-one | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 4994-16-5 | 4-phenylcyclohexene | YES | 1st ITSL |  | 33 | annual | YES |  |  |  |  |
| 5131-66-8 | propylene glycol n-butyl ether (alpha isomer) | YES | 1st ITSL |  | 77 | annual | YES |  |  |  |  |
| 5314-55-6 | ethyltrimethoxysilane | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 5329-14-6 | sulfamic acid | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 5385-75-1 | Dibenzo[a,e]fluoranthene | YES | EPA Carc |  |  |  |  |  |  |  |  |
| 5436-21-5 | 4,4-dimethoxy-2-butanone | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 5507-44-8 | vinylmethyldiethoxysilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 5509-65-9 | 2,6-difluoroaniline | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 5779-94-2 | 2,5-dimethylbenzaldehyde | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 5888-33-5 | iso-bornyl acrylate | YES | 1st ITSL |  | 14 | annual | YES |  |  |  |  |
| 5906-75-2 | vinyl dimethylsilanol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 5989-27-5 | d-limonene | NO |  | >75th\% | 6250 | annual* |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 6004-24-6 | cetylpyridinium chloride monohydrate | YES | 1st ITSL |  | 1.8 | annual | YES |  |  |  |  |
| 6166-86-5 | cyclic methylhydrogensiloxane, d5 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6192-52-5 | p-toluenesulfonic acid monohydrate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6408-78-2 | C.I. acid blue 25 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6419-19-8 | aminotrimethylene phosphonic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6574-99-8 | 3,4-dichlorobenzonitrile | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 6674-22-2 | 1,8-diazabicyclo[5.4.0]undec-7-ene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6700-34-1 | dextromethorphan hydrochloride | YES | 1st ITSL |  | 0.4 | annual | YES |  |  |  |  |
| 6713-03-7 | 1-(2-hydroxyethylthio)propane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6846-50-0 | 2,2,4-trimethylpentanediol-1,3diisobutyrate | NO |  | >75th\% | 106 | annual |  |  |  |  |  |
| 6904-66-1 | tetraphenylhexamethyltetrasiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 6915-15-7 | malic acid | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 6975-71-9 | cyclohexenylacetonitrile | YES | 1st ITSL |  | 3.5 | annual | YES |  |  |  |  |
| 7085-85-0 | ethyl 2-cyanoacrylate | YES | 1st ITSL |  | 10 | 8 hr | YES |  |  |  |  |
| 7439-93-2 | lithium | YES | 1st ITSL |  | 35 | annual* | YES |  |  |  |  |
| 7439-95-4 | magnesium | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 7439-96-5 | manganese and compounds | YES | 1st ITSL |  | 0.05 | annual | YES |  |  |  |  |
| 7439-97-6 | Mercury | YES | < $75 \%$ |  |  |  |  |  |  |  |  |
| 7439-98-7 | molybdenum | YES | 1st ITSL |  | 30 | 8 hr | YES |  |  |  |  |
| 7440-02-0 | nickel | YES | Carc |  |  |  |  |  |  |  | 0.0042 |
| 7440-05-3 | palladium | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 7440-06-4 | platinum soluble salt | YES | 1st ITSL |  | 0.02 | 8 hr | YES |  |  |  |  |
| 7440-22-4 | silver - soluble | YES | 1st ITSL |  | 0.1 | 8 hr | YES |  |  |  |  |
| 7440-24-6 | strontium | NO |  | >75th\% | 2000 | annual* |  |  |  |  |  |
| 7440-28-0 | thallium | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |
| 7440-31-5 | tin | YES | 1st ITSL |  | 20 | 8 hr | YES |  |  |  |  |
| 7440-36-0 | antimony | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. |  | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ |  | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 7440-38-2 | arsenic | YES | Carc |  |  |  |  |  |  |  | 0.0002 |
| 7440-39-3 | barium | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 7440-41-7 | beryllium | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \\ \hline \end{gathered}$ |  | 0.02 | annual* | YES |  |  |  | 0.0004 |
| 7440-43-9 | cadmium | YES | Carc |  |  |  |  |  |  |  | 0.0006 |
| 7440-45-1 | cerium | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 7440-48-4 | cobalt | YES | 1st ITSL |  | 0.2 | 8 hr | YES |  |  |  |  |
| 7440-50-8 | copper | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 7440-65-5 | yttrium | YES | 1st ITSL |  | 10 | 8 hr | YES |  |  |  |  |
| 7446-11-9 | sulfur trioxide | YES | 1st ITSL, 2nd ITSL |  | 1 | annual | YES | 120 | 1 hr | YES |  |
| 7446-70-0 | aluminum chloride | YES | 1st ITSL |  | 20 | 8 hr | YES |  |  |  |  |
| 7473-98-5 | 2-hydroxy-2-methyl-1-phenyl-1propanone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 7525-62-4 | ethylvinyl benzene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 7553-56-2 | iodine | YES | 1st ITSL |  | 1 | 8 hr | YES |  |  |  |  |
| 7558-79-4 | disodium hydrogen phosphate | YES | 1st ITSL |  | 10 | annual* | YES |  |  |  |  |
| 7580-85-0 | 2-tert-butoxyethanol | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 7631-90-5 | sodium bisulfite | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 7631-95-0 | sodium molybdate | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 7632-00-0 | sodium nitrite | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 7632-04-4 | sodium perborate | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 7637-07-2 | boron trifluoride | YES | 1st ITSL |  | 0.7 | annual* | YES |  |  |  |  |
| 7647-01-0 | hydrogen chloride | YES | 1st ITSL |  | 20 | annual | YES | 2100 | 1 hr |  |  |
| 7647-15-6 | sodium bromide | NO |  | >75th\% | 140 | annual* |  |  |  |  |  |
| 7664-38-2 | phosphoric acid | YES | 1st ITSL |  | 10 | annual* | YES |  |  |  |  |
| 7664-39-3 | hydrogen fluoride | YES | 1st ITSL, 2nd ITSL |  | 14 | annual | YES | 240 | 1 hr | YES |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. |  | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | 2nd <br> ITSL <br> $<=75$ th $\%$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 7664-41-7 | ammonia | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 7664-93-9 | sulfuric acid | YES | 1st ITSL, 2nd ITSL |  | 1 | annual | YES | 120 | 1 hr | YES |  |
| 7681-49-4 | sodium fluoride | YES | 1st ITSL |  | 60 | 8 hr | YES |  |  |  |  |
| 7681-52-9 | sodium hypochlorite | YES | 1st ITSL |  | 16 | 8 hr | YES |  |  |  |  |
| 7681-82-5 | sodium iodide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 7691-02-3 | tetramethyldivinyldisila | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 7697-37-2 | nitric acid | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 7704-34-9 | sulfur (elemental) | YES | 1st ITSL |  | 30 | 8 hr | YES |  |  |  |  |
| 7722-64-7 | potassium permanganate | YES | 1st ITSL |  | 0.1 | annual* | YES |  |  |  |  |
| 7722-76-1 | ammonium dihydrogen phosphate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 7722-84-1 | hydrogen peroxide | YES | 1st ITSL |  | 14 | 8 hr | YES |  |  |  |  |
| 7723-14-0 | phosphorus (total) | YES | 1st ITSL |  | 1 | 8 hr | YES |  |  |  |  |
| 7726-95-6 | bromine | YES | 1st ITSL |  | 7 | 8 hr | YES |  |  |  |  |
| 7727-43-7 | barium sulfate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 7757-83-7 | sodium sulfite | YES | 1st ITSL |  | 0.028 | annual | YES |  |  |  |  |
| 7758-05-6 | potassium iodate | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 7758-98-7 | copper sulfate, anhydrous | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 7758-99-8 | copper sulfate pentahydrate | YES | 1st ITSL |  | 10 | 8 hr | YES |  |  |  |  |
| 7782-49-2 | selenium | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 7782-50-5 | chlorine | YES | 1st ITSL, 2nd ITSL |  | 0.3 | annual | YES | 500 | 8 hr | YES |  |
| 7782-65-2 | germanium tetrahydride | YES | 1st ITSL |  | 6 | 8 hr | YES |  |  |  |  |
| 7783-06-4 | hydrogen sulfide | YES | 1st ITSL, 2nd ITSL |  | 2 | annual | YES | 100 | 24 hr | YES |  |
| 7783-28-0 | diammonium hydrogen phosphate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 7783-54-2 | nitrogen trifluoride | YES | 1st ITSL |  | 290 | 8 hr | YES |  |  |  |  |
| 7783-61-1 | silicon tetrafluoride | YES | 1st ITSL |  | 0.2 | annual | YES |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 7784-42-1 | arsine | YES | 1st ITSL |  | 0.05 | annual* | YES |  |  |  |  |
| 7786-30-3 | magnesium chloride | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 7789-23-3 | potassium fluoride | YES | 1st ITSL |  | 76 | 8 hr | YES |  |  |  |  |
| 7789-82-4 | calcium molybdate | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 7803-51-2 | phosphine | YES | 1st ITSL |  | 0.3 | annual* | YES |  |  |  |  |
| 7803-52-3 | stibine | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 7803-62-5 | silicon tetrahydride | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 8001-35-2 | toxaphene | YES | Carc |  |  |  |  |  |  |  | 0.003 |
| 8001-79-4 | castor oil | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 8002-09-3 | yarmor pine oil | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 8002-74-2 | paraffin wax fume | YES | 1st ITSL |  | 20 | 8 hr | YES |  |  |  |  |
| 8005-02-5 | solvent black | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 8006-61-9 | gasoline | YES | Carc |  |  |  |  |  |  |  | 2 |
| 8006-64-2 | turpentine | YES | 1st ITSL |  | 1120 | 8 hr | YES |  |  |  |  |
| 8007-45-2 | coke oven emissions | YES | Carc |  |  |  |  |  |  |  | 0.0016 |
| 8012-95-1 | mineral oil | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 8014-95-7 | oleum | YES | 1st ITSL, 2nd ITSL |  | 1 | annual | YES | 120 | 1 hr | YES |  |
| 8020-83-5 | deodorized kerosene | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 8030-30-6 | naphtha | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 8032-32-4 | VM \& P naphtha | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 8042-47-5 | white mineral oil | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 8050-09-7 | colophony | YES | 1st ITSL |  | 1 | 1 hr | YES |  |  |  |  |
| 8052-41-3 | stoddard solvent | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 8052-42-4 | Asphalt fumes | YES | Carc |  |  |  |  |  |  |  |  |
| 9000-90-2 | alpha-amylase | YES | 1st ITSL |  | 0.02 | 1 hr | YES |  |  |  |  |
| 9001-92-7 | bacillus subtilis neutral protease | YES | 1st ITSL |  | 0.02 | 1 hr | YES |  |  |  |  |
| 9002-86-2 | polyvinyl chloride | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |

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| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 9002-92-0 | polyoxyethylene lauryl ether | YES | 1st ITSL |  | 12 | annual | YES |  |  |  |  |
| 9002-93-1 | triton $\times 100$ | YES | 1st ITSL |  | 0.15 | annual | YES |  |  |  |  |
| 9003-11-6 | methyl oxirane (pluronic p103) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 9003-13-8 | polyalkylene glycol monobutyl ether/ butoxypolypropylene glycol | NO |  | >75th\% | 160 | annual |  |  |  |  |  |
| 9003-22-9 | polyvinylchloride/polyvinylacetate | YES | 1st ITSL |  | 50 | annual | YES |  |  |  |  |
| 9003-39-8 | polyvinyl pyrrolidone | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 9003-55-8 | styrene-butadiene polymer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 9004-32-4 | carboxymethyl cellulose | NO |  | >75th\% | 300 | annual |  |  |  |  |  |
| 9004-58-4 | ethylhydroxyethyl cellulose | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 9004-74-4 | polyethylene glycol methyl ether | YES | 1st ITSL |  | 13 | annual | YES |  |  |  |  |
| 9011-17-0 | polyvinylidine fluoride | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 9014-85-1 | tetramethyl decyndiol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 9014-92-0 | t-det dd-14 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 9016-45-9 | igepal co-630 | YES | 1st ITSL |  | 18 | annual | YES |  |  |  |  |
| 9016-87-9 | polmeric methylene diphenyl diisocyanate | YES | 1st ITSL |  | 0.6 | annual* | YES |  |  |  |  |
| 9036-19-5 | t-det c08 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 9063-06-3 | oxirane, methyl-, polymer with oxirane, monomethyl ether | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 10025-78-2 | trichlorosilane | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 10025-91-9 | antimony trichloride | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 10026-04-7 | silicon tetrachloride | NO |  | >75th\% | 1100 | annual |  |  |  |  |  |
| 10034-93-2 | hydrazine sulfate | YES | Carc |  |  |  |  |  |  |  | 0.0008 |
| 10034-96-5 | manganese sulfate monohydrate | YES | 1st ITSL |  | 0.15 | annual* | YES |  |  |  |  |
| 10035-10-6 | hydrogen bromide | YES | 1st ITSL |  | 70 | 1 hr | YES |  |  |  |  |
| 10039-56-2 | sodium hypophosphite monohydrate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 10049-04-4 | chlorine dioxide | YES | 1st ITSL |  | 0.2 | annual* | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{gathered} \text { 1st ITSL } \\ <=75 \text { th\% } \end{gathered}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | $\begin{aligned} & \text { 2nd } \\ & \text { ITSL } \\ & \text { AvgT } \\ & \hline \end{aligned}$ | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 10096-91-0 | hydroxyphenylbenzotriazole | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 10097-09-3 | bis-urea accelerator | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 10190-55-3 | lead molybdate | YES | 1st ITSL |  | 30 | 8 hr | YES |  |  |  |  |
| 10215-30-2 | 2-propoxy-1-propanol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 10377-60-3 | magnesium nitrate | YES | 1st ITSL |  | 100 | 8 hr | YES |  |  |  |  |
| 10431-98-8 | 2-ethyl-2-oxazoline | YES | 1st ITSL |  | 53 | annual | YES |  |  |  |  |
| 10469-09-7 | tetrachloropicolinic acid | YES | 1st ITSL |  | 21 | annual | YES |  |  |  |  |
| 10482-56-1 | alpha-terpineol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 10551-21-0 | phenethyl alpha picolinium bromide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 12021-95-3 | hexafluorozirconium acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 12035-72-2 | nickel subsulfide | YES | Carc |  |  |  |  |  |  |  | 0.0021 |
| 12037-29-5 | praseodymium oxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 12054-85-2 | ammonium molybdate | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 12070-12-1 | tungsten carbide | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 12136-45-7 | potassium oxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 12262-58-7 | cyclohexanone peroxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 12401-86-4 | sodium monoxide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 12789-03-6 | chlordane (technical) | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 0.7 | annual* | YES |  |  |  | 0.01 |
| 13007-85-7 | sodium glucoheptonate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 13209-41-1 | 17,21-dihydroxy-16 alpha-methylpregna-1,4,9(11)-triene-3,20dione | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 13465-77-5 | hexachlorodisilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 13466-78-9 | carene, delta | YES | 1st ITSL |  | 1120 | 8 hr | YES |  |  |  |  |
| 13528-93-3 | bis(me2clsilyl)ethane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 13701-59-2 | barium metaborate monohydrate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 13879-32-8 | 1,1'[methylenebis(oxyethane-1,2diloxy)]bisbenzene | YES | 1st ITSL |  | 0.7 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd <br> ITSL <br> AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 13952-84-6 | sec-butylamine | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 14579-03-4 | cyclopentyltrichlorosilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 14807-96-6 | talc | YES | Carc |  |  |  |  |  |  |  | 0.8 |
| 14808-60-7 | Crystalline silica | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 14960-06-6 | sodium lauriminodipropionate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 15096-52-3 | sodium aluminum fluoride | YES | 1st ITSL |  | 270 | 8 hr | YES |  |  |  |  |
| 15245-12-2 | nitric acid, ammonium calcium salt | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 15321-61-6 | iron oxalate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 15821-83-7 | propylene glycol n-butyl ether (beta isomer) | YES | 1st ITSL |  | 77 | annual | YES |  |  |  |  |
| 15956-58-8 | manganese 2-ethylhexanoate | YES | 1st ITSL |  | 0.3 | annual* | YES |  |  |  |  |
| 16065-83-1 | Chromium, trivalent | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 16079-88-2 | 1-bromo-3-chloro-5,5dimethylhydantoin | YES | 1st ITSL |  | 2 | 8 hr | YES |  |  |  |  |
| 16369-21-4 | n-propylethanolamine | YES | 1st ITSL |  | 28 | annual | YES |  |  |  |  |
| 16691-43-3 | 3-amino-5-mercapto-1,2,4-triazole | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 16753-62-1 | methylvinyldimethoxysilane | YES | 1st ITSL |  | 100 | annual | YES |  |  |  |  |
| 16881-77-9 | methyldimethoxysilane | YES | 1st ITSL |  | 92 | annual | YES |  |  |  |  |
| 16883-83-3 | 1,3-pentanediol-2,2,4-trimethyl-3(benzyl phthalate)-isobutyrate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 16893-85-9 | sodium silicofluoride | YES | 1st ITSL |  | 250 | 8 hr | YES |  |  |  |  |
| 16919-31-6 | ammonium hexafluorozirconate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 17557-23-2 | neopentyl glycol diglycidyl ether | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 17639-93-9 | methyl chloroproprionate | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 18063-03-1 | 2,6-difluorobenzamide | YES | 1st ITSL |  | 11 | annual | YES |  |  |  |  |
| 18300-89-5 | cinnamate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 18395-30-7 | isobutyltrimethoxysilane | NO |  | >75th\% | 200 | annual |  |  |  |  |  |
| 18540-29-9 | chromium, hexavalent - mist | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 0.008 | annual* | YES |  |  |  | 8E-05 |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 18540-29-9 | chromium, hexavalent - particulate | YES | $\begin{gathered} \text { 1st ITSL, } \\ \text { Carc } \end{gathered}$ |  | 0.1 | annual* | YES |  |  |  | 8E-05 |
| 18868-43-4 | molybdenum dioxide | YES | 1st ITSL |  | 30 | 8 hr | YES |  |  |  |  |
| 19089-47-5 | propylene glycol monoethyl ether (alpha) | YES | 1st ITSL |  | 23 | annual | YES |  |  |  |  |
| 19430-93-4 | perfluorobutylethylene | NO |  | >75th\% | 340 | annual |  |  |  |  |  |
| 19549-80-5 | 4,6-dimethyl-2-heptanone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 19666-30-9 | oxadiazon | YES | Carc |  |  |  |  |  |  |  | 0.05 |
| 20324-33-8 | tripropylene glycol methyl ether, dowanol 62b | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 20536-16-7 | tetrachlorodisilane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 21324-40-3 | lithium hexafluorophosphate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 21348-59-4 | niobium oxalate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 22407-51-8 | tetramethylchlorovinyldisiloxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 22431-89-6 | 3,3,6,6-tetramethyl-1,2-dioxane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 23410-40-4 | 1,2-ethanediamine, $n$-(3-(dimethoxymethylsilyl)-2methylpropyl) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 24304-00-5 | aluminum nitride | YES | 1st ITSL |  | 0.03 | annual | YES |  |  |  |  |
| 24510-87-0 | flumethasone 5 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 24729-96-2 | clindamycin phosphate | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 24937-79-9 | polyvinylidene fluoride | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 24938-91-8 | polyglycol 59-13 | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 25013-15-4 | vinyl toluene | YES | 1st ITSL |  | 6 | annual* | YES |  |  |  |  |
| 25036-25-3 | diglycidyl ether of bisphenol a | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 25068-38-6 | bisphenol a/epichlorohydrin resin | NO |  | >75th\% | 160 | annual |  |  |  |  |  |
| 25085-99-8 | bisphenol epoxy resin | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 25154-52-3 | nonyl phenol (mixed isomers) | YES | 1st ITSL |  | 30 | annual* | YES |  |  |  |  |
| 25168-26-7 | 2,4-D, isooctyl ester | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future <br> TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 25265-71-8 | dipropylene glycol | NO |  | >75th\% | 242 | annual |  |  |  |  |  |
| 25265-77-4 | texanol | YES | 1st ITSL |  | 55 | annual | YES |  |  |  |  |
| 25322-68-3 | polyethylene glycol | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 25322-69-4 | polypropylene glycol | YES | 1st ITSL |  | 49 | annual | YES |  |  |  |  |
| 25340-17-4 | diethylbenzene mixture | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 25498-49-1 | tripropylene glycol methyl ether | YES | 1st ITSL |  | 11 | annual | YES |  |  |  |  |
| 25550-14-5 | ethyl toluene -mixture | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 25551-13-7 | trimethylbenzenes (mixed isomers) | YES | 1st ITSL, 2nd ITSL |  | 50 | annual | YES | 1200 | 8 hr | YES |  |
| 25973-55-1 | benzotriazol dimethylpropyl phenol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 25988-97-0 | dimethylamine-epichlorohydrin polymer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 26062-79-3 | polydimethyl diallyl ammonium chloride | NO |  | >75th\% | 1000 | annual* |  |  |  |  |  |
| 26142-30-3 | diglycidyl ether of polyglycol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 26172-55-4 | 5-chloro-2-methyl-4-isothiazolin-3-one | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 26447-40-5 | 1,1'-methylene bisisocyanatobenzene | YES | 1st ITSL |  | 0.6 | annual* | YES |  |  |  |  |
| 26471-62-5 | toluene diisocyanate | YES | 1st ITSL, Carc |  | 0.07 | annual* | YES |  |  |  | 0.03 |
| 26530-20-1 | octylisothiazolone | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 26544-20-7 | mcpa 2-ehe (2-methyl-4chlorophenoxyacetic acid 2-ethylhexyl ester) | YES | 1st ITSL |  | 90 | annual | YES |  |  |  |  |
| 26761-40-0 | diisodecyl ester phthalate | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 26780-96-1 | poly(1,2-dihydro-2,2,4trimethylquinoline) | YES | 1st ITSL |  | 35 | annual* | YES |  |  |  |  |
| 26952-20-5 | picloram, isooctyl ester | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 26952-21-6 | isooctanol | NO |  | >75th\% | 2700 | 8 hr |  |  |  |  |  |
| 27078-75-7 | 4,6-difluoro-2-methoxypyrimidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 27208-37-3 | Cyclopenta[c,d]pyrene | YES | EPA Carc |  |  |  |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants
*NEW AvgT "annual*" Some ITSLs with 24-hr avg. time were converted to annual avg. because ITSL was derived to protect for chronic effects.

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th } \% \\ & \hline \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd <br> ITSL <br> AvgT | $\begin{array}{c\|} \hline \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{array}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 27253-31-2 | cobalt neodecanoate | YES | 1st ITSL |  | 1.4 | 8 hr | YES |  |  |  |  |
| 27253-32-3 | manganese neodecanoate | YES | 1st ITSL |  | 0.3 | annual* | YES |  |  |  |  |
| 27274-31-3 | polyethylene glycol monoallyl ether | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 27646-80-6 | 2-methylamino-2-methyl-1-propanol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 27668-52-6 | octadecyldimethyl (3- <br> (trimethoxysilyl)propyl) ammonium chloride | NO |  | >75th\% | 170 | annual |  |  |  |  |  |
| 28300-74-5 | antimony potassium tartrate | YES | 1st ITSL |  | 5 | 8 hr | YES |  |  |  |  |
| 28476-83-7 | 2-butenedioic acid (z)-dibutyl ester, polymer with chloroethene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 28553-12-0 | diisononyl phthalate | YES | 1st ITSL |  | 75 | annual | YES |  |  |  |  |
| 28729-52-4 | dimethylcyclopentane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 28729-54-6 | m-propyl toluene | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 28961-43-5 | triacrylate ester | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 28984-69-2 | 4,4-(5h)-oxazoledimethanol, 2(hepadecanyl) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 29387-86-8 | propylene glycol, n-butyl ether (mixed isomers) | YES | 1st ITSL |  | 77 | annual | YES |  |  |  |  |
| 29733-18-4 | diisodecyl glutarate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 29911-27-1 | dipropylene glycol monopropyl ether | YES | 1st ITSL |  | 5 | annual | YES |  |  |  |  |
| 29911-28-2 | dipropylene glycol monobutyl ether | YES | 1st ITSL |  | 11 | annual | YES |  |  |  |  |
| 30030-25-2 | vinylbenzylchloride | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 30705-14-7 | SR 1153 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 31138-65-5 | sodium glucoheptonate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 31726-34-8 | poly(oxy-1,2-ethanediyl),alpha-hexyl-omega-hydroxy | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 34375-28-5 | hydroxymethylamino ethanol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 34590-94-8 | dipropylene glycol methyl ether | NO |  | >75th\% | 720 | annual* |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd <br> ITSL <br> AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 35176-78-4 | polyethylene terephthalate (uncoated) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 35794-11-7 | 3,5-dimethylpiperidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 35884-42-5 | dowanol dpnb | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 37251-67-5 | polyethylene polypropylene glycol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 38436-16-7 | perfluorobutylethylmethyldichlorosilan e | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 39464-66-9 | lauryl alcohol, phosphated | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 40758-65-4 | 4,6-dichloro-2-ethoxypyrimidine | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 41556-26-7 | bis(pentamethylpiperdinyl)sebacate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 41593-38-8 | propylene glycol monophenyl ether | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 42978-66-5 | tripropylene glycol diacrylate | YES | 1st ITSL |  | 22 | annual | YES |  |  |  |  |
| 44992-01-0 | acryloyloxyethyltrimethyl ammonium chloride | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 46438-39-5 | monobutyl monophenyl phosphoric acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 50791-87-2 | methylvinylbis(N-methylace | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 51200-87-4 | dimethyloxazolidine | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 51730-94-0 | dipropylene glycol phenyl ether | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 51811-38-2 | tryfac 5556 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 52125-53-8 | propylene glycol monoethyl ether (mixture) | YES | 1st ITSL |  | 23 | annual | YES |  |  |  |  |
| 53880-05-0 | isophorone diisocyanate polymer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 55818-57-0 | phenol, 4,4-(1-methylethylidene)bis, polymer with (chloromethyl)oxiran | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 55934-93-5 | tripropylene glycol n-butyl ether | NO |  | >75th\% | 116 | annual* |  |  |  |  |  |
| 56539-66-3 | 3-methoxy-3methyl-1butanol | YES | 1st ITSL |  | 13 | annual | YES |  |  |  |  |
| 56741-95-8 | bropirimine | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 56780-58-6 | 2-hydroxy-3-trimethylammoniopropyl ether starch | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants

| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 57018-52-7 | propylene glycol tert-butyl ether | NO |  | >75th\% | 329 | annual |  |  |  |  |  |
| 60304-36-1 | aluminum potassium fluoride | YES | 1st ITSL |  | 0.2 | annual | YES |  |  |  |  |
| 60676-86-0 | amorphous silica - fused silica | YES | 1st ITSL |  | 60 | 8 hr | YES |  |  |  |  |
| 60966-36-1 | bisnoralcohol | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 61477-94-9 | pirmenol hydrochloride | YES | 1st ITSL |  | 3 | annual | YES |  |  |  |  |
| 61788-93-0 | coco alkyldimethyl amines | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 61790-33-8 | tallow alkylamines | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 61790-53-2 | amorphous silica - diatomaceous earth | YES | 1st ITSL |  | 60 | 8 hr | YES |  |  |  |  |
| 61791-28-4 | ethoxy, tallow alcohol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 63148-57-2 | Dow Corning Fluid 1107 | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |
| 63148-62-9 | high molecular wt. silicon | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 63148-65-2 | polyvinyl butyral | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 63449-39-8 | chlorinated paraffins | YES | Carc |  |  |  |  |  |  |  | 0.03 |
| 63716-40-5 | n-butoxy propanol (mixed isomers) | YES | 1st ITSL |  | 77 | annual | YES |  |  |  |  |
| 63937-30-4 | anhydro-dimethylamino hexose reductone | YES | 1st ITSL |  | 0.6 | annual | YES |  |  |  |  |
| 64248-62-0 | 3,4-difluorobenzonitrile | YES | 1st ITSL |  | 0.6 | annual | YES |  |  |  |  |
| 64265-57-2 | crosslinker cx100 | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 64475-85-0 | mineral spirits | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64485-82-1 | thiazole ester | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 64741-41-9 | naphtha heavy straight run | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64741-42-0 | naphtha, full range straight run | YES | 1st ITSL |  | 18 | annual | YES |  |  |  |  |
| 64741-44-2 | straight run middle distillate | YES | 1st ITSL |  | 36 | annual | YES |  |  |  |  |
| 64741-54-4 | naphtha, heavy catalytic cracked | NO |  | >75th\% | 115 | annual |  |  |  |  |  |
| 64741-55-5 | naphtha (petroleum), light catalytic cracked | NO |  | >75th\% | 5600 | annual* |  |  |  |  |  |
| 64741-56-6 | residues, (petroleum), vacuum | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |

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| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd <br> ITSL <br> AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 64741-59-9 | distillates, (petroleum), light catalytic cracked | YES | 1st ITSL |  | 93 | annual | YES |  |  |  |  |
| 64741-62-4 | clarified oils (petroleum), catalytic cracked | YES | 1st ITSL |  | 12 | annual | YES |  |  |  |  |
| 64741-63-5 | naphtha, light catalytic reformed | YES | 1st ITSL |  | 100 | annual | YES |  |  |  |  |
| 64741-64-6 | naphtha, full range alkylate | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64741-65-7 | heavy alkylate naphtha | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64741-66-8 | light alkylate naphtha | NO |  | >75th\% | 138 | annual |  |  |  |  |  |
| 64741-68-0 | heavy catalytic reformed naphtha | YES | 1st ITSL |  | 70 | annual | YES |  |  |  |  |
| 64741-81-7 | distillates (petroleum), heavy thermal cracked | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |
| 64741-82-8 | distillates (petroleum), light thermal cracked | YES | 1st ITSL |  | 93 | annual | YES |  |  |  |  |
| 64741-83-9 | naphtha, heavy thermal cracked | NO |  | >75th\% | 5600 | annual* |  |  |  |  |  |
| 64741-86-2 | sweetened middle distillate | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 64741-88-4 | solvent refined heavy paraffnic distillate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64741-89-5 | distillates (petroleum) solvent-refined light paraffinic | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-06-9 | extracts (petroleum), middle distillate solvent | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 64742-14-9 | petroleum distillates, acid treated | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 64742-30-9 | distillates (petroleum), chemically neutralized middle | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 64742-46-7 | hydrotreated middle distillate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-47-8 | hydrotreated light distillate | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 64742-48-9 | hydrotreated heavy napht | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64742-49-0 | hydrotreated light naphtha | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64742-52-5 | hydrotreated heavy naphthenic | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | distillate |  |  |  |  |  |  |  |  |  |  |
| 64742-53-6 | hydrotreated light naphthenic distillate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-54-7 | hydrotreated heavy paraffinic mineral oil | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-55-8 | hydrotreated light paraffinic distillate | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-62-7 | residual oils (petroleum) solventdewaxed | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-65-0 | dewaxed heavy paraffinic mineral oil | YES | 1st ITSL |  | 50 | 8 hr | YES |  |  |  |  |
| 64742-80-9 | hydrodesulfurized middle distillate | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 64742-81-0 | hydrodesulfurized kerosene | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 64742-82-1 | naphtha (petroleum) hydrodesulfurized heavy | YES | 1st ITSL |  | 14 | annual | YES |  |  |  |  |
| 64742-88-7 | solvent naphtha medium aliphatic | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64742-89-8 | solvent naphtha light aliphatic | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 64742-94-5 | heavy aromatic solvent naphtha | YES | 1st ITSL |  | 70 | annual* | YES |  |  |  |  |
| 64742-95-6 | light aromatic solvent naphtha (petroleum) | YES | 1st ITSL |  | 61 | annual | YES |  |  |  |  |
| 64742-96-7 | solvent naphtha (petroleum) heavy aliphatic | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 64771-72-8 | norpar 12 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 65402-65-5 | 4-hydroxytetramethyl piperadine free radical (4-oh-tempo) | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 66071-86-1 | LV 837/821 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 67701-10-4 | sodium soap 903923 | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 67701-11-5 | sodium soap 900602 | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 67762-41-8 | linear primary alcohol | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 67762-90-7 | siloxanes and silicones(silica filled polydimethylsiloxane) | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 67784-80-9 | soybean oil, methyl esters | YES | 1st ITSL |  | 15 | annual | YES |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{gathered} \text { 1st ITSL } \\ <=75 \text { th\% } \end{gathered}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 67812-17-3 | 3-trimethoxysilyl propylmethyl methylphosphonate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68002-20-0 | 1,2,3-triazine-2,4,6-triamine polymer with methylated formaldehyde | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68003-28-1 | polyamide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68037-58-1 | high molecular wt. silicon | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68037-76-3 | alphamethylstyrene(dodecyl)polysilox ane | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68037-77-4 | ethylmethylsiloxane, 2phenylpropylmethylsiloxane copolymer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68037-88-7 | high molecular weight sili | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68071-85-2 | Spenkel F34 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68083-19-2 | high molecular wt. silicon | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68083-20-5 | linear methylvinylsiloxane ppolymer hydroxl endblock | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68083-40-9 | 2-hydroxy-4(2'-hydroxy- <br> 3'octoxypropoxy)-benzophenone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68092-49-9 | 2-hydroxy-4(2'-hydroxy- <br> 3'dacyloxypropoxy)-benzophenone | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68131-40-8 | tergitol 15-s-3 | NO |  | >75th\% | 290 | annual* |  |  |  |  |  |
| 68132-02-5 | coumarone indene resin | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68309-52-4 | Nylen 5 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68334-30-5 | diesel fuel | YES | 1st ITSL |  | 70 | annual | YES |  |  |  |  |
| 68390-56-7 | diketene hydrogenated fatty acids | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68410-00-4 | distillates (petroleum), crude oil | YES | 1st ITSL |  | 19 | annual | YES |  |  |  |  |
| 68410-23-1 | polyethylenepolyamine reaction products with c18-unsat. fatty acids | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68439-49-6 | ethoxylated c16-18 alcohols | YES | 1st ITSL |  | 4 | annual | YES |  |  |  |  |
| 68458-91-3 | Solvar \& LV 820 | NO |  | default | 0.1 | annual |  |  |  |  |  |

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| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 68459-31-4 | fatty acids c9-11 branched glycidyl esters polymer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68476-86-8 | petroleum gases, liquefied, sweetened | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68477-31-6 | aromatic petroleum derivative solvent | YES | 1st ITSL |  | 13 | annual | YES |  |  |  |  |
| 68515-40-2 | alkyl benzyl phthalate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68515-44-6 | branched and linear diheptyl phthalate ester | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68516-16-5 | sulfuric acid c6-10 alkyl esters | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68526-86-3 | tridecanol | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 68551-17-7 | heavy naphtha | NO |  | >75th\% | 3500 | 8 hr |  |  |  |  |  |
| 68575-36-0 | 3,5-dichloro-a-methyl st | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 68608-26-4 | sodium petroleum sulfonate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68610-11-7 | diethylenetriamine reaction product with bisphenol a | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68783-24-4 | di-tallow alkylamines | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68918-22-9 | high molecular wt. silicon | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68955-35-1 | naphtha, catalytic reformed | NO |  | >75th\% | 350 | annual* |  |  |  |  |  |
| 68956-56-9 | hydrocarbons, terpene processing byproducts | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 68987-42-8 | ethylenated benzene residues | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 68990-79-4 | oils, vegetable, mixed with animal oil methylesters, polymerized, oxidixed | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 69012-64-2 | amorphous silica - silica fume | YES | 1st ITSL |  | 60 | 8 hr | YES |  |  |  |  |
| 69013-18-9 | alcohols c8-18 ethoxylated propoxylated | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 69029-39-6 | polyglycol 26-2 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 69102-90-5 | butadiene homopolymer | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 69430-24-6 | high molecular wt. silicon | YES | 1st ITSL |  | 30 | annual | YES |  |  |  |  |

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| CAS Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 69696-98-6 | hexane 1,6-bis(tributyl ammonium bromi | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 69991-67-9 | fomblin perfluorpolyether | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 70131-67-8 | high molecular wt. silicon | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 70657-70-4 | 2-methoxy-1-propanol acetate | NO |  | >75th\% | 500 | annual* |  |  |  |  |  |
| 70914-20-4 | c6-8 branched alcohols | YES | 1st ITSL |  | 13 | annual | YES |  |  |  |  |
| 71888-89-6 | diisoheptyl phthalate | YES | 1st ITSL |  | 100 | annual* | YES |  |  |  |  |
| 71945-54-5 | 3-(1,1-dimethylethoxy)-heptane | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 75782-86-4 | alcohols c12-13 | YES | 1st ITSL |  | 31 | annual | YES |  |  |  |  |
| 77820-58-7 | 2-amino-3-chlorobenzoic acid methyl ester | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 78330-21-9 | c11-c14 isoalcohols, c14 rich, ethoxylated alcohol | YES | 1st ITSL |  | 8 | annual | YES |  |  |  |  |
| 82586-54-7 | quinapril step 8 | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 82919-37-7 | methyl pentamethyl-4-piperidinyl ester of decanedioic acid | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 84632-65-5 | pyrrolo[3,4-c]pyrrole-1,4-dione,3,6-bis(4-chlorophenyl)-2,5-dihydro | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 86753-78-8 | Solsperse 5000 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 88230-35-7 | oxo-hexyl acetate | YES | 1st ITSL |  | 81 | annual | YES |  |  |  |  |
| 88851-61-0 | trospectomycin sulfate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 88917-22-0 | dipropylene glycol methyl ether acetate | NO |  | >75th\% | 930 | annual* |  |  |  |  |  |
| 90438-79-2 | oxo-heptyl acetate | YES | 1st ITSL |  | 41 | annual | YES |  |  |  |  |
| 90622-57-4 | isopar h | NO |  | >75th\% | 128 | annual |  |  |  |  |  |
| 95481-62-2 | dibasic ester | YES | 1st ITSL |  | 1 | annual | YES |  |  |  |  |
| 97658-80-5 | 5-bp-bisenamine | YES | 1st ITSL |  | 10 | annual | YES |  |  |  |  |
| 98516-30-4 | propanol, 1(or 2) ethoxy, acetate isoparaffinic petroleum hydrocarbon | NO |  | default | 0.1 | annual |  |  |  |  |  |

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| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{gathered} \text { 1st ITSL } \\ <=75 \text { th\% } \end{gathered}$ |  | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th } \% \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 98967-40-9 | flumetsulam | YES | 1st ITSL |  | 26 | annual | YES |  |  |  |  |
| 98967-55-6 | n-(2,6-difluorophenyl)-7-methyl-1h-1,2,4-triazolo(1,5a)pyrimidine-2-su | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 102054-10-4 | bis(2-methoxy-1-methylethy | YES | 1st ITSL |  | 6 | annual | YES |  |  |  |  |
| 103335-54-2 | 4-aza acid | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 103429-90-9 | 3-methoxy-3methyl-1butyl acetate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 103980-44-5 | ceftiofur hydrochloride | NO |  | >75th\% | 166 | annual |  |  |  |  |  |
| 106917-31-1 | sanduvor 3068 liquid | YES | 1st ITSL |  | 52 | annual | YES |  |  |  |  |
| 108419-32-5 | exxate 800 -octyl acetate | NO |  | >75th\% | 110 | annual* |  |  |  |  |  |
| 108419-33-6 | exxate 900 | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 108419-34-7 | exxate 1000 | YES | 1st ITSL |  | 17 | annual | YES |  |  |  |  |
| 108419-35-8 | c11-14 branched alkyl acetates | NO |  | >75th\% | 300 | annual* |  |  |  |  |  |
| 109265-71-6 | Solsperse 12000 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 110839-13-9 | 1,3-benzenedimethanamine polymer with 2,2'-((1-methylethylidene) bis(4 | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 110888-15-8 | 4-chloro-3-fluorobenzonitrile | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 111109-77-4 | dipropylene glycol dimethyl ether | YES | 1st ITSL |  | 59 | annual* | YES |  |  |  |  |
| 111381-89-6 | branched and linear heptyl nonyl phthalate ester | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 112926-00-8 | amorphous silica - precipitated silica and silica gel | YES | 1st ITSL |  | 60 | 8 hr | YES |  |  |  |  |
| 112945-52-5 | amorphous silica - pyrogenic or fumed silica | YES | 1st ITSL |  | 60 | 8 hr | YES |  |  |  |  |
| 113171-12-3 | n-(2,6-difluorophenyl)-5-amino-1h-1,2,4-triazole-3-sulfonamide | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 117482-84-5 | 3-chloro-4-fluorobenzonitrile | YES | 1st ITSL |  | 2 | annual | YES |  |  |  |  |
| 123312-54-9 | distearyldimethylammonium bisulfate | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 123333-53-9 | 1-hydroxy benzotriazole | NO |  | default | 0.1 | annual |  |  |  |  |  |

Table 4. Current and Proposed MDEQ Air Quality Division Screening Level List of Toxic Air Contaminants
*NEW AvgT "annual*" Some ITSLs with 24-hr avg. time were converted to annual avg. because ITSL was derived to protect for chronic effects.

| CAS <br> Number | Chemical Name | Future TAC? | Basis for Yes | Reason for No, etc. | $\begin{gathered} \text { 1st } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | *NEW AvgT | $\begin{aligned} & \text { 1st ITSL } \\ & <=75 \text { th\% } \end{aligned}$ | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ \left(\mu \mathrm{g} / \mathrm{m}^{3}\right) \end{gathered}$ | 2nd ITSL AvgT | $\begin{gathered} \text { 2nd } \\ \text { ITSL } \\ <=75 \text { th\% } \end{gathered}$ | IRSL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 126803-73-4 | n-(2,6-dichloro-3-methylphenyl)-5,7dimethyoxy (1,2,4)triazo...[de-511] | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 129879-84-1 | 5-amino-1,2,4-triazole-3-sulfonyl chloride | YES | 1st ITSL |  | 7 | annual | YES |  |  |  |  |
| 130014-38-9 | trifluoropropylsilsesquioxane, dimethylhydrogensilyoxy-terminated | NO |  | default | 0.1 | annual |  |  |  |  |  |
| 136797-56-3 | FC-247 | YES | 1st ITSL |  | 24 | annual | YES |  |  |  |  |
| 136816-75-6 | atevirdane mesylate | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 144669-03-4 | hexenylsiloxane | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 144669-04-5 | hexenylsiloxanes | YES | 1st ITSL |  | 16 | annual | YES |  |  |  |  |
| 166524-65-8 | 2-ethoxy-4,6-difluoropyrimidine | YES | 1st ITSL |  | 20 | annual | YES |  |  |  |  |
| 166524-75-0 | 2,2'-dithiobis(5-ethoxy-7-fluoro[1,2,4]triazolo(1,5-c)pyrimidine | YES | 1st ITSL |  | 0.8 | annual | YES |  |  |  |  |
| 170557-43-4 | dowanol tmh-deg borate ester | YES | 1st ITSL |  | 32 | annual | YES |  |  |  |  |

## ApPENDIX K:

## COMPARISON OF HAP AND TAC Screening Level Lists

# Appendix K - Comparison of HAP and TAC Screening Level Lists 

MDNRE-AQD Toxic Air Contaminants List Compared to the EPA Hazardous Air
Pollutants List
February 16, 2010
Robert Sills, Toxics Unit Supervisor, MDNRE-Air Quality Division
Michigan's Air Pollution Control Rules (under Part 55 of NREPA) to regulate the emission of toxic air contaminants (TACs) have been in place since 1992. TACs are defined (Rule 336. I120(f)) as any air contaminants for which there are no national - ambient air quality standard and which are or may become harmful to the environment when present in the outdoor atmosphere in sufficient quantities and duration. The TAC definition lists 41 substances which are not TACs. This list includes the six pollutants that have national ambient air quality standards and 35 other substances.

The original air toxics rules (1992) included the current definition of TACs based on the Michigan Air Toxics Policy Committee (1989) recommendation that the AQD should address a large list of TACs plus any other substances which the AQD determines to be of concern at a specific site. The TAC definition was re-visited again in 1997 by the AQD Air Toxics Subcommittee. Based on the Subcommittee's discussion and recommendations, the AQD made revisions to the air toxics rules in 1998, retaining the open-ended TAC definition but providing greater flexibility in the rules and adopting a small quantity exemption.

The regulatory programs of the AQD and the EPA are intended to provide a level of protection against the potential risks of air toxics and therefore ensure the public that facility emissions are safe. However, the federal regulations for air toxics have significant limitations. These limitations include the specific air toxics that are regulated, types of facilities that are regulated, the quantity of emissions that are subject to regulation, and the risk assessment requirements.

EPA lists 187 substances as hazardous air pollutants (HAPs) that are subject to federal regulation. Major sources are any facility that emits 10 tons per year of any HAP or 25 tons per year of any combination of HAPs. EPA has made progress in developing pollution control technology requirements for categories of major sources. EPA is also required to assess the need for standards to protect public health and the environment. However, EPA has completed very few of these residual risk assessments on their listed HAPs. In those instances where a technology standard is established and a residual risk assessment has been completed, the source category is exempted from the AQD air toxics regulations so there is no regulatory redundancy.

Michigan's program is broader than the federal program to better ensure public health protection from air toxics emissions from proposed new or modified sources, while also including a number of exemptions for sources and air toxics emission levels which have been specifically determined to pose no unacceptable risks to the public health. Michigan's program is designed to supplement and complement (without redundancy) the federal air toxics regulations.

Although the EPA HAP list captures many substances recognized as high-concern air contaminants, there are many non-HAP air toxics which can potentially pose health risks to the public who are exposed to them. These substances include pharmaceuticals, pesticides,
metals, inorganic compounds, and organic compounds. In Michigan, approximately 1200 TACs listed in the attached table were identified in Permit to Install applications for proposed facilities. As part of the permit application review, the TACs were evaluated by AQD toxicologist staff and health-based screening levels were developed, which provide a level of protection from adverse health effects. The AQD frequently provides assurances to the concerned public about the safety of existing or proposed facility air emissions, and is able to do so because of the healthbased screening levels and the open-ended TAG definition.

As indicated in the 3rd and 4th columns of Table 2, there are many TACs which are not HAPs but which are carcinogenic. TACs also pose concerns for potential acute toxicity, developmental effects, sensitization, respiratory effects such as asthma, liver or kidney effects, neurological effects, etc. Table 1 shows some specific examples of non-HAP TACs and their primary public health concerns.

Table 1. Example non-HAP TACs and their primary public health concerns.

| Toxic Air Contaminants (TACs) | Primary Public Health Concern |
| :--- | :--- |
| Aldrin, benzaldehyde, bromodichloromethane, <br> dimethylvinyl chloride, hydrazine sulfate, <br> molybdenum trioxide, nitromethane, <br> tetrahydrofuran, etc. | Carcinogens |
| Ammonia, Glutaraldehyde, Hydrogen sulfide | Irritation of the eyes and respiratory tract |
| $2,4,6$-trinitrotoluene | Liver toxicity, anemia |
| Barium | Muscle toxicity; environmental persistence |
| Bromine | Respiratory irritation, headache |
| Chlorine dioxide | Lung toxicity |
| Chlormadinone acetate | Reproductive effects |
| Chlorpyrifos | Nervous system toxicity |
| Colophony, Isophorone diisocyanate | Asthma exacerbation, sensitizer |
| Dibutyltin oxide | Immune function and central nervous system <br> toxicity |
| Methylene diphenyl isocyanate | Respiratory tract toxicity |
| Melengesterol acetate | Reproductive toxicity; menstruation blockage <br> Osmium tetroxideIrritant to the eyes, nose and throat; pulmonary <br> edema and bronchitis |
| Sulfuric acid | Eye and respiratory irritancy and <br> corrosiveness, shortness of breath |
| Tetrachlorobenzene | Liver and kidney toxicity; environmental <br> persistence |
| Thallium | Developmental, respiratory, and <br> gastrointestinal effects; environmental <br> persistence |
| Vanadium pentaoxide | Bronchitis, emphysema, respiratory tract <br> irritation |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 3052-70-8 | (1-methylethylidene)bis(1,1-dimethylpropyl)peroxide | no | no |
| 6713-03-7 | 1-(2-hydroxyethylthio)propane | no | no |
| 630-20-6 | 1,1,1,2-tetrachloroethane | no | yes |
| 811-97-2 | 1,1,1,2-tetrafluoroethane | no | no |
| 460-73-1 | 1,1,1,3,3-pentafluoropropane | no | no |
| 79-34-5 | 1,1,2,2-tetrachloroethane | yes | yes |
| --0 | 1,1,2,4-tetramethyl-1-1-1-sila-2-aza-cyclopentane | no | no |
| 76-13-1 | 1,1,2-trichloro-1,2,2-trifluoroethane | no | no |
| 79-00-5 | 1,1,2-trichloroethane | yes | yes |
| 3006-86-8 | 1,1-di-(tert-buytlperoxy)cyclohexane | no | no |
| 1717-00-6 | 1,1-dichloro-1-fluoroethane | no | no |
| 75-34-3 | 1,1-dichloroethane | yes | no |
| 75-37-6 | 1,1-difluoroethane | no | no |
| 612-00-0 | 1,1-diphenylethane | no | no |
| 26447-40-5 | 1,1'-methylene bisisocyanatobenzene | no | no |
| 2403-89-6 | 1,2,2,6,6-pentamethyl-4-piperidinol | no | no |
| 634-66-2 | 1,2,3,4-tetrachlorobenzene | no | no |
| 634-90-2 | 1,2,3,5-tetrachlorobenzene | no | no |
| 68002-20-0 | 1,2,3-triazine-2,4,6-triamine polymer with methylated formaldehyde | no | no |
| 87-61-6 | 1,2,3-trichlorobenzene | no | no |
| 96-18-4 | 1,2,3-trichloropropane | no | no |
| 526-73-8 | 1,2,3-trimethylbenzene | no | no |
| 95-94-3 | 1,2,4,5-tetrachlorobenzene | no | no |
| 95-93-2 | 1,2,4,5-tetramethylbenzene | no | no |
| 120-82-1 | 1,2,4-trichlorobenzene | yes | no |
| 95-63-6 | 1,2,4-trimethylbenzene | no | no |
| 106-88-7 | 1,2-butylene oxide | yes | yes |
| 95-50-1 | 1,2-dichlorobenzene | no | no |
| 107-06-2 | 1,2-dichloroethane | yes | yes |
| 540-59-0 | 1,2-dichloroethylene | no | no |
| 23410-40-4 | 1,2-ethanediamine, n-(3-(dimethoxymethylsilyl)-2-methylpropyl) | no | no |
| 87-90-1 | 1,3,5-trichloroisocyanuric acid | no | no |
| 108-67-8 | 1,3,5-trimethylbenzene | no | no |
| 110839-13-9 | 1,3-benzenedimethanamine polymer with 2,2'-((1-methylethylidene) bis(4 | no | no |
| 1477-55-0 | 1,3-bis(aminomethyl)benzenen | no | no |
| 106-99-0 | 1,3-butadiene | yes | yes |
| 77-48-5 | 1,3-dibromo-5,5-dimethylhydantoin | no | no |
| 118-52-5 | 1,3-dichloro-5,5-dimethylhydantoin | no | no |
| 541-73-1 | 1,3-dichlorobenzene | no | no |
| 542-75-6 | 1,3-dichloropropene | yes | yes |
| 646-06-0 | 1,3-dioxolane | no | no |
| 16883-83-3 | 1,3-pentanediol-2,2,4-trimethyl-3-(benzyl phthalate)-isobutyrate | no | no |
| 110-63-4 | 1,4 butanediol | no | no |
| 106-46-7 | 1,4-dichlorobenzene | yes | yes |
| 123-91-1 | 1,4-dioxane | yes | yes |
| 592-42-7 | 1,5-hexanediene | no | no |
| 629-11-8 | 1,6-hexanediol | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 6674-22-2 | 1,8-diazabicyclo[5.4.0]undec-7-ene | no | no |
| 58-36-6 | 10,10'-oxybisphenoxarsine oxide | no | no |
| --0 | 100 sxl | unknown | no |
| 13209-41-1 | 17,21-dihydroxy-16 alpha-methylpregna-1,4,9(11)-triene-3,20-dione | no | no |
| 16079-88-2 | 1-bromo-3-chloro-5,5-dimethylhydantoin | no | no |
| 109-70-6 | 1-bromo-3-chloropropane | no | no |
| 109-65-9 | 1-bromobutane | no | no |
| 75-68-3 | 1-chloro-1,1-difluoroethane | no | no |
| 88-73-3 | 1-chloro-2-nitrobenzene | no | yes |
| 611-14-3 | 1-ethyl-2-methylbenzene | no | no |
| 592-76-7 | 1-heptene | no | no |
| 629-73-2 | 1-hexadecene | no | no |
| 947-19-3 | 1-hydroxcyclohexyl phenyl ketone | no | no |
| 123333-53-9 | 1-hydroxy benzotriazole | no | no |
| 90-12-0 | 1-methyl naphthalene | yes (POM) | yes |
| 108-03-2 | 1-nitropropane | no | no |
| 1569-01-3 | 1-propoxy-2-propanol | no | no |
| 1072-63-5 | 1-vinylimidazol | no | no |
| --0 | 2-(1-ethoxyethoxy)-6-(trifluroromethyl)-benzenethiol | no | no |
| 3731-51-9 | 2-(aminomethyl)pyridine | no | no |
| 540-84-1 | 2,2,4-trimethyl pentane | yes | no |
| 6846-50-0 | 2,2,4-trimethylpentanediol-1,3-diisobutyrate | no | no |
| 366-18-7 | 2,2'-bipyridyl | no | no |
| 77-76-9 | 2,2-dimethoxypropane | no | no |
| 166524-75-0 | 2,2'-dithiobis(5-ethoxy-7-fluoro[1,2,4]triazolo(1,5-c)pyrimidine | no | no |
| 1746-01-6 | 2,3,7,8-tetrachlorodibenzo(p)dioxin | yes | yes |
| 513-85-9 | 2,3-butanediol | no | no |
| 526-75-0 | 2,3-dimethyl phenol | no | no |
| 79-29-8 | 2,3-dimethylbutane | no | no |
| 565-59-3 | 2,3-dimethylpentane | no | no |
| 95-95-4 | 2,4,5-trichlorophenol | yes | no |
| 90-72-2 | 2,4,6-tri(dimethylaminomethyl)phenol | no | no |
| 88-06-2 | 2,4,6-trichlorophenol | yes | yes |
| 3764-01-1 | 2,4,6-trichloropyrimidine | no | no |
| 696-82-2 | 2,4,6-trifluoropyrimidine | no | no |
| 118-96-7 | 2,4,6-trinitrotoluene | no | no |
| 94-75-7 | 2,4-Dichlorophenoxyacetic Acid (2,4-d) | yes | no |
| 25168-26-7 | 2,4- Dichlorophenoxyacetic Acid (2,4-d) isooctyl ester | yes | no |
| 120-83-2 | 2,4-dichlorophenol | no | no |
| 108-08-7 | 2,4-dimethylpentane | no | no |
| 105-67-9 | 2,4-dimethylphenol | no | no |
| 51-28-5 | 2,4-dinitrophenol | yes | no |
| 121-14-2 | 2,4-dinitrotoluene | yes | yes |
| 123-54-6 | 2,4-pentanedione | no | no |
| 548-84-9 | 2,4-toluene diisocyanate | yes | yes |
| 5779-94-2 | 2,5-dimethylbenzaldehyde | no | no |
| 95-87-4 | 2,5-dimethylphenol | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 608-31-1 | 2,6-dichlorobenzenamine | no | no |
| 5509-65-9 | 2,6-difluoroaniline | no | no |
| 18063-03-1 | 2,6-difluorobenzamide | no | no |
| 385-00-2 | 2,6-difluorobenzoic acid | no | no |
| 1897-52-5 | 2,6-difluorobenzonitrile | no | no |
| 141-91-3 | 2,6-dimethyl morpholine | no | no |
| 576-26-1 | 2,6-dimethyl phenol | no | no |
| 606-20-2 | 2,6-dinitrotoluene | no | yes |
| 128-37-0 | 2,6-di-tert-butyl-p-cresol | no | yes |
| 87-62-7 | 2,6-xylidine | no | yes |
| 124-68-5 | 2-amino-2-methyl-1-propanol | no | no |
| 77820-58-7 | 2-amino-3-chlorobenzoic acid methyl ester | no | no |
| 359-07-9 | 2-bromo-1,1-difluoroethane | no | no |
| 103-63-9 | 2-bromoethylbenzene | no | no |
| 28476-83-7 | 2-butenedioic acid (z)-dibutyl ester, polymer with chloroethene | no | no |
| 111-76-2 | 2-butoxyethanol | yes | no |
| 111-75-1 | 2-butylaminoethanol | no | no |
| 2837-89-0 | 2-chloro-1,1,1,2-tetrafluoroethane | no | no |
| 363-51-9 | 2-chloro-6-fluorobenzenamine | no | no |
| 95-51-2 | 2-chloroaniline | no | no |
| 95-57-8 | 2-chlorophenol | no | no |
| 75-29-6 | 2-chloropropane | no | no |
| 100-37-8 | 2-diethylaminoethanol | no | no |
| 100-36-7 | 2-diethylaminoethylamine | no | no |
| 166524-65-8 | 2-ethoxy-4,6-difluoropyrimidine | no | no |
| 110-80-5 | 2-ethoxyethanol | yes | no |
| 97-95-0 | 2-ethyl butanol | no | no |
| 94-96-2 | 2-ethyl-1,3-hexanediol | no | no |
| 1758-88-9 | 2-ethyl-1,4-dimethyl benzene | no | no |
| 10431-98-8 | 2-ethyl-2-oxazoline | no | no |
| 110-73-6 | 2-ethylaminoethanol | no | no |
| 3814-34-4 | 2-ethylbutyl bromide | no | no |
| 123-05-7 | 2-ethylhexanal | no | no |
| 149-57-5 | 2-ethylhexanoic acid | no | no |
| 104-76-7 | 2-ethylhexanol | no | no |
| 103-09-3 | 2-ethylhexyl acetate | no | no |
| 103-11-7 | 2-ethylhexyl acrylate | no | no |
| 104-75-6 | 2-ethylhexylamine | no | no |
| 1070-10-6 | 2-ethylhexyltitanate | no | no |
| 7473-98-5 | 2-hydroxy-2-methyl-1-phenyl-1-propanone | no | no |
| 56780-58-6 | 2-hydroxy-3-trimethylammoniopropyl ether starch | no | no |
| 68092-49-9 | 2-hydroxy-4(2'-hydroxy-3'dacyloxypropoxy)-benzophenone | unknown | no |
| 68083-40-9 | 2-hydroxy-4(2'-hydroxy-3'octoxypropoxy)-benzophenone | unknown | no |
| 868-77-9 | 2-hydroxyethyl methacrylate | no | no |
| --0 | 2-mercapto-3-(trifluoromethyl)-phenol | no | no |
| 1589-47-5 | 2-methoxy-1-propanol | no | no |
| 70657-70-4 | 2-methoxy-1-propanol acetate | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 116-11-0 | 2-methoxy-1-propene | no | no |
| 109-86-4 | 2-methoxyethanol | yes | no |
| 78-78-4 | 2-methyl butane | no | no |
| 624-41-9 | 2-methyl butyl acetate | no | no |
| 137-32-6 | 2-methyl-1-butanol | no | no |
| 2682-20-4 | 2-methyl-4-isothiazolin-3-one | no | no |
| 27646-80-6 | 2-methylamino-2-methyl-1-propanol | no | no |
| 109-83-1 | 2-methylaminoethanol | no | no |
| 591-76-4 | 2-methylhexane | no | no |
| 91-57-6 | 2-methylnaphthalene | yes (POM) | no |
| 107-83-5 | 2-methylpentane | no | no |
| 91-59-8 | 2-naphthylamine | yes (POM) | yes |
| 102-81-8 | 2-n-dibutylaminoethanol | no | no |
| 79-46-9 | 2-nitropropane | yes | yes |
| 111-13-7 | 2-octanone | no | no |
| 10215-30-2 | 2-propoxy-1-propanol | no | no |
| 7580-85-0 | 2-tert-butoxyethanol | yes | no |
| 71945-54-5 | 3-(1,1-dimethylethoxy)-heptane | no | no |
| 677-21-4 | 3,3,3-trifluoropropene | no | no |
| 22431-89-6 | 3,3,6,6-tetramethyl-1,2-dioxane | no | no |
| 6574-99-8 | 3,4-dichlorobenzonitrile | no | no |
| 64248-62-0 | 3,4-difluorobenzonitrile | no | no |
| 95-65-8 | 3,4-dimethyl phenol | no | no |
| 68575-36-0 | 3,5-dichloro-a-methyl st | no | no |
| 107-54-0 | 3,5-dimethyl-1-hexyn-3-ol | no | no |
| 108-68-9 | 3,5-dimethylphenol | no | no |
| 35794-11-7 | 3,5-dimethylpiperidine | no | no |
| 591-22-0 | 3,5-lutidine | no | no |
| 16691-43-3 | 3-amino-5-mercapto-1,2,4-triazole | no | no |
| 126-06-7 | 3-bromo-1-chloro-5,5-dimethylhydantoin | no | no |
| 627-30-5 | 3-chloro-1-propanol | no | no |
| 563-47-3 | 3-chloro-2-methylpropene | no | yes |
| 117482-84-5 | 3-chloro-4-fluorobenzonitrile | no | no |
| 95-74-9 | 3-chloro-p-toluidine | no | no |
| 4420-74-0 | 3-mercaptopropyltrimethoxysilane | no | no |
| 56539-66-3 | 3-methoxy-3methyl-1butanol | no | no |
| 103429-90-9 | 3-methoxy-3methyl-1butyl acetate | no | no |
| 589-34-4 | 3-methylhexane | no | no |
| 96-14-0 | 3-methylpentane | no | no |
| 108-99-6 | 3 -picoline | no | no |
| 67812-17-3 | 3-trimethoxysilyl propylmethyl methylphosphonate | no | no |
| 28984-69-2 | 4,4-(5h)-oxazoledimethanol, 2-(hepadecanyl) | no | no |
| 5436-21-5 | 4,4-dimethoxy-2-butanone | no | no |
| 101-14-4 | 4,4-methylenebis(2-chloroaniline) | yes | yes |
| 40758-65-4 | 4,6-dichloro-2-ethoxypyrimidine | no | no |
| 1074-40-4 | 4,6-dichloro-2-methoxypyrimidine | no | no |
| 27078-75-7 | 4,6-difluoro-2-methoxypyrimidine | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 19549-80-5 | 4,6-dimethyl-2-heptanone | no | no |
| 103335-54-2 | 4-aza acid | no | no |
| --0 | 4-chloro-2-ethoxy-6-fluoropyrimidine | no | no |
| 110888-15-8 | 4-chloro-3-fluorobenzonitrile | no | no |
| 65402-65-5 | 4-hydroxytetramethyl piperadine free radical(4-oh-tempo) | no | no |
| 4652-27-1 | 4-methoxy-3-buten-2-one | no | no |
| 100-06-1 | 4-methoxyacetophenone | no | no |
| 622-97-9 | 4-methylstyrene | no | no |
| 100-02-7 | 4-nitrophenol | yes | no |
| 4994-16-5 | 4-phenylcyclohexene | no | no |
| 100-40-3 | 4-vinylcyclohexene | no | no |
| 129879-84-1 | 5-amino-1,2,4-triazole-3-sulfonyl chloride | no | no |
| 97658-80-5 | 5-bp-bisenamine | no | no |
| 26172-55-4 | 5-chloro-2-methyl-4-isothiazolin-3-one | no | no |
| 91-44-1 | 7-diethylamino-4-methyl coumarin | no | no |
| 83-32-9 | acenaphthene | yes (POM) | no |
| 208-96-8 | acenaphthylene | yes (POM) | no |
| 75-07-0 | acetaldehyde | yes | yes |
| 64-19-7 | acetic acid | no | no |
| 108-24-7 | acetic anhydride | no | no |
| 67-64-1 | acetone | no | no |
| 75-05-8 | acetonitrile | yes | no |
| 98-86-2 | acetophenone | yes | no |
| 75-36-5 | acetyl chloride | no | no |
| 50-78-2 | acetylsalicylic acid | no | no |
| 107-02-8 | acrolein | yes | no |
| 79-06-1 | acrylamide | yes | yes |
| 79-10-7 | acrylic acid | yes | no |
| 107-13-1 | acrylonitrile | yes | yes |
| 814-68-6 | acryloyl chloride | no | no |
| 44992-01-0 | acryloyloxyethyltrimethyl ammonium chloride | no | no |
| 126-86-3 | actylenic diol | no | no |
| --0 | ad acid | no | no |
| 1330-86-5 | adipate plasticizer | no | no |
| 75782-86-4 | alcohols c12-13 | no | no |
| 69013-18-9 | alcohols c8-18 ethoxylated propoxylated | no | no |
| 309-00-2 | aldrin | no | yes |
| 3779-63-3 | aliphatic polyisocyanate-1 | no | no |
| 68515-40-2 | alkyl benzyl phthalate | no | no |
| 107-18-6 | allyl alcohol | no | no |
| 300-57-2 | allyl benzene | no | no |
| 107-05-1 | allyl chloride | yes | no |
| 106-92-3 | allyl glycidyl ether | no | yes |
| 532-27-4 | alpha chloroacetophenone | yes | no |
| 9000-90-2 | alpha-amylase | no | no |
| 552-45-4 | alpha-chloro-ortho-xylene | no | no |
| 319-84-6 | alpha-hexachlorocyclohexane | yes | yes |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 98-83-9 | alpha-methyl styrene | no | no |
| 68037-76-3 | alphamethylstyrene(dodecyl)polysiloxane | no | no |
| 109-06-8 | alpha-picoline | no | no |
| 7446-70-0 | aluminum chloride | no | no |
| 24304-00-5 | aluminum nitride | no | no |
| 60304-36-1 | aluminum potassium fluoride | no | no |
| 6419-19-8 | aminotrimethylene phosphonic acid | no | no |
| 7664-41-7 | ammonia | no | no |
| 12125-02-9 | ammonium chloride | no | no |
| 16919-31-6 | ammonium hexafluorozirconate | no | no |
| 1336-21-6 | ammonium hydroxide | no | no |
| 12054-85-2 | ammonium molybdate | no | no |
| 60676-86-0 | amorphous fused silica | no | no |
| --0 | amyl acetate (mixture) | no | no |
| 71-41-0 | amyl alcohol | no | no |
| 110-58-7 | amylamine | no | no |
| 513-35-9 | amylene | no | no |
| 63-05-8 | androstenedione | no | no |
| 63937-30-4 | anhydro-dimethylamino hexose reductone | no | no |
| 62-53-3 | aniline | yes | no |
| 120-12-7 | anthracene | yes (POM) | no |
| 7440-36-0 | antimony | yes (Sb comps.) | no |
| 28300-74-5 | antimony potassium tartrate | yes (Sb comps.) | no |
| 10025-91-9 | antimony trichloride | yes (Sb comps.) | no |
| 1309-64-4 | antimony trioxide | yes (Sb comps.) | no |
| 1345-04-6 | antimony trisulfide | yes (Sb comps.) | no |
| 68477-31-6 | aromatic petroleum derivative solvent | unknown | no |
| 7440-38-2 | arsenic | yes (As comps) | yes |
| 7784-42-1 | arsine | yes | no |
| 1332-21-4 | asbestos | yes | yes |
| 136816-75-6 | atevirdane mesylate | no | no |
| --0 | atlox 848 | no | no |
| 103-33-3 | azobenzene | no | yes |
| 9001-92-7 | bacillus subtilis neutral protease | no | no |
| 7440-39-3 | barium | no | no |
| 13701-59-2 | barium metaborate monohydrate | no | no |
| 7727-43-7 | barium sulfate | no | no |
| 147-24-0 | benadryl hcl | unknown | no |
| 56-55-3 | benz(a)anthracene | yes (POM) | no |
| 100-52-7 | benzaldehyde | no | yes |
| 71-43-2 | benzene | yes | yes |
| 91-01-0 | benzhydrol | no | no |
| 92-87-5 | benzidine | yes | yes |
| 50-32-8 | benzo(a)pyrene | yes (POM) | yes |
| 205-99-2 | benzo(b)fluoranthene | yes (POM) | no |
| 191-24-2 | benzo(g,h,i)perylene | yes (POM) | no |
| 207-08-9 | benzo(k)fluoranthene | yes (POM) | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 694-87-1 | benzocyclobutene | no | no |
| 119-53-9 | benzoin | no | no |
| 95-16-9 | benzothiazole | no | no |
| 25973-55-1 | benzotriazol dimethylpropyl phenol | unknown | no |
| 100-51-6 | benzyl alcohol | no | no |
| 100-44-7 | benzyl chloride | yes | yes |
| 103-83-3 | benzyl dimethylamine | no | no |
| 100-46-9 | benzylamine | no | no |
| 100-85-6 | benzyltrimethylammonium hydroxide | no | no |
| 7440-41-7 | beryllium | yes (Be comps) | yes |
| 64-04-0 | beta phenylethylamine | no | no |
| 126-99-8 | beta-chloroprene | yes | yes |
| 981-34-0 | betamethasone 11 | no | no |
| --0 | biosam tp-1.5 | no | no |
| 92-52-4 | biphenyl | yes | no |
| 3033-62-3 | bis (2-dimethylaminoethyl) ether | no | no |
| 108-60-1 | bis(2-chloroisopropyl)ether | no | no |
| 102054-10-4 | bis(2-methoxy-1-methylethy | no | no |
| 542-88-1 | bis(chloromethyl)ether | yes | yes |
| 13528-93-3 | bis(me2clsilyl)ethane | no | no |
| 41556-26-7 | bis(pentamethylpiperdinyl)sebacate | no | no |
| 111-44-4 | bis-2-chloroethylether | yes | yes |
| 60966-36-1 | bisnoralcohol | no | no |
| 25068-38-6 | bisphenol a/epichlorohydrin resin | no | no |
| 25085-99-8 | bisphenol epoxy resin | no | no |
| 2467-02-9 | bisphenolf | no | no |
| 10097-09-3 | bis-urea accelerator | no | no |
| 7637-07-2 | boron trifluoride | no | no |
| 68515-44-6 | branched and linear diheptyl phthalate ester | no | no |
| 111381-89-6 | branched and linear heptyl nonyl phthalate ester | no | no |
| 7726-95-6 | bromine | no | no |
| 108-86-1 | bromobenzene | no | no |
| 75-27-4 | bromodichloromethane | no | yes |
| 75-25-2 | bromoform | yes | yes |
| 56741-95-8 | bropirimine | no | no |
| 69102-90-5 | butadiene homopolymer | no | no |
| 106-97-8 | butane | no | no |
| 4435-53-4 | butoxyl | no | no |
| 141-32-2 | butyl acrylate | no | no |
| 85-68-7 | butyl benzyl phthalate | no | no |
| 112-34-5 | butyl carbitol | yes | no |
| 592-84-7 | butyl formate | no | no |
| 138-22-7 | butyl lactate | no | no |
| 143-29-3 | butylcarbitol formal | no | no |
| 102-79-4 | butyldiethanolamine | no | no |
| 123-72-8 | butyraldehyde | no | no |
| 107-92-6 | butyric acid | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 6408-78-2 | c.i. acid blue 25 | no | no |
| 108419-35-8 | c11-14 branched alkyl acetates | no | no |
| 78330-21-9 | c11-c14 isoalcohols, c14 rich, ethoxylated alcohol | no | no |
| 70914-20-4 | c6-8 branched alcohols | no | no |
| 7440-43-9 | cadmium | yes (Cd comps.) | yes |
| 7789-82-4 | calcium molybdate | no | no |
| 1592-23-0 | calcium stearate | no | no |
| 79-92-5 | camphene | no | no |
| 105-60-2 | caprolactam | no | no |
| 86-74-8 | carbazole | yes (POM) | yes |
| 1333-86-4 | carbon black | no | no |
| 75-15-0 | carbon disulfide | yes | no |
| 56-23-5 | carbon tetrachloride | yes | yes |
| 353-50-4 | carbonyl fluoride | no | no |
| 463-58-1 | carbonyl sulfide | yes | no |
| 9004-32-4 | carboxymethyl cellulose | no | no |
| 13466-78-9 | carene, delta | no | no |
| 8001-79-4 | castor oil | no | no |
| 120-80-9 | catechol | yes | no |
| 103980-44-5 | ceftiofur hydrochloride | no | no |
| 7440-45-1 | cerium | no | no |
| 1306-38-3 | cerium oxide | no | no |
| 123-03-5 | cetylpyridinium chloride | no | no |
| 6004-24-6 | cetylpyridinium chloride monohydrate | no | no |
| 12789-03-6 | chlordane (technical) | unknown | yes |
| 63449-39-8 | chlorinated paraffins | no | yes |
| 7782-50-5 | chlorine | yes | no |
| 10049-04-4 | chlorine dioxide | no | no |
| 302-22-7 | chlormadinone acetate | no | no |
| 108-90-7 | chlorobenzene | yes | no |
| 74-97-5 | chlorobromomethane | no | no |
| 57-15-8 | chlorobutanol | no | no |
| 124-48-1 | chlorodibromomethane | no | yes |
| 75-45-6 | chlorodifluoromethane | no | no |
| 668-45-1 | chlorofluorobenzonitrile | no | no |
| 67-66-3 | chloroform | yes | yes |
| 2921-88-2 | chlorpyrifos | no | no |
| 1308-14-1 | chromium (+3) hydroxide | no | no |
| 1308-38-9 | chromium 3 oxide | no | no |
| 218-01-9 | chrysene | yes (POM) | no |
| 18300-89-5 | cinnamate | no | no |
| 156-59-2 | cis-1-2, dichloroethylene | no | no |
| 627-20-3 | cis-2-pentene | no | no |
| 64741-62-4 | clarified oils (petroleum), catalytic cracked | no | no |
| 24729-96-2 | clindamycin phosphate | no | no |
| 1702-17-6 | clopyralid | no | no |
| 7440-48-4 | cobalt | yes (Co comps.) | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 136-52-7 | cobalt 2-ethylhexanoate | no | no |
| 27253-31-2 | cobalt neodecanoate | yes (Co comps.) | no |
| 61788-93-0 | coco alkyldimethyl amines | no | no |
| 8007-45-2 | coke oven emissions | yes | yes |
| 8050-09-7 | colophony | no | no |
| 7440-50-8 | copper | no | no |
| 3251-23-8 | copper nitrate | no | no |
| 147-14-8 | copper phthalocyanine | no | no |
| 7758-99-8 | copper sulfate pentahydrate | no | no |
| 7758-98-7 | copper sulfate, anhydrous | no | no |
| 68132-02-5 | coumarone indene resin | no | no |
| 7440-47-3 | Cr | yes ( Cr comps.) | no |
| 18540-29-9 | Cr , hexavalent - mist | yes (Cr comps.) | yes |
| 18540-29-9 | Cr , hexavalent - particulate | yes (Cr comps.) | yes |
| 16065-83-1 | Cr , trivalent | yes (Cr comps.) | no |
| 1319-77-3 | cresol (mixed isomers) | yes | no |
| 64265-57-2 | crosslinker cx100 | no | no |
| 4170-30-3 | crotonaldehyde | no | no |
| 98-82-8 | cumene | yes | yes |
| 80-15-9 | cumene hydroperoxide | no | no |
| 142-71-2 | cupric acetate | no | no |
| 1317-38-0 | cupric oxide (dust) | no | no |
| 57-12-5 | cyanide | yes as cyanides | no |
| 461-58-5 | cyanoguanidine | no | no |
| --0 | cyclic (phme)2(me)2, d4 | no | no |
| 2370-88-9 | cyclic methylhydrogensiloxane, d4 | no | no |
| 6166-86-5 | cyclic methylhydrogensiloxane, d5 | no | no |
| 2374-14-3 | cyclic methyltrifluoropropylsiloxane, d3 | no | no |
| 110-82-7 | cyclohexane | no | no |
| 108-94-1 | cyclohexanone | no | no |
| 12262-58-7 | cyclohexanone peroxide | no | no |
| 110-83-8 | cyclohexene | no | no |
| 6975-71-9 | cyclohexenylacetonitrile | no | no |
| 3399-73-3 | cyclohexenylethylamine | no | no |
| 1122-82-3 | cyclohexyl isothiocyanate | no | no |
| 287-92-3 | cyclopentane | no | no |
| 142-29-0 | cyclopentene | no | no |
| --0 | cyclopentyldichlorosilane | no | no |
| 14579-03-4 | cyclopentyltrichlorosilane | no | no |
| 147-94-4 | cytarabine | no | no |
| 72-54-8 | DDD (p, p'-dichlorodiphenyl dichloroethane) | no | yes |
| 72-55-9 | DDE ( $\mathrm{p}, \mathrm{p}$ '-dichlorodiphenyl dichloroethylene) | no | yes |
| 50-29-3 | DDT (p, $\mathrm{p}^{\prime}$-dichlorodiphenyl trichloroethane) | no | yes |
| 1163-19-5 | decabromodiphenyl oxide | unknown | yes |
| 91-17-8 | decahydronaphthalene | no | yes |
| 541-02-6 | decamethylcyclopentasiloxane | no | no |
| 141-62-8 | decamethyltetrasiloxane | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 8020-83-5 | deodorized kerosene | no | no |
| 64742-65-0 | dewaxed heavy paraffinic mineral oil | no | no |
| 6700-34-1 | dextromethorphan hydrochloride | no | no |
| 103-23-1 | di (2-ethylhexyl) adipate | no | yes |
| 6422-86-2 | di(ethylhexyl)terephthalate | no | no |
| 123-42-2 | diacetone alcohol | no | no |
| 131-17-9 | diallyl phthalate | no | yes |
| 2050-92-2 | diamylamine | no | no |
| 95481-62-2 | dibasic ester | no | no |
| 53-70-3 | dibenz(a,h)anthracene | yes (POM) | no |
| 132-64-9 | dibenzofuran | yes | no |
| 96-12-8 | dibromochloropropane | yes | no |
| 107-66-4 | dibutyl phosphate | no | no |
| 107-66-4 | dibutyl phosphate | no | no |
| 84-74-2 | dibutyl phthalate | yes | no |
| 77-58-7 | dibutyl tin dilaurate | no | no |
| 111-92-2 | dibutylamine | no | no |
| 818-08-6 | dibutyltin oxide | no | no |
| 91-94-1 | dichlorobenzidine | yes | yes |
| 75-71-8 | dichlorodifluoromethane | no | no |
| 75-43-4 | dichlorofluoromethane | no | no |
| 4109-96-0 | dichlorosilane | no | no |
| 76-14-2 | dichlorotetrafluoroethan | no | no |
| 62-73-7 | dichlorvos | yes | no |
| 80-43-3 | dicumyl peroxide | no | no |
| 5124-30-1 | dicyclohexylmethane-4,4'-diisocyanate | no | no |
| --0 | dicyclopentyldichlorosilane | no | no |
| 60-57-1 | dieldrin | no | yes |
| 68334-30-5 | diesel fuel | no | no |
| 111-42-2 | diethanolamine | yes | no |
| 117-81-7 | diethyl hexyl phthalate | yes | yes |
| 84-66-2 | diethyl phthalate | no | no |
| 64-67-5 | diethyl sulfate | yes | no |
| 109-89-7 | diethylamine | no | no |
| 25340-17-4 | diethylbenzene mixture | no | no |
| 111-46-6 | diethylene glycol | yes | no |
| 1559-36-0 | diethylene glycol mono-2-ethylhexyl ether | no | no |
| 124-17-4 | diethylene glycol monobutyl ether acetate | yes | no |
| 111-90-0 | diethylene glycol monoethyl ether | yes | no |
| 112-15-2 | diethylene glycol monoethyl ether acetate | yes | no |
| 111-77-3 | diethylene glycol monomethyl ether | yes | no |
| 104-68-7 | diethylene glycol monophenyl ether | yes | no |
| 111-40-0 | diethylene triamine | no | no |
| 68610-11-7 | diethylenetriamine reaction product with bisphenol a | no | no |
| 105-53-3 | diethylmalonate | no | no |
| 2238-07-5 | diglycidyl ether | no | no |
| 25036-25-3 | diglycidyl ether of bisphenol a | unknown | no |

# Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels 

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 26142-30-3 | diglycidyl ether of polyglycol | no | no |
| 108-83-8 | diisobutyl ketone | no | no |
| 107-39-1 | diisobutylene | no | no |
| 26761-40-0 | diisodecyl ester phthalate | no | no |
| 29733-18-4 | diisodecyl glutarate | no | no |
| 71888-89-6 | diisoheptyl phthalate | no | no |
| 28553-12-0 | diisononyl phthalate | no | no |
| 110-97-4 | diisopropanolamine | no | no |
| 108-20-3 | diisopropyl ether | no | no |
| 108-18-9 | diisopropylamine | no | no |
| 96-80-0 | diisopropylaminoethanol | no | no |
| 68390-56-7 | diketene hydrogenated fatty acids | no | no |
| 57-41-0 | dilantin | unknown | yes |
| 624-92-0 | dimethyldisulfide | no | no |
| 627-93-0 | dimethyl adipate | no | no |
| 106-79-6 | dimethyl decanedioate | no | no |
| 115-10-6 | dimethyl ether | no | no |
| 1119-40-0 | dimethyl glutarate | no | no |
| 756-79-6 | dimethyl methyl phosphonate | no | no |
| 106-65-0 | dimethyl succinate | no | no |
| 77-78-1 | dimethyl sulfate | yes | no |
| 124-40-3 | dimethylamine | no | no |
| 25988-97-0 | dimethylamine-epichlorohydrin polymer | no | no |
| 108-16-7 | dimethylamino-2-propanol | no | no |
| 121-69-7 | dimethylaniline | yes | yes |
| 1066-35-9 | dimethylchlorosilane | no | no |
| 28729-52-4 | dimethylcyclopentane | no | no |
| 75-78-5 | dimethyldichlorosilane | no | no |
| 1112-39-6 | dimethyldimethoxysilane | no | no |
| 2627-97-6 | dimethyldiphenydivinylsiloxane | no | no |
| 108-01-0 | dimethylethanolamine | no | no |
| 996-35-0 | dimethylisopropylamine | no | no |
| 51200-87-4 | dimethyloxazolidine | no | no |
| 131-11-3 | dimethylphthalate | yes | no |
| 1111-74-6 | dimethylsilane | no | no |
| 75-18-3 | dimethylsulfide | no | no |
| 67-68-5 | dimethylsulfoxide | no | no |
| 513-37-1 | dimethylvinyl chloride | no | yes |
| 1719-58-0 | dimethylvinylchlorosilane | no | no |
| 117-84-0 | di-n-octyl phthalate | no | no |
| 88-85-7 | dinoseb | no | no |
| 142-84-7 | di-n-propylamine | no | no |
| 838-85-7 | diphenyl phosphoric acid | no | no |
| 122-39-4 | diphenylamine | unknown | no |
| 144-79-6 | diphenylmethylchlorosilane | unknown | no |
| 778-25-6 | diphenylmethylsilanol | unknown | no |
| 101-84-8 | diphenyloxide | unknown | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 1231-93-0 | dipropyl ketone | no | no |
| 25265-71-8 | dipropylene glycol | no | no |
| 111109-77-4 | dipropylene glycol dimethyl ether | no | no |
| 34590-94-8 | dipropylene glycol methyl ether | no | no |
| 88917-22-0 | dipropylene glycol methyl ether acetate | no | no |
| 29911-28-2 | dipropylene glycol monobutyl ether | no | no |
| 29911-27-1 | dipropylene glycol monopropyl ether | no | no |
| 51730-94-0 | dipropylene glycol phenyl ether | yes | no |
| 4444-67-1 | di-sec-butylamine | no | no |
| 1590-87-0 | disilane | no | no |
| --0 | disiloxane | no | no |
| 123312-54-9 | distearyldimethylammonium bisulfate | no | no |
| 3843-16-1 | distearyldimethylammonium methosulfate | no | no |
| 64741-89-5 | distillates (petroleum) solvent-refined light paraffinic | no | no |
| 64742-30-9 | distillates (petroleum), chemically neutralized middle | no | no |
| 68410-00-4 | distillates (petroleum), crude oil | no | no |
| 64741-81-7 | distillates (petroleum), heavy thermal cracked | no | no |
| 64741-82-8 | distillates (petroleum), light thermal cracked | no | no |
| 64741-59-9 | distillates, (petroleum), light catalytic cracked | no | no |
| 68783-24-4 | di-tallow alkylamines | no | no |
| 330-54-1 | diuron | no | no |
| 1321-74-0 | divinyl benzene | no | no |
| 98-84-0 | dl-alpha phenylethylamine | no | no |
| 5989-27-5 | d-limonene | no | no |
| 540-97-6 | dodecamethylcyclohexasiloxane | no | no |
| 63148-57-2 | dow corning fluid 1107 | no | no |
| 35884-42-5 | dowanol dpnb | no | no |
| 170557-43-4 | dowanol tmh-deg borate ester | no | no |
| 145-73-3 | endothall | no | no |
| 106-89-8 | epichlorohydrin | yes | yes |
| --0 | epoxy resin solution | no | no |
| 50-28-2 | estradiol | no | no |
| 313-06-4 | estradiol cypionate | no | no |
| 141-43-5 | ethanolamine | no | no |
| --0 | ethomeen t/30 | no | no |
| 61791-12-6 | ethoxylated castor oil | no | no |
| 7085-85-0 | ethyl 2-cyanoacrylate | no | no |
| 141-78-6 | ethyl acetate | no | no |
| 141-97-9 | ethyl acetoacetate | no | no |
| 140-88-5 | ethyl acrylate | yes | no |
| 64-17-5 | ethyl alcohol | no | no |
| 541-85-5 | ethyl amyl ketone | no | no |
| 75-00-3 | ethyl chloride | yes | no |
| 105-39-5 | ethyl chloroacetate | no | no |
| 105-56-6 | ethyl cyanoacetate | no | no |
| 60-29-7 | ethyl ether | no | no |
| 109-94-4 | ethyl formate | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 97-64-3 | ethyl lactate | no | no |
| 106-68-3 | ethyl sec-amyl ketone | no | no |
| 78-10-4 | ethyl silicate | no | no |
| 637-92-3 | ethyl tertiary butyl ether | no | no |
| 25550-14-5 | ethyl toluene -mixture | no | no |
| 109-92-2 | ethyl vinyl ether | no | no |
| 763-69-9 | ethyl-3-ethyloxypropionate | no | no |
| 107-00-6 | ethylacetylene | no | no |
| 75-04-7 | ethylamine | no | no |
| 100-41-4 | ethylbenzene | yes | yes |
| 68987-42-8 | ethylenated benzene residues | no | no |
| 74-85-1 | ethylene | no | no |
| 142-59-6 | ethylene bisthiocarbamate disodium | no | no |
| 107-15-3 | ethylene diamine | no | no |
| 106-93-4 | ethylene dibromide | yes | yes |
| 107-21-1 | ethylene glycol | yes | no |
| 112-48-1 | ethylene glycol dibutyl ether | yes | no |
| 110-71-4 | ethylene glycol dimethyl ether | no | no |
| 1559-35-9 | ethylene glycol mono-2-ethylhexyl ether | no | no |
| 112-07-2 | ethylene glycol monobutyl ether acetate | yes | no |
| 111-15-9 | ethylene glycol monoethyl ether acetate | yes | no |
| 112-25-4 | ethylene glycol monohexyl ether | yes | no |
| 110-49-6 | ethylene glycol monomethyl ether acetate | yes | no |
| 122-99-6 | ethylene glycol monophenyl ether | yes | no |
| 2807-30-9 | ethylene glycol monopropyl ether | yes | no |
| 75-21-8 | ethylene oxide | yes | yes |
| 96-45-7 | ethylene thiourea | yes | yes |
| 64-02-8 | ethylenediamine tetra-acetic acid, tetrasodium salt | no | no |
| 9004-58-4 | ethylhydroxyethyl cellulose | no | no |
| 78-07-9 | ethyltriethoxysilane | no | no |
| 5314-55-6 | ethyltrimethoxysilane | no | no |
| 7525-62-4 | ethylvinyl benzene | no | no |
| 64742-06-9 | extracts (petroleum), middle distillate solvent | no | no |
| 108419-34-7 | exxate 1000 | no | no |
| 108419-32-5 | exxate 800 - octyl acetate | no | no |
| 108419-33-6 | exxate 900 | no | no |
| 68459-31-4 | fatty acids c9-11 branched glycidyl esters polymer | no | no |
| 136797-56-3 | fluorochemical-247 | unknown | no |
| 7705-08-0 | ferric chloride | no | no |
| 24510-87-0 | flumethasone 5 | no | no |
| 2476-74-6 | flumethasone 6 | no | no |
| 98967-40-9 | flumetsulam | no | no |
| 16872-11-0 | fluoboric acid | no | no |
| 206-44-0 | fluoranthene | yes (POM) | no |
| 86-73-7 | fluorene | yes (POM) | no |
| 69991-67-9 | fomblin perfluorpolyether | no | no |
| 50-00-0 | formaldehyde | yes | yes |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 75-12-7 | formamide | no | no |
| 64-18-6 | formic acid | no | no |
| 110-00-9 | furan | no | yes |
| 98-01-1 | furfural | no | yes |
| 98-00-0 | furfuryl alcohol | no | yes |
| --0 | fyre-zyme | no | no |
| 96-48-0 | gamma-butyrolactone | no | no |
| 8006-61-9 | gasoline | yes some comps | yes |
| 1310-53-8 | germanium dioxide | no | no |
| 7782-65-2 | germanium tetrahydride | no | no |
| 111-30-8 | glutaraldehyde | no | no |
| 56-81-5 | glycerol | no | no |
| 106-91-2 | glycidyl methacrylate | no | no |
| 93-14-1 | guaifenesin | no | no |
| 64741-65-7 | heavy alkylate naphtha | unknown | no |
| 64742-94-5 | heavy aromatic solvent naphtha | unknown | no |
| 64741-68-0 | heavy catalytic reformed naphtha | no | no |
| 68551-17-7 | heavy naphtha | no | no |
| 76-44-8 | heptachlor | yes | yes |
| --0 | heptamethyl-1-vinyl-1,7-dichlorotetrasilazane | no | no |
| 1873-88-7 | heptamethyltrisiloxane | no | no |
| 142-82-5 | heptane | no | no |
| 118-74-1 | hexachlorobenzene | yes | yes |
| 87-68-3 | hexachlorobutadiene | yes | yes |
| 77-47-4 | hexachlorocyclopentadiene | yes | no |
| 13465-77-5 | hexachlorodisilane | no | no |
| 67-72-1 | hexachloroethane | yes | yes |
| 12021-95-3 | hexafluorozirconium acid | no | no |
| 1009-93-4 | hexamethylcyclotrisilazane | no | no |
| 541-05-9 | hexamethylcyclotrisiloxane | no | no |
| 107-46-0 | hexamethyldisiloxane | no | no |
| 822-06-0 | hexamethylene diisocyanate | yes | no |
| 100-97-0 | hexamethylenetetramine | no | no |
| 66-25-1 | hexanaldehyde | no | no |
| 69696-98-6 | hexane 1,6-bis(tributyl ammonium bromi | no | no |
| 144669-03-4 | hexenylsiloxane | no | no |
| 144669-04-5 | hexenylsiloxanes | no | no |
| 107-41-5 | hexylene glycol | no | no |
| 431-89-0 | hfc-227ea | no | no |
| 68037-88-7 | high molecular weight sili | no | no |
| 63148-62-9 | high molecular wt. silicon | no | no |
| 68037-58-1 | high molecular wt. silicon | no | no |
| 68083-19-2 | high molecular wt. silicon | no | no |
| 68918-22-9 | high molecular wt. silicon | no | no |
| 69430-24-6 | high molecular wt. silicon | no | no |
| 70131-67-8 | high molecular wt. silicon | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 302-01-2 | hydrazine | yes | yes |
| 10034-93-2 | hydrazine sulfate | no | yes |
| 68956-56-9 | hydrocarbons, terpene processing by-products | no | no |
| 50-03-3 | hydrocortisone acetate | no | no |
| 64742-81-0 | hydrodesulfurized kerosene | no | no |
| 64742-80-9 | hydrodesulfurized middle distillate | unknown | no |
| 10035-10-6 | hydrogen bromide | no | no |
| 7647-01-0 | hydrogen chloride | yes | no |
| 74-90-8 | hydrogen cyanide | yes | no |
| 7664-39-3 | hydrogen fluoride | yes | no |
| 7722-84-1 | hydrogen peroxide | no | no |
| 7783-06-4 | hydrogen sulfide | no | no |
| 64742-48-9 | hydrotreated heavy napht | unknown | no |
| 64742-52-5 | hydrotreated heavy naphthenic distillate | unknown | no |
| 64742-54-7 | hydrotreated heavy paraffinic mineral oil | unknown | no |
| 64742-47-8 | hydrotreated light distillate | unknown | no |
| 64742-49-0 | hydrotreated light naphtha | no | no |
| 64742-53-6 | hydrotreated light naphthenic distillate | no | no |
| 64742-55-8 | hydrotreated light paraffinic distillate | unknown | no |
| 64742-46-7 | hydrotreated middle distillate | unknown | no |
| 79-14-1 | hydroxyacetic acid/ glycolic acid | no | no |
| 34375-28-5 | hydroxymethylamino ethanol | no | no |
| 10096-91-0 | hydroxyphenylbenzotriazole | no | no |
| 999-61-1 | hydroxypropyl acrylate | no | no |
| 9016-45-9 | igepal co-630 | no | no |
| 193-39-5 | indeno(1,2,3-cd)pyrene | yes (POM) | no |
| 7553-56-2 | iodine | no | no |
| 123-92-2 | isoamyl acetate | no | no |
| 123-51-3 | isoamyl alcohol | no | no |
| 5888-33-5 | iso-bornyl acrylate | no | no |
| 75-28-5 | isobutane | no | no |
| 110-19-0 | isobutyl acetate | no | no |
| 78-83-1 | isobutyl alcohol | no | no |
| 97-85-8 | isobutyl isobutyrate | no | no |
| 97-86-9 | isobutyl methacrylate | no | no |
| 115-11-7 | isobutylene | no | no |
| 18395-30-7 | isobutyltrimethoxysilane | no | no |
| 78-84-2 | isobutyraldehyde | no | no |
| 79-31-2 | isobutyric acid | no | no |
| 338-98-7 | isoflupredone acetate | no | no |
| 26952-21-6 | isooctanol | no | no |
| 90622-57-4 | isopar h | no | no |
| 78-59-1 | isophorone | yes | yes |
| 4098-71-9 | isophorone diisocyanate | no | no |
| 53880-05-0 | isophorone diisocyanate polymer | no | no |
| 78-79-5 | isoprene | no | yes |
| 108-21-4 | isopropyl acetate | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 67-63-0 | isopropyl alcohol | no | no |
| 75-31-0 | isopropylamine | no | no |
| 121-93-7 | isopropyldiethanolamine | no | no |
| 109-56-8 | isopropylethanolamine | no | no |
| 590-86-3 | isovaleraldehyde | no | no |
| 1332-58-7 | kaolin | no | no |
| 50-21-5 | lactic acid | no | no |
| 2627-86-3 | I-alpha-phenylethylamine | no | no |
| 39464-66-9 | lauryl alcohol, phosphated | no | no |
| 10190-55-3 | lead molybdate | no | no |
| 64741-66-8 | light alkylate naphtha | no | no |
| 64742-95-6 | light aromatic solvent naphtha (petroleum) | unknown | no |
| 859-18-7 | lincomycin hydrochloride | no | no |
| 141-63-9 | linear dimethylsiloxanes,md3m(\&higher) | no | no |
| 68083-20-5 | linear methylvinylsiloxane ppolymer hydroxl endblock | no | no |
| 67762-41-8 | linear primary alcohol | no | no |
| 1345-05-7 | lithopone | no | no |
| 66071-86-1 | Iv 837/821 | no | no |
| 7439-95-4 | magnesium | no | no |
| 546-93-0 | magnesium carbonate | no | no |
| 7786-30-3 | magnesium chloride | no | no |
| 1309-42-8 | magnesium hydroxide | no | no |
| 10377-60-3 | magnesium nitrate | no | no |
| 1309-48-4 | magnesium oxide | no | no |
| 557-04-0 | magnesium stearate | no | no |
| 108-31-6 | maleic anhydride | yes | no |
| 6915-15-7 | malic acid | no | no |
| 591-27-5 | m-aminophenol | no | no |
| 7439-96-5 | manganese | yes (Mn comps.) | no |
| 15956-58-8 | manganese 2-ethylhexanoate | yes (Mn comps.) | no |
| 27253-32-3 | manganese neodecanoate | yes (Mn comps.) | no |
| 1317-35-7 | manganese oxide | yes (Mn comps.) | no |
| 10034-96-5 | manganese sulfate monohydrate | yes (Mn comps.) | no |
| 26544-20-7 | mcpa 2-ehe (2-methyl-4-chlorophenoxyacetic acid 2-ethylhexyl ester) | no | no |
| 644-62-2 | meclofenamic acid | unknown | no |
| 108-78-1 | melamine | no | yes |
| 2919-66-6 | melengesterol acetate | no | no |
| 7439-97-6 | mercury | yes (Hg comps.) | no |
| 141-79-7 | mesityl oxide | no | no |
| 79-41-4 | methacrylic acid | no | no |
| 75-75-2 | methane sulfonic acid | no | no |
| 3144-09-0 | methanesulfonamide | no | no |
| 67-56-1 | methanol | yes | no |
| 79-20-9 | methyl acetate | no | no |
| 74-99-7 | methyl acetylene | no | no |
| 96-33-3 | methyl acrylate | no | no |
| 108-11-2 | methyl amyl alcohol | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 93-58-3 | methyl benzoate | no | no |
| 74-83-9 | methyl bromide | yes | no |
| 115-19-5 | methyl butynol | no | no |
| 74-87-3 | methyl chloride | yes | yes |
| 71-55-6 | methyl chloroform | yes | no |
| 17639-93-9 | methyl chloroproprionate | no | no |
| 78-93-3 | methyl ethyl ketone | yes | no |
| 1338-23-4 | methyl ethyl ketone peroxide | no | no |
| 107-31-3 | methyl formate | no | no |
| 60-34-4 | methyl hydrazine | yes | yes |
| 110-12-3 | methyl isoamyl ketone | no | no |
| 108-10-1 | methyl isobutyl ketone | yes | no |
| 624-83-9 | methyl isocyanate | yes | no |
| 74-93-1 | methyl mercaptan | no | no |
| 80-62-6 | methyl methacrylate | yes | no |
| 1184-85-6 | methyl methane sulfonamide | no | no |
| 110-43-0 | methyl $n$-amyl ketone | no | no |
| 591-78-6 | methyl n-butyl ketone | no | no |
| 9003-11-6 | methyl oxirane (pluronic p103) | no | no |
| 82919-37-7 | methyl pentamethyl-4-piperidinyl ester of decanedioic acid | no | no |
| 53-36-1 | methyl predisolone acetate | no | no |
| 107-87-9 | methyl propyl ketone | no | no |
| 124-63-0 | methyl sulfonyl chloride | no | no |
| 1634-04-4 | methyl t-butyl ether | yes | no |
| 109-87-5 | methylal | no | no |
| 74-89-5 | methylamine | no | no |
| 593-51-1 | methylamine hydrochloride | no | no |
| 108-87-2 | methylcyclohexane | no | no |
| 75-54-7 | methyldichlorosilane | no | no |
| 105-59-9 | methyldiethanolamine | no | no |
| 16881-77-9 | methyldimethoxysilane | no | no |
| 75-09-2 | methylene chloride | yes | yes |
| 101-68-8 | methylene diphenyl isocyanate | yes | no |
| 96-29-7 | methylethylketoxime | no | yes |
| 992-94-9 | methylsilane | no | no |
| 999-97-3 | methylsilazane | no | no |
| 4253-34-3 | methyltriacetoxysilane | no | no |
| 75-79-6 | methyltrichlorosilane | no | no |
| 2031-67-6 | methyltriethoxysilane | no | no |
| 50791-87-2 | methylvinylbis(n-methylace | no | no |
| 124-70-9 | methylvinyldichlorosilane | no | no |
| 16753-62-1 | methylvinyldimethoxysilane | no | no |
| 8012-95-1 | mineral oil | no | no |
| 64475-85-0 | mineral spirits | unknown | no |
| 1330-20-7 | mixed xylenes | yes | no |
| 7439-98-7 | molybdenum | no | no |
| 18868-43-4 | molybdenum dioxide | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 1317-33-5 | molybdenum disulfide | no | no |
| 1313-27-5 | molybdenum trioxide | no | yes |
| 46438-39-5 | monobutyl monophenyl phosphoric acid | no | no |
| 1623-15-0 | monobutyl phosphoric acid | no | no |
| 95-49-8 | monochlorotoluene | no | no |
| 78-96-6 | monoisopropanolamine | no | no |
| 701-64-4 | monophenyl phosphoric acid | no | no |
| 110-91-8 | morpholine | no | no |
| 28729-54-6 | m-propyl toluene | no | no |
| 620-23-5 | m-tolualdehyde | no | no |
| 98-17-9 | m-trifluoromethylphenol | no | no |
| 108-38-3 | m-xylene | yes | no |
| 126803-73-4 | n-(2,6-dichloro-3-methylphenyl)-5,7-dimethyoxy(1,2,4)triazo...[de-511] | no | no |
| 113171-12-3 | n -(2,6-difluorophenyl)-5-amino-1h-1,2,4-triazole-3-sulfonamide | no | no |
| 98967-55-6 | n-(2,6-difluorophenyl)-7-methyl-1h-1,2,4-triazolo(1,5a)pyrimidine-2-su | no | no |
| 1760-24-3 | n -(3-(trimethoxysilyl)propyl)-ethylenediamine | no | no |
| 104-78-9 | $\mathrm{n}, \mathrm{n}$-diethyl-1,3-propanediamine | no | no |
| 606-46-2 | $n, \mathrm{n}$-diethyl-o-toluine | no | no |
| 613-48-9 | $n, n$-diethyl-p-toludine | no | no |
| 124-28-7 | n ,n-dimethyl octadecylamine | no | no |
| 80-73-9 | $\mathrm{n}, \mathrm{n}$ '-dimethylethyleneurea | no | no |
| 68-12-2 | n,n-dimethylformamide | yes | no |
| 99-97-8 | n,n-dimethyl-p-toluidine | no | no |
| 110-30-5 | n , n --ethylene bis-octadecanamide | no | no |
| 628-63-7 | n -amyl acetate | no | no |
| 8030-30-6 | naphtha | no | no |
| 64742-82-1 | naphtha (petroleum) hydrodesulfurized heavy | no | no |
| 64741-55-5 | naphtha (petroleum), light catalytic cracked | no | no |
| 64741-41-9 | naphtha heavy straight run | unknown | no |
| 68955-35-1 | naphtha, catalytic reformed | no | no |
| 64741-64-6 | naphtha, full range alkylate | no | no |
| 64741-42-0 | naphtha, full range straight run | no | no |
| 64741-54-4 | naphtha, heavy catalytic cracked | no | no |
| 64741-83-9 | naphtha, heavy thermal cracked | no | no |
| 64741-63-5 | naphtha, light catalytic reformed | no | no |
| 91-20-3 | naphthalene | yes | yes |
| 71-36-3 | n-butanol | no | no |
| 63716-40-5 | n-butoxy propanol (mixed isomers) | no | no |
| 123-86-4 | n-butyl acetate | no | no |
| 109-69-3 | n-butyl chloride | no | no |
| 2426-08-6 | n-butyl glycidyl ether | no | no |
| 97-88-1 | n -butyl methacrylate | no | no |
| 590-01-2 | n-butyl propionate | no | no |
| 109-73-9 | n-butylamine | no | no |
| 104-51-8 | n-butylbenzene | no | no |
| --0 | n-butylglucamine | no | no |
| --0 | n -chloro-2,6-difluorobenzamide | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 112-55-0 | n-dodecyl mercaptan | no | no |
| 1405-10-3 | neomycin sulfate | no | no |
| 126-30-7 | neopentyl glycol | no | no |
| 17557-23-2 | neopentyl glycol diglycidyl ether | no | no |
| 112-06-1 | n-heptyl acetate | no | no |
| 110-54-3 | n-hexane | yes | no |
| 7440-02-0 | nickel | yes (Ni comps.) | yes |
| 12035-72-2 | nickel subsulfide | yes (Ni comps.) | yes |
| 7697-37-2 | nitric acid | no | no |
| 98-95-3 | nitrobenzene | yes | yes |
| 79-24-3 | nitroethane | no | no |
| 77835-42-0 | nitrogen trifluoride | no | no |
| 75-52-5 | nitromethane | no | yes |
| 107-68-6 | n-methyl taurine | no | no |
| 626-67-5 | n-methylpiperidine | no | no |
| 872-50-4 | n -methylpyrrolidone | no | no |
| 836-30-6 | n -nitrodiphenylamine | no | no |
| 1116-54-7 | n -nitrosodiethanolamine | no | yes |
| 621-64-7 | n-nitroso-di-n-propylamine | no | yes |
| 86-30-6 | n -nitrosodiphenylamine | unknown | yes |
| 684-93-5 | n-nitroso-n-methylurea | yes | yes |
| 111-84-2 | n -nonane | no | no |
| 25154-52-3 | nonyl phenol (mixed isomers) | no | no |
| 64771-72-8 | norpar 12 | no | no |
| 303-81-1 | novobiocin | unknown | no |
| 624-54-4 | n -pentyl proprionate | no | no |
| 109-60-4 | n-propyl acetate | no | no |
| 71-23-8 | n-propyl alcohol | no | no |
| 16369-21-4 | n-propylethanolamine | no | no |
| 103-99-1 | n -stearoyl-4-aminophenol | no | no |
| 88-12-0 | n -vinylpyrrolidinone | no | yes |
| 68309-52-4 | nylen 5 | unknown | no |
| --0 | o-(1-ethoxyethyl)-2-(propylthio)-3-(trifluoromethyl)phenol | no | no |
| --0 | o-(1-ethoxyethyl)-3-(trifluoromethyl)phenol | no | no |
| 134-29-2 | o-ansidine hydrochloride | no | yes |
| 88-65-3 | o-bromobenzoic acid | no | no |
| 118-91-2 | o-chlorobenzoic acid | no | no |
| 95-48-7 | o-cresol | yes | no |
| 124-26-5 | octadecanamide | no | no |
| 627-83-8 | octadecanoic acid, 1,2-ethanediyl ester | no | no |
| 27668-52-6 | octadecyldimethyl (3-(trimethoxysilyl)propyl) ammonium chloride | no | no |
| 556-67-2 | octamethylcyclotetrasilo | no | no |
| 107-51-7 | octamethyltrisiloxane | no | no |
| 124-07-2 | octanoic acid | no | no |
| 26530-20-1 | octylisothiazolone | no | no |
| 68990-79-4 | oils, vegetable, mixed with animal oil methylesters, polymerized, oxidixed | no | no |
| 112-80-1 | oleic acid | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 93-83-4 | oleoyl diethanolamine | no | no |
| 8014-95-7 | oleum | no | no |
| 90-43-7 | o-phenylphenol | unknown | yes |
| 2530-85-0 | organofunctional silane | no | no |
| 20816-12-0 | osmium tetroxide | no | no |
| 95-53-4 | o-toluidine | yes | yes |
| 19666-30-9 | oxadiazon | no | yes |
| 9063-06-3 | oxirane, methyl-, polymer with oxirane, monomethyl ether | no | no |
| 90438-79-2 | oxo-heptyl acetate | no | no |
| 88230-35-7 | oxo-hexyl acetate | no | no |
| 95-47-6 | o-xylene | yes | no |
| 95-38-5 | oyel hydroxyethylimidazoline | no | no |
| 7440-05-3 | palladium | no | no |
| 8002-74-2 | paraffin wax fume | no | no |
| 98-56-6 | p-chlorobenzotrifluoride | no | no |
| 82-68-8 | pentachloronitrobenzene | yes | no |
| 87-86-5 | pentachlorophenol | yes | yes |
| 109-66-0 | pentane | no | no |
| 19430-93-4 | perfluorobutylethylene | no | no |
| 38436-16-7 | perfluorobutylethylmethyldichlorosilane | no | no |
| 382-21-8 | perfluoroisobutylene | no | no |
| 93-59-4 | peroxybenzoic acid | no | no |
| 8002-05-9 | petroleum | no | no |
| 64742-14-9 | petroleum distillates, acid treated | unknown | no |
| 68476-86-8 | petroleum gases, liquefied, sweetened | no | no |
| 1194-02-1 | p-fluorobenzonitrile | no | no |
| 85-01-8 | phenanthrene | yes (POM) | no |
| 10551-21-0 | phenethyl alpha picolinium bromide | no | no |
| 108-95-2 | phenol | yes | no |
| 122-79-2 | phenyl acetate | no | no |
| 617-94-7 | phenyl isopropanol (2-phenyl-2-propanol) | no | no |
| 120-07-0 | phenyldiethanolamine | no | no |
| 98-13-5 | phenyltrichlorosilane | no | no |
| 2996-92-1 | phenyltrimethoxysilane | no | no |
| 75-44-5 | phosgene | yes | no |
| 7803-51-2 | phosphine | yes | no |
| 7664-38-2 | phosphoric acid | no | no |
| 7723-14-0 | phosphorus (total) | yes | no |
| 10025-87-3 | phosphorus oxychloride | no | no |
| 10026-13-8 | phosphorus pentachloride | no | no |
| 7719-12-2 | phosphorus trichloride | no | no |
| 1328-53-6 | phthalocyanine pigment green | no | no |
| 26952-20-5 | picloram, isooctyl ester | no | no |
| 80-56-8 | pinene, alpha | no | no |
| 127-91-3 | pinene, beta | no | no |
| 2981-10-4 | piperdinocyclohexene | no | no |
| 110-89-4 | piperidine | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 61477-94-9 | pirmenol hydrochloride | no | no |
| 99-87-6 | p-isopropyltoluene | no | no |
| 7440-06-4 | platinum soluble salt | no | no |
| 4221-98-1 | p-mentha-1,5-diene | no | no |
| 9016-87-9 | polmeric methylene diphenyl diisocyanate | no | no |
| 26780-96-1 | poly(1,2-dihydro-2,2,4-trimethylquinoline) | no | no |
| 9003-13-8 | polyalkylene glycol monobutyl ether/ butoxypolypropylene glycol | no | no |
| 68003-28-1 | polyamide | no | no |
| 1336-36-3 | polychlorinated biphenyls | yes | yes |
| --0 | polycyclic aromatic hydrocarbons (pahs) | yes (POM) | no |
| 26062-79-3 | polydimethyl diallyl ammonium chloride | no | no |
| 25322-68-3 | polyethylene glycol | no | no |
| 9004-74-4 | polyethylene glycol methyl ether | no | no |
| 27274-31-3 | polyethylene glycol monoallyl ether | no | no |
| 37251-67-5 | polyethylene polypropylene glycol | no | no |
| 68410-23-1 | polyethylenepolyamine reaction products with c18-unsat. fatty acids | no | no |
| 69029-39-6 | polyglycol 26-2 | no | no |
| --0 | polyglycol 26-3 | no | no |
| 24938-91-8 | polyglycol 59-13 | no | no |
| 9002-92-0 | polyoxyethylene lauryl ether | no | no |
| 25322-69-4 | polypropylene glycol | no | no |
| 9002-86-2 | polyvinyl chloride | no | no |
| 9003-39-8 | polyvinyl pyrrolidone | no | no |
| 9003-22-9 | polyvinylchloride/polyvinylacetate | no | no |
| 7789-23-3 | potassium fluoride | no | no |
| 1310-58-3 | potassium hydroxide | no | no |
| 7758-05-6 | potassium iodate | no | no |
| 12136-45-7 | potassium oxide | no | no |
| 7722-64-7 | potassium permanganate | yes (Mn comps.) | no |
| 12037-29-5 | praseodymium oxide | no | no |
| 57-83-0 | progesterone | no | no |
| 3986-89-8 | progesterone 4 | no | no |
| 98516-30-4 | propanol, 1(or 2) ethoxy, acetate isoparaffinic petroleum hydrocarbon | no | no |
| 123-38-6 | propionaldehyde | no | no |
| 79-09-4 | propionic acid | no | no |
| 106-94-5 | propyl bromide | no | no |
| 106-36-5 | propyl propionate | no | no |
| 107-10-8 | propylamine | no | no |
| 103-65-1 | propylbenzene | no | no |
| 115-07-1 | propylene | no | no |
| 108-32-7 | propylene carbonate | no | no |
| 78-87-5 | propylene dichloride | yes | no |
| 57-55-6 | propylene glycol | no | no |
| 19089-47-5 | propylene glycol monoethyl ether (alpha) | no | no |
| 1569-02-4 | propylene glycol monoethyl ether (beta) | no | no |
| 52125-53-8 | propylene glycol monoethyl ether (mixture) | no | no |
| 107-98-2 | propylene glycol monomethyl ether | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 1320-67-8 | propylene glycol monomethyl ether | no | no |
| 108-65-6 | propylene glycol monomethyl ether acetate | no | no |
| 41593-38-8 | propylene glycol monophenyl ether | no | no |
| 5131-66-8 | propylene glycol n-butyl ether (alpha isomer) | no | no |
| 15821-83-7 | propylene glycol n -butyl ether (beta isomer) | no | no |
| 770-35-4 | propylene glycol phenyl ether | no | no |
| 57018-52-7 | propylene glycol tert-butyl ether | no | no |
| 29387-86-8 | propylene glycol, n-butyl ether (mixed isomers) | no | no |
| 75-56-9 | propylene oxide | yes | yes |
| 1067-25-0 | propyltrimethoxysilane | no | no |
| 104-87-0 | p-tolualdehyde | no | no |
| 104-15-4 | p-toluenesulfonic acid | no | no |
| 6192-52-5 | p-toluenesulfonic acid monohydrate | no | no |
| 106-49-0 | p-toluidine | no | yes |
| --0 | purafect 4000 g | no | no |
| 106-42-3 | p-xylene | yes | no |
| 129-00-0 | pyrene | yes (POM) | no |
| 110-86-1 | pyridine | no | no |
| 84632-65-5 | pyrrolo[3,4-c]pyrrole-1,4-dione,3,6-bis(4-chlorophenyl)-2,5-dihydro | no | no |
| 1047-16-1 | quinacridone pigment | no | no |
| 82586-54-7 | quinapril step 8 | no | no |
| 91-22-5 | quinoline | yes | yes |
| 106-51-4 | quinone | yes | no |
| 64742-62-7 | residual oils (petroleum) solvent-dewaxed | no | no |
| 64741-56-6 | residues, (petroleum), vacuum | no | no |
| 108-46-3 | resorcinol | no | no |
| 1314-28-9 | rhenium oxide | no | no |
| 90-02-8 | salicylaldehyde | no | no |
| 106917-31-1 | sanduvor 3068 liquid | no | no |
| 3081-01-4 | santoflex 14 | no | no |
| 626-38-0 | sec-amyl acetate | no | no |
| 78-92-2 | sec-butyl alcohol | no | no |
| 13952-84-6 | sec-butylamine | no | no |
| 135-98-8 | sec-butylbenzene | no | no |
| 7782-49-2 | selenium | yes (Se comps.) | no |
| 112926-00-8 | silica - precipitated | no | no |
| 69012-64-2 | silica amorphous fume | no | no |
| 112945-52-5 | silica, amorphous, crystalline free, fumed | no | no |
| 10026-04-7 | silicon tetrachloride | no | no |
| 7783-61-1 | silicon tetrafluoride | no | no |
| 7803-62-5 | silicon tetrahydride | no | no |
| 67762-90-7 | siloxanes and silicones(silica filled polydimethylsiloxane) | no | no |
| 7440-22-4 | silver - soluble | no | no |
| 15096-52-3 | sodium aluminum fluoride | no | no |
| 7631-90-5 | sodium bisulfite | no | no |
| 7647-15-6 | sodium bromide | no | no |
| 630-93-3 | sodium dilantin | unknown | yes |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 128-04-1 | sodium dimethyl dithiocarbamate | no | no |
| 7681-49-4 | sodium fluoride | no | no |
| 13007-85-7 | sodium glucoheptonate | no | no |
| 31138-65-5 | sodium glucoheptonate | no | no |
| 1310-73-2 | sodium hydroxide | no | no |
| 7681-52-9 | sodium hypochlorite | no | no |
| 10039-56-2 | sodium hypophosphite monohydrate | no | no |
| 7681-82-5 | sodium iodide | no | no |
| 14960-06-6 | sodium lauriminodipropionate | no | no |
| 124-41-4 | sodium methylate | no | no |
| 7631-95-0 | sodium molybdate | no | no |
| 12401-86-4 | sodium monoxide | no | no |
| 7632-00-0 | sodium nitrite | no | no |
| 7632-04-4 | sodium perborate | no | no |
| 68608-26-4 | sodium petroleum sulfonate | unknown | no |
| 16893-85-9 | sodium silicofluoride | no | no |
| 67701-11-5 | sodium soap 900602 | no | no |
| 67701-10-4 | sodium soap 903923 | no | no |
| 7757-83-7 | sodium sulfite | no | no |
| 1300-72-7 | sodium xylenesulfonate | no | no |
| 109265-71-6 | solsperse 12000 | unknown | no |
| 86753-78-8 | solsperse 5000 | unknown | no |
| 68458-91-3 | solvar \& Iv 820 | no | no |
| 8005-02-5 | solvent black | no | no |
| 64742-96-7 | solvent naphtha (petroleum) heavy aliphatic | no | no |
| 64742-89-8 | solvent naphtha light aliphatic | unknown | no |
| 64742-88-7 | solvent naphtha medium aliphatic | unknown | no |
| 64741-88-4 | solvent refined heavy paraffnic distillate | unknown | no |
| 67784-80-9 | soybean oil, methyl esters | no | no |
| 68071-85-2 | spenkel f34 | unknown | no |
| --0 | sponto 11 | no | no |
| --0 | sponto 723 | no | no |
| 30705-14-7 | sr 1153 | no | no |
| 1912-83-0 | stannous octoate | no | no |
| 57-11-4 | stearic acid | no | no |
| 7803-52-3 | stibine | yes (Sb comps.) | no |
| 8052-41-3 | stoddard solvent | no | no |
| 64741-44-2 | straight run middle distillate | no | no |
| 100-42-5 | styrene | yes | yes |
| 9003-55-8 | styrene-butadiene polymer | no | no |
| 505-48-6 | suberic acid | no | no |
| 5329-14-6 | sulfamic acid | no | no |
| 7704-34-9 | sulfur (elemental) | no | no |
| 7446-11-9 | sulfur trioxide | no | no |
| 7664-93-9 | sulfuric acid | no | no |
| 68516-16-5 | sulfuric acid c6-10 alkyl esters | no | no |
| 64741-86-2 | sweetened middle distillate | unknown | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 14807-96-6 | talc | no | yes |
| 61790-33-8 | tallow alkylamines | no | no |
| 75-65-0 | t-butanol | no | no |
| 107-71-1 | t-butyl peroxyacetate | no | no |
| 75-64-9 | t-butylamine | no | no |
| 4620-70-6 | t-butylaminoethanol | no | no |
| 98-29-3 | t-butylcatechol | no | no |
| 2160-93-2 | t-butyldiethanolamine | no | no |
| 3006-82-4 | t-butylperoxy-2-ethylhexanoate | no | no |
| 9036-19-5 | t-det c08 | no | no |
| --0 | t-det c-40 | no | no |
| 9014-92-0 | t-det dd-14 | no | no |
| 68131-40-8 | tergitol 15-s-3 | no | no |
| 540-88-5 | tert-butyl acetate | no | no |
| 98-06-6 | tert-butylbenzene | no | no |
| 1333-13-7 | tert-butyl-m-cresol | no | no |
| 994-05-8 | tertiary amyl methyl ether | no | no |
| 2157-45-1 | tetra-2-methoxyethoxy-silane | no | no |
| 136-47-0 | tetracaine hyrochloride | no | no |
| 20536-16-7 | tetrachlorodisilane | no | no |
| 127-18-4 | tetrachloroethylene | yes | yes |
| 10469-09-7 | tetrachloropicolinic acid | no | no |
| 116-14-3 | tetrafluoroethylene | no | yes |
| 109-99-9 | tetrahydrofuran | no | yes |
| 97-99-4 | tetrahydrofuryl methanol | no | no |
| 9014-85-1 | tetramethyl decyndiol | no | no |
| 632-22-4 | tetramethyl urea | no | no |
| 22407-51-8 | tetramethylchlorovinyldisiloxane | no | no |
| 3277-26-7 | tetramethyldihydrogendisiloxane | no | no |
| 7691-02-3 | tetramethyldivinyldisila | no | no |
| 2627-95-4 | tetramethyldivinyldisiloxane | no | no |
| 75-76-3 | tetramethylsilane | no | no |
| 509-14-8 | tetranitromethane | no | yes |
| 3982-82-9 | tetraphenyldimethyl-2-dimethyltrisiloxane | unknown | no |
| 3390-61-2 | tetraphenyldimethyl-2-phenylmethyltrisiloxane | unknown | no |
| 807-28-3 | tetraphenyldimethyldisiloxane | unknown | no |
| 6904-66-1 | tetraphenylhexamethyltetrasiloxane | unknown | no |
| 25265-77-4 | texanol | no | no |
| 1314-32-5 | thallic oxide | no | no |
| 7440-28-0 | thallium | no | no |
| 64485-82-1 | thiazole ester | no | no |
| 7719-09-7 | thionyl chloride | no | no |
| 137-26-8 | thiram | no | no |
| 7440-31-5 | tin | no | no |
| 13463-67-7 | titanium dioxide | no | no |
| 1643-19-2 | t-n-butyl ammonium bromide | no | no |
| 1156-19-0 | tolazamide | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

 January 25, 2010(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| :---: | :---: | :---: | :---: |
| 108-88-3 | toluene | yes | no |
| 26471-62-5 | toluene diisocyanate | yes | yes |
| 8001-35-2 | toxaphene | yes | yes |
| 156-60-5 | trans-1-2,dichloroethylene | no | no |
| 102-76-1 | triacetin | no | no |
| 28961-43-5 | triacrylate ester | no | no |
| 621-77-2 | triamylamine | no | no |
| 126-73-8 | tributyl phosphate | no | no |
| 102-82-9 | tributylamine | no | no |
| 79-01-6 | trichloroethylene | yes | yes |
| 75-69-4 | trichlorofluoromethane | no | no |
| 10025-78-2 | trichlorosilane | no | no |
| 68526-86-3 | tridecanol | no | no |
| 102-71-6 | triethanolamine | no | no |
| 77-93-0 | triethyl citrate | no | no |
| 121-44-8 | triethylamine | yes | no |
| --0 | triethylammonium suleptanate | no | no |
| 1559-37-1 | triethylene glycol mono-2-ethyhexyl ether | no | no |
| 112-50-5 | triethylene glycol monoethyl ether | no | no |
| 112-24-3 | triethylene tetramine | no | no |
| 280-57-9 | triethylenediamine | no | no |
| 76-05-1 | trifluoroacetic acid | no | no |
| 358-67-8 | trifluoropropylmethyl dimethoxysilane | no | no |
| 130014-38-9 | trifluoropropylsilsesquioxane, dimethylhydrogensilyoxy-terminated | no | no |
| 592-09-6 | trifluoropropyltrichlorosilane | no | no |
| 122-20-3 | triisopropanolamine (tipa) | no | no |
| 121-43-7 | trimethoxyborine | no | no |
| 1185-55-3 | trimethoxymethylsilane | no | no |
| 3236-53-1 | trimethyl hexamethylenediamine | no | no |
| 75-50-3 | trimethylamine | no | no |
| 25551-13-7 | trimethylbenzenes (mixed isomers) | no | no |
| 75-77-4 | trimethylchlorosilane | no | no |
| 1445-45-0 | trimethyl-o-acetate | no | no |
| 3290-92-4 | trimethylolpropane trimethacrylate | no | no |
| 149-73-5 | trimethylorthoformate | no | no |
| 993-07-7 | trimethylsilane | no | no |
| 1066-40-6 | trimethylsilanol | no | no |
| 76-83-5 | triphenyl methyl chloride | unknown | no |
| 102-69-2 | tripropylamine | no | no |
| 42978-66-5 | tripropylene glycol diacrylate | no | no |
| 25498-49-1 | tripropylene glycol methyl ether | no | no |
| 20324-33-8 | tripropylene glycol methyl ether, dowanol 62b | no | no |
| 55934-93-5 | tripropylene glycol n-butyl ether | no | no |
| 126-72-7 | tris(2,3-dibromopropyl) phosphate | no | yes |
| 9002-93-1 | triton $\times 100$ | no | no |
| 88851-61-0 | trospectomycin sulfate | no | no |
| 51811-38-2 | tryfac 5556 | no | no |

## Appendix K - Comparison of HAP and TAC Screening Level Lists MDNRE AQD Toxic Air Contaminants with Health-Based Screening Levels

January 25, 2010
(EPA HAPs appear in BOLD type)

| CAS No. | Chemical Name | HAP? | Carcinogen? |
| ---: | :--- | :---: | :---: |
| $12070-12-1$ | tungsten carbide | no | no |
| $8006-64-2$ | turpentine | no | no |
| $110-62-3$ | valeraldehyde | no | no |
| $3153-26-2$ | vanadium oxide bis $(2,4-$ pentanedionate) | no | no |
| $1314-62-1$ | vanadium pentoxide | no | no |
| $68990-52-3$ | vegetable oil fatty acid methyl ester | no | no |
| $108-05-4$ | vinyl acetate | yes | no |
| $593-60-2$ | vinyl bromide | yes | no |
| $75-01-4$ | vinyl chloride | yes | yes |
| $5906-75-2$ | vinyl dimethylsilanol | no | no |
| $25013-15-4$ | vinyl toluene | no | no |
| $30030-25-2$ | vinylbenzylchloride | no | no |
| $75-35-4$ | vinylidene chloride (1,1-dichloroethylene) | no | no |
| $75-38-7$ | vinylidene fluoride | no | no |
| $5507-44-8$ | vinylmethyldiethoxysilane | no | no |
| $75-94-5$ | vinyltrichlorosilane | no | no |
| $2768-02-7$ | vinyltrimethoxysilane | no | no |
| $8032-32-4$ | vm \& p naphtha | no | no |
| $8042-47-5$ | white mineral oil | no | no |
| --0 | witconol al $69-66$ | no | no |
| $8002-09-3$ | yarmor pine oil | no | no |
| $1314-13-2$ | zinc oxide | no | no |
| $557-05-1$ | zinc stearate |  |  |

## ApPENDIX L:

## Benchmarking Survey of State Air Toxics Assessments in New Source Permitting

# Benchmarking Survey of State Air Toxics Assessments in New Source Permitting 

Robert Sills, Supervisor, Toxics Unit, Air Quality Division, Michigan Department of Natural Resources and Environment

February 25, 2010

## Background and Introduction

The Michigan Department of Natural Resources and Environment (MDNRE) Air Quality Division (AQD) implements the "air toxics rules" (Rules 224-232) of Part 55 of the Natural Resources and Environmental Protection Act (NREPA) as part of the New Source Review (NSR) permitting program. Because the federal government has not required air toxics risk assessment in NSR, except for the limited and long-delayed requirements of the Clean Air Act under Section 112(f), many states have developed their own requirements to better ensure public health protection. Recently AQD has become aware of interest regarding the scope and basis for the MDNRE air toxics regulatory requirements, and how they compare to other state's programs. In particular, there is interest in comparing the issue of "the list", i.e., the scope of the air toxics included in the state's programs.
Previous "benchmarking" surveys have been conducted, however, they do not provide sufficient detail on this particular issue. For example, previous surveys by MDEQ (2009) and the Louisville (2005) local air pollution control agency are helpful for many purposes, but do not provide sufficient and current program details regarding the key question about "the list" which is the present interest. And, given the broad variety of state air toxics programs, and the many nuances in their scope and applicability, some surveys only provide a simple "yes" or "no" indication of the requirement for air toxics risk assessment.

Proper framing of the survey questions is critical to obtaining the desired information. The present survey sought to find if state air permitting programs go beyond the federal technologybased requirements and address public health concerns for ambient air impacts of air toxics emissions. Care was taken to avoid "false-negative" responses. For example, "false negative" responses could result if a question is phrased, "Is air toxics risk assessment required as part of New Source Review?" In response to that question, a state representative may unfortunately reply "no", if only because, 1) they evaluate modeled ambient air impacts in comparison to some health-based criteria such as TLV/100, but they consider that "screening" rather than "risk assessment"; 2) they have established permissible emission rate limits, which were derived based on assumed facility parameters (e.g., building and stack height and distance to fenceline), dispersion modeling, and health-based ambient air exposure criteria, which they may not think of as being essentially "risk assessment based"; or, 3) they don't perform the assessment as a requirement of their rules, but as a matter of policy. With regard to this 3rd point, the present survey found that there are many states which do not have air toxics risk assessment-based requirements in state statutes or rules per se, however, they do conduct air toxics impact and risk assessment as a policy under broad "safety net" language in statute or rule. The "safety net" language cited by many states generally requires that air emissions shall not pose a threat to the public health (similar to Michigan’s Rule 901 under NREPA Part 55).

Some states have air toxics impact assessment requirements which are fairly unusual or unique. For example, some state programs specifically evaluate (or exclude from evaluation) selected
source categories, or, they utilize air toxics monitoring data for targeted geographic areas to drive initiatives to reduce emissions of selected air toxics. The present benchmarking survey attempted to note some of these significant program nuances, while primarily attempting to clarify if the air toxics addressed were limited to a specific list or not. As indicated in the "reference/contact" column of the table below, the results of the previous surveys by MDEQ (2009) and Louisville (2005) were relied upon in many cases, while in many other cases an appropriate state contact person was interviewed. It should also be noted that many state air permitting programs, like Michigan's, have a number of permit exemptions, permits by rule, or allowable emission thresholds, which would circumvent the need to perform modeling of ambient air impacts for air toxics to determine acceptability. Those program nuances have not been compiled in the present exercise, but are a significant and relevant aspect of state program comparisons nevertheless.

## Results

| State | Reference / contact | For proposed new/modified air emission sources, are ambient air impacts of any air toxics evaluated? If yes, what is the regulatory basis? | What air toxics are included? | What are the ambient air impacts compared to in order to determine acceptability? |
| :---: | :---: | :---: | :---: | :---: |
| Alabama | Wes Thornhill 334-271-7887 | Yes, by policy but not in rules. | All air toxics with TLVs or other OELs. | If the substance has an OEL AND is emitted at >0.1 $\mathrm{lb} / \mathrm{hr}$, then the modeled ambient air impact cannot exceed TLV/40 (8 hr AT) or TLV/420 (annual AT). |
| Alaska | $\begin{aligned} & \text { MDEQ (2009); } \\ & \text { Louisville } \\ & \text { (2005) } \end{aligned}$ | No. |  |  |
| Arizona | MDEQ (2009); <br> Louisville <br> (2005) | No. |  |  |
| Arkansas | MDEQ (2009); Louisville (2005) | No. |  |  |
| California | $\begin{aligned} & \text { Louisville } \\ & \text { (2005) } \end{aligned}$ | Yes, by Hot Spots regulation; sources causing fenceline or community monitored levels of excess risk addressed via control measures (existing; point, area or mobile); modeling done for new sources. | 748 total air toxics; 438 must be quantified in risk assessment (as of 2005 survey) | CA-OEHHA Reference Exposure Levels (RELs), or, one in 1 million cancer risk. |
| Colorado | $\begin{aligned} & \hline \text { MDEQ (2009); } \\ & \text { Louisville } \\ & \text { (2005) } \end{aligned}$ | No. |  |  |
| Connecticut | Jim Grillo 860- <br> 424-4152; <br> Louisville <br> (2005) survey. | Yes. In rules. New and existing sources; major and area sources. | The HAPs list (187). Hazardous Limiting Values (HLVs) were derived for the HAPs based on modified occupational standards. | The rules provide 2 equations (one for under 20 m stacks, one for over 20 m stacks) relating air emissions to ambient impacts, which are compared to HLVs; it is a pass/fail standard for all permits. |


| Delaware | $\begin{aligned} & \text { Jim Snead 302- } \\ & 323-4542 \end{aligned}$ | Yes, by policy but not in rules. Policy is under a general "safety net" provision (regulation 1102). | All substances; no discrete list. | Maximum ambient air impacts cannot exceed TLV/100 if there is a TLV available; if not, then impact cannot exceed the default value of $100 \mu \mathrm{~g} / \mathrm{m}^{3}$. This is the same approach for carcinogens as well as noncarcinogens. |
| :---: | :---: | :---: | :---: | :---: |
| Florida | MDEQ (2009); Louisville (2005) | No. |  |  |
| Georgia | Eric Cornwell 404-363-7020 | Yes, in guidance only; not by rule; under "safety net" rule provisions. | No discrete list; any substance with IRIS value or OEL. | Hierarchy used; 1) most stringent value between cancer-based value (one in 1 million if " A " carcinogen, otherwise, 1 in 100,000 ) or RfC; 2) TLV/100 (or, TLV/300 if "A" carcinogen), then scaled by $40 \mathrm{hrs} / 168 \mathrm{hrs}$ (approx. a factor of 4) to derive acceptable ambient concentration (AAC) with 24 hr AT; for OELs which are ceiling limits or STELs, divide by 10 and also scale by a factor of 1.32 to account for 15" AT of OEL (per SCREEN3). |
| Hawaii | MDEQ (2009); Louisville (2005) | Yes; new/modified sources only; major and area sources. | HAPs only. |  |
| Idaho | $\begin{aligned} & \text { Carl Brown } \\ & \text { 208-373-0206 } \end{aligned}$ | Yes. In rules. New/modified sources only. Does not apply if a MACT rule applies. | Approximately 350 toxic air pollutants; list was developed before the 1990 HAPs list | Utilize conservative pph emission thresholds; if exceeded, then ambient air impacts modeled; acceptable ambient concentrations (AACs) are based on 1E-06 cancer risk, and for noncarcinogens, OEL/UF. |
| Illinois | Jeff Sprague 217-524-4692 | No, unless there are public concerns. Do have an internal screening for ethanol plants. |  |  |
| Indiana | Bryan Wolff 317-234-3499 | Yes. By policy; air toxics impacts are assessed only if requested by citizen or applicant. No routine screening. | No discrete list; any substance with any state or federal criteria or any health data may be included. | Commission has discretionary basis for permit denial if impacts are deemed adverse. |


| Iowa | MDEQ (2009); Louisville (2005) | No. |  |  |
| :---: | :---: | :---: | :---: | :---: |
| Kansas | MDEQ (2009); Louisville (2005) | No. |  |  |
| Kentucky | $\begin{aligned} & \text { Taimur Shaikh } \\ & \text { 502-564-3999 } \\ & \text { x4480 } \end{aligned}$ | Yes, as a policy regarding new/modified source permitting, under a general "safety net" regulation regarding public health protection. | EPA HAPs plus all substances regulated by EPA under the chemical accident prevention provisions (CAAA Section 112(r)). | Risk assessment based levels associated with $\mathrm{HQ}=1$ or one in 1 million incremental cancer risk. |
| Louisiana | Louisville (2005) | Yes. | HAPs plus other air toxics. | Ambient impacts cannot exceed TLV/factor, or one in 10,000 cancer risk. |
| Maine | Lisa Higgins 207-287-7023; Louisville (2005) survey | Yes. | Have ambient air quality guidelines for HAPs plus additional compounds. | Have calculated healthbased guideline values. Have a State statute mercury emission limit of $25 \mathrm{lbs} / \mathrm{yr}$ for any new or existing facility. |
| Maryland | Louisville (2005) | Yes. | All HAPs plus others; database of 6329 substances as of 2005 survey. | Maximum ambient air impacts cannot exceed TLV/100 or one in 1 million cancer risk. |
| Massachusetts | Marc Wolman 617-292-5515 | Yes, as ambient air guidelines. Apply to only: incinerators, WWTPs and residuals mgmt., major remedial actions, and PSD projects. | Discrete list of air toxics ( $\mathrm{n} \sim 120$ ) which pre-dates the EPA 1990 HAPs list | They have derived threshold effects exposure limits (TELs; 24 hr AT) and allowable ambient limits (AALs; annual AT) for all the targeted air toxics. |
| Michigan | Robert Sills 517-335-6973 | Yes. Required by air toxics rules. New / modified sources only. | There is an open-ended definition of Toxic Air Contaminants (TACs); includes all substances other than 41 listed nonTACs. Health-based screening levels have been developed for approx. 1200 TACs. | Screening levels (SLs) for carcinogens are at 1E-06 risk per chemical for the proposed process; or, 1E-05 is acceptable for facilitywide emissions per chemical. Noncancer SLs are derived from RfCs, RfDs, OELs, or other data; default $=0.1 \mu \mathrm{~g} / \mathrm{m}^{3}$. SLs on website. |

[^1]| Minnesota | Mary Dymond 651-757-2327 | Yes. By policy, an Air Emissions Risk Analysis (AERA) is needed for proposed new/modified sources exceeding emission thresholds, or if "flexible air permit", or if needed per MPCAs discretion; existing sources may also need an AERA if significant public interest. | All substances which have a health benchmark value from MN Dept of Health, EPA-IRIS, or California-OEHHA. | Facility-wide emissions, multi-media impacts: risk guidelines are for a cancer risk of 1E-05 and cumulative hazard index of 1 for pollutants with the same toxic endpoint. |
| :---: | :---: | :---: | :---: | :---: |
| Mississippi | $\begin{aligned} & \text { Danny Jackson } \\ & \text { 601-961-5225 } \end{aligned}$ | No; risk provisions are only implemented as needed, and are not being triggered by anything at present. |  |  |
| Missouri | MDEQ (2009) | No. |  |  |
| Montana | MDEQ (2009); Louisville (2005) | No, except incinerators must demonstrate negligible risk. |  |  |
| Nebraska | MDEQ (2009); Louisville (2005) | No. |  |  |
| Nevada | MDEQ (2009); Louisville (2005) | No. |  |  |
| New Hampshire | $\begin{aligned} & \text { Pat North 603- } \\ & \text { 271-0901 } \end{aligned}$ | Yes; by rule; new and existing sources of all types. | Utilize a discrete list of $\sim 800$ air toxics, including all HAPs plus substances with ACGIH TLVs or IRIS values. | OELs are divided by UFs depending on the OEL type. Three cancer classifications are recognized. |
| New Jersey | $\begin{aligned} & \text { Olga Boyko } \\ & \text { 609-633-1108 } \end{aligned}$ | Yes; by regulations. | Regulations reference the HAPs list, and also an older pre-HAPs list of air toxics. Risk screening is done for ALL compounds with health benchmarks from EPA, CA, etc. | They utilize permit reporting thresholds which trigger a reporting requirement; utilize $\mathrm{HI}=1$, and one in 1 million cancer risk for a process (one in 100,000 for facility-wide emissions). |
| New Mexico | Ted Schooley 505-476-4334; Louisville (2005) | Yes. New/modified sources only. | HAPs plus substances with OELs. | Use chemical-specific pph emission thresholds; if exceeded, then modeled ambient air impacts cannot exceed OEL/100 or MDL if carcinogenic. |


| New York | $\begin{aligned} & \hline \text { Tom Gentile } \\ & 518-402-8402 \end{aligned}$ | Yes. Required in rules. New and existing sources, excluding fossil fuel combustion sources (which are regulated separately). | Regulated air pollutants (RAPs) defined as criteria pollutants, HAPs, and CAA 112(r) compounds. | Guideline values derived via risk assessment. Currently considering draft rulemaking to restrict RAPs to a shorter list of high priority cpds., due to limited r.a. staffing. |
| :---: | :---: | :---: | :---: | :---: |
| North Carolina | MDEQ (2009); Louisville (2005) | Yes. | HAPs plus a discrete list of other air toxics. | Acceptable ambient pollutant levels established. |
| North Dakota | MDEQ (2009); Louisville (2005) | Yes; new/modified major and area sources. | 700 air toxics, including HAPs, as of 2005 survey. | TLV/100 or one in 1 million cancer risk cannot be exceeded in ambient air. |
| Ohio | Paul Koval 614-644-2270 | Yes. Per rules. For new or existing sources with over 1 ton/yr emissions of TAPs. | Toxic Air Pollutants $($ TAPs $)=303$ substances . | TLV/42 for noncarcinogens. |
| Oklahoma | MDEQ (2009); Louisville (2005) | Yes. | 1500 air toxics as of 2005 survey. | TLV divided by a factor which depends on the degree of toxicity. |
| Oregon | Patricia Huback 503-229-6932 | No. Development of a program is under consideration. | Have 3 strategies in place to address air toxics concerns: 1) geographic approach based on NATA to identify areas of concern and develop strategies to reduce risks; 2) statewide source sector strategy approach (e.g., wood stoves); 3) safety net program, to address concerns identified by fenceline monitoring or source modeling. | Their Air Toxics Advisory Committee has established public health protective levels ("ambient benchmark concentrations") for 51 air toxics. Diesel, benzene, manganese, formaldehyde, steel foundry emissions, and wood stoves are among the higher priorities. |
| Pennsylvania | Dean Van Orden 717-787-1455 | No, not routinely or as a broad policy. State statute does have a "safety net" provision, and under that, permit engineers have discretion to evaluate air toxics impacts and risks. Landfill gases, combustors, and cement kiln emissions have been evaluated. | HAPs plus other air toxics of concern (source-specific). |  |
| Rhode Island | MDEQ (2009); <br> Louisville (2005) | Yes. | HAPs plus a discrete list of other air toxics. | RfCs and other noncancer benchmarks; one in 1 million to one in 100,000 cancer risk. |


| South Carolina | $\begin{aligned} & \text { Louisville } \\ & \text { (2005) } \end{aligned}$ | Yes; new/modified and existing. | 257 toxic air pollutants (TAPs), as of 2005 survey. |  |
| :---: | :---: | :---: | :---: | :---: |
| South Dakota | MDEQ (2009); Louisville (2005) | No. |  |  |
| Tennessee | MDEQ (2009); Louisville (2005) | No, except in a few cases where public interest is high. |  |  |
| Texas | Manuel Reina 512-239-1816 | Yes. "Safety-net" rule for the protection of the public; policy under that for the modeling and assessment procedure. New / modified sources only. | All substances are subject; list of substances identified in air emissions with Effect Screening Levels (ESLs) developed has grown since 1980's to over 3000 substances. | Target cancer risk $=1 \mathrm{E}-05$ per substance, facilitywide emissions. For noncarcinogens, TLV/100 (1 hr AT) and TLV/1000 (annual AT); default=1 $\mu \mathrm{g} / \mathrm{m}^{3}$. Draft ESLs and justifications public noticed. All appear on website. |
| Utah |  | No. |  |  |
| Vermont | MDEQ (2009); Louisville (2005) | Yes; new/modified and existing sources; major and area sources. | 382 hazardous air pollutants, all HAPs, plus any new air toxic if toxicological information is available. | TLV divided by UF; one in 1 million incremental cancer risk. |
| Virginia | Patricia Buonviri 804-698-4016 | Yes, unless source is covered by a MACT standard; requirement is in regulations. | HAPs list with a couple of exceptions. | TLV divided by UF. No cancer risk-based criteria. Currently considering rule revisions to adopt a more risk-based program. |
| Washington | MDEQ (2009); Louisville (2005) | No. |  |  |
| West Virginia | MDEQ (2009); Louisville (2005) | Yes. | HAPs plus substances with OELs. |  |
| Wisconsin | $\begin{aligned} & \text { Jeff Myers 608- } \\ & 266-2879 \end{aligned}$ | Yes. By rule; applies to new and existing sources, except for HAPs covered by a MACT std., or if chemical-specific health-based emission thresholds are not exceeded. |  | Noncarcinogens: use RfCs or TLV/42 as ambient standards not to be exceeded by aggregate impacts of the source, bkgd. levels, and impacts from other sources. <br> Carcinogens: technologybased control only <br> (LAER), or, can use lowrisk modeling demonstration (1E-06 per cpd., or 1E-05 facilitywide) as a compliance option. |
| Wyoming | MDEQ (2009); Louisville (2005) | No. |  |  |

## Discussion

Thirty states evaluate and regulate air toxics emissions in their permit reviews, based on public health exposure concerns, although there are many state-specific nuances regarding the regulatory basis, the types of sources included, the air toxics included, the acceptability criteria, and exemptions. Of the six states in EPA Region 5, five states generally and routinely evaluate air toxics ambient air impacts for public health acceptability; only Illinois generally does not (but may in exceptional cases). Of the eight Great Lakes states, six states generally and routinely evaluate air toxics ambient air impacts for public health acceptability; only Illinois and Pennsylvania generally do not (but they may in exceptional cases).

## Acronyms and abbreviations not defined in text or table:

1E-05= one in 100,000 incremental cancer risk
1E-06= one in 1 million incremental cancer risk
$\mathrm{AT}=$ averaging time
bkgd.= background
CAA= clean air act
cpd.= compound
HAPs= hazardous air pollutants
HI= hazard index
$\mathrm{HQ}=$ hazard quotient
LAER= lowest achievable emission rate
MDL= method detection limit
NATA = National-Scale Air Toxics Assessment
OEL= occupational exposure level
pph= pounds per hour
RfC= reference concentration
RfD $=$ reference dose
TLV= threshold limit value
UF= uncertainty factor
$\mu \mathrm{g} / \mathrm{m}^{3}=$ micrograms per cubic meter

## References

Louisville Air Pollution Control. 2005. Summary of State Air Toxics Programs. Unpublished report.

Michigan Department of Environmental Quality (MDEQ; currently MDNRE). 2009. Survey of State Air Permitting Programs. By Doreen Lehner, MDNRE-AQD. Unpublished report.

## Appendix M:

## EPA Region 5 States Benchmarking Comparison Table

## January 17, 2013 R. Sills EPA Region 5 States Benchmarking Comparison Table

HRA = Health risk assessment; i.e., modeling of ambient air impacts and comparison to health-protective benchmark values $N / M=$ New or modified sources. E = Existing sources.

| Air Toxics Program Characteristic |  | MI | MN | OH | WI | IN | IL |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Impetus for HRA of air emission sources | Required by statute or rules | yes | yes (statute requires cumulative RA for certain sources and locations) | yes | yes |  |  |
|  | By policy |  | yes (except as noted above) |  |  | yes | yes |
|  | If significant interest by public or applicant (i.e., not routine) |  | yes (for existing sources) |  |  | yes | Not routine, done only if significant public concerns. |
|  | Discretionary by agency |  |  |  |  | No criteria for max. ambient air impacts, but permit may be denied if "adverse." |  |
| HRA done for new/modified (N/M) or existing (E) sources |  | N/M | N/M, and also E if significant public interest | N/M. Also existing sources are evaluated on a case-bycase basis | N/M or E | N/M | N/M |


| Source types or emission rate exemptions from HRA? |  | Yes | Yes | Yes (i.e., exempt if each TAP emission is $\leq 1$ ton/yr) | Yes (i.e., HAPs exempt if covered by a MACT, but only if chem-specific emission thresholds not exceeded) |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| What air toxics are subject to HRA? | Any |  |  |  |  | yes |  |
|  | All except 41 exemptions | yes |  |  |  |  |  |
|  | Unique list beyond HAPs (how many CPDs/Groups) |  | yes (any with a benchmark value from IRIS, Cal or MDH) | yes ( $\mathrm{n}=303$ <br> TAPs; includes all HAPs plus others) | yes ( $\mathrm{n}=535$; 26 <br> HAPs not included) |  |  |
|  | HAPs only |  |  |  |  |  |  |
| How are cumulative air toxics impacts accounted for? | Generally not accounted for in permit review. | yes |  | yes | yes | yes |  |
|  | Can be considered in permit review | yes (Rule 228) | Done under statutory requirements for Minneapolis. | yes (combined impacts; not background conc.) |  |  |  |
|  | Routinely accounted for in permit review. |  |  |  |  |  |  |


| How are cumulative air toxics impacts accounted for? (continued) | ? Routinely evaluated via statewide monitoring or modeling initiative, $\pm$ risk reduction targets? | Detroit ambient air evaluated in 2005 and 2010 Detroit Air Toxics Initiative reports; no risk $\downarrow$ target; several facilityspecific monitors are in operation. |  | Specific <br> monitoring or <br> modeling <br> studies have <br> been <br> conducted to <br> evaluate <br> specific <br> concern <br> sources/areas. <br> No risk $\downarrow$ <br> target. | RAIMI <br> statewide HAPs modeling of cumulative impacts of all sources; goal of 50\% reduction (from 2002 to 2012) of people at >1E-6 CA risk. | Focused studies of monitoring and risk assessment completed for Indianapolis, and underway for the lakeshore area; statewide RAIMI modeling; are evaluating high-risk NATA'05 facilities. No risk $\downarrow$ target. |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| Acceptable risk | chmarks | 1E-6 per cpd for the process. 1E-5 per cpd for the facility. 10X higher for roads and indus. areas. EPA or other agency values; TLV/100; or derived from short-term studies. Default ITSL= 0.1 $\mu \mathrm{g} / \mathrm{m}^{3}$. | Provided to MPCA by MDH, based on values from EPA or other agencies, or derived by MDH. | IRIS values; modeled 1 hr AT max impacts not to exceed TLV/42. | 1E-6 per cpd. $1 \mathrm{E}-5$ all cpds. EPA RfCs etc. TLV-TWA/42 with $24-\mathrm{hr}$ AT. TLV-Ceiling/10 with 1-hr. AT. | Use various EPA approved sources and databases. |  |

## Appendix N:

## Consistency with Other States

# Air Toxics Workgroup "Consistency with Other States" Discussion Paper 

April 10, 2013

## ORR (2011) Report Recommendation A-1(7):

R 336.1225 should be amended and specifically include the following:
Make the acceptable exposure limits consistent with other nearby states.

## ATW discussion

Discussion with the ATW indicates that some members have concerns for a lack of consistency between MDEQ and the nearby states with regard to the air toxics screening level values and/or averaging times, which can contribute to an un-level playing field. AQD staff committed to developing some information and comparisons to help inform the discussion.

## AQD impressions

There are differences between states' air toxics health-based screening levels for several possible reasons, which may be summarized as follows:

1. States may use different target risk level for carcinogens (e.g., 1E-5 vs. 1E-6). WDNR applies a 1E-6 target risk per chemical (and 1E-5 for cumulative risk), while MPCA and MDEQ allow 1E-5 per chemical (see Table 3 cancer risk values and risk levels).
2. States may use different methods for deriving a benchmark. For example, OEL/100 vs. OEL/42. States have different methods to address (or not address) data-poor situations; see the discussion below.
3. States may adopt their screening levels from benchmarks provided by other recognized sources. Many substances have multiple applicable benchmarks already available from recognized sources, such as EPA-IRIS values, EPA-PPRTVs (from the Superfund program office), ATSDR MRLs, and CaIOEHHA RELs, and Texas TCEQ ESLs. The benchmarks available from these sources are often different. A state may review all of those available, or utilize a hierarchy, and choose to adopt any one of these available benchmarks as-is or with modification. States may vary in their choices. Also, states establish their screening levels at different points in time, when different key studies and different benchmarks may be available. Many of DEQ's screening levels were developed in the 1990s and 2000s. See Table 1 for general hierarchies utilized by States for establishing chronic inhalation screening levels. See the Table 3 manganese example; the DEQ ITSL is based on the EPA-IRIS RfC (1993), while the MPCA screening level was derived in the 2000s by MDH.
4. Different critical effects may be addressed by the different state's benchmarks. For example, see the styrene example in Table 3: DEQ regulates it as a carcinogen, while MPCA and WDNR do not.
5. States may establish acute screening levels in addition to chronic noncancer screening levels. These can be derived by the agency or adopted from a recognized agency source; as with \#3 above, such values may differ. There are some widely accepted sources of acute benchmarks: acute inhalation Minimum Risk Levels (MRLs) from the ATSDR; Acute Exposure Guidance Levels (AEGLs) from EPA's National Advisory Committee; and, California OEHHA's Acute Reference Exposure Levels (ARELs). Texas TCEQ also derives acute ESLs. Occupational Exposure Levels
(TLVs, Ceiling Limits, Short-term Exposure Limits) are also used by MDEQ and other agencies to derive acute benchmarks, with the application of an uncertainty factor to help ensure protection of sensitive individuals.
6. States may have different conventions for setting averaging times for their screening levels.

Table 3 has examples of different states having the same screening level value, but different ATs for this reason.

## Data-Poor Situations

One of the most significant programmatic differences between DEQ and the other R5 State agencies is in the treatment of data-poor situations for noncancer risk assessment. Based on the recommendations from the 1981, 1989, and 1997 stakeholder workgroup reports, MDEQ has adopted rules and algorithms for utilizing short-term study results (short-term NOAELs and LOAELs; LC50s and LD50s) to derive ITSLs (with annual ATs) that are presumptively protective from chronic noncancer exposure and adverse effects, when the preferred studies or ITSL bases are not available (Rule 232). Ohio, Wisconsin and Minnesota would not extrapolate to derive chronic benchmarks, although they may address such limited datasets by setting acute screening levels. Texas TCEQ is an example of another state agency that utilizes LC50 data to derive acute and chronic benchmarks; their acute benchmark method is more restrictive than the DEQ approach.

Table 1. General Hierarchy of Basis for Chronic Inhalation Health Benchmarks

| Hierarchy I rank ${ }^{1}$ | Michigan DEQ | Minnesota PCA | Ohio EPA | Wisconsin DNR |
| :---: | :---: | :---: | :---: | :---: |
| Relatively higher | IRIS RfC value. Rules have default AT of 24 hours, which can be overridden by staff for an annual AT. | MDH healthbased value (hbv) | IRIS or other available appropriate benchmark from reputable agency. 1 hr AT. | EPA values and ACGIH TLVs. |
| $\downarrow$ | EPA RfD, ATSDR MRL, EPA PPRTV, Cal REL, or staff-derived $\mathrm{RfC}^{2}$. AT may be 24 hours (default in rules for RfD). | MDH health risk value (hrv) |  |  |
| $\downarrow$ | OEL (TLV/100). AT is 8 hrs. |  |  |  |
| $\downarrow$ | Subchronic study (e.g., 2week) with extrapolation to chronic. Annual AT | Cal REL, EPA HEAST, ATSDR MRL | Compare to other chemicals with similar structures, apply SAR. |  |
| $\downarrow$ | $\mathrm{LC}_{50}$ value with extrapolation to chronic. Annual AT. | EPA Superfund PPRTV |  |  |
| $\downarrow$ | $L_{50}$ value with extrapolation to chronic. Annual AT. |  |  |  |
| Relatively lower | Default ITSL $=0.1 \mu \mathrm{~g} / \mathrm{m}^{3}$ (annual AT). | No default | No default | No default |
| Comments | Methods for deriving ITSLs from very limited data are protective, and have a long history at AQD. | Rarely use OELs (exception: ethanol facilities). Do not use short-term bioassay data to derive screening values. <br> Chemicals with inadequate data are evaluated qualitatively in context with the entire facility. | Chemicals with inadequate data may be evaluated by comparison to similar compounds with better tox data (computational toxicology). | They do not address air toxics without benchmarks available from other reputable sources. |

[^2]${ }^{2}$ Depending on the age and basis for the available benchmarks from other reputable agencies, AQD toxicologist staff may perform an updated literature review and utilize key studies differently than other available benchmarks in deriving an ITSL utilizing EPA's RfC methodology.

Table 2. Access to R5 State's Air Toxics Information and Screening Levels

| State Agency | Location |
| :--- | :--- |
| Michigan DEQ | $\underline{\text { http://www.michigan.gov/deq/0,4561,7-135-3310 4105---,00.html }}$ |
| Ohio EPA | $\frac{\text { http://epa.ohio.gov/dapc/regs/3745 114.aspx (Toxics compound }}{\text { data sheets ONLY; NOT a list of benchmarks.) }}$ |

Table 3. Comparison of R5 States' Health-Based Screening Levels for Select Air Toxics (acute and chronic noncancer; cancer at specified risk level, with annual AT; all values in $\mu \mathrm{g} / \mathrm{m}^{3}$ ).

| Chemical | MDEQ-AQD | MPCA | Ohio EPA ${ }^{1}$ | WDNR |
| :---: | :---: | :---: | :---: | :---: |
| Acetaldehyde \#75-07-0 | $\begin{aligned} & 9 \text { (24 hr AT) } \\ & 5 \text { (1E-5 cancer) } \end{aligned}$ | $\begin{array}{\|l\|} \hline 470(1 \mathrm{hr} \text { AT }) \\ 9 \text { (chronic) } \\ 4.5 \text { (1E-5 cancer) } \\ \hline \end{array}$ |  | $\begin{aligned} & 4504(1 \mathrm{hr} \text { AT }) \\ & 0.45(1 \mathrm{E}-6 \\ & \text { cancer) } \\ & \hline \end{aligned}$ |
| Acrolein \#107- $02-8$ | $\begin{aligned} & \hline 5 \text { ( } 1 \mathrm{hr} \text { AT) } \\ & 0.02 \text { (annual } \\ & \text { AT) } \\ & \hline \end{aligned}$ | $\begin{aligned} & 5 \text { (1 hr AT) } \\ & 0.4 \text { (chronic) } \end{aligned}$ |  | 22.9 (1 hr AT) |
| Ammonia \#7664-41-7 | 100 (24 hr AT) | $\begin{aligned} & 3200 \text { (1 hr AT) } \\ & 80 \text { (chronic) } \end{aligned}$ |  | $\begin{aligned} & 418 \text { (24 hr AT) } \\ & 100 \text { (annual } \\ & \text { AT) } \\ & \hline \end{aligned}$ |
| $\begin{aligned} & \text { Benzene \#71- } \\ & \text { 43-2 } \end{aligned}$ | $\begin{aligned} & \hline 30 \text { (24 hr AT) } \\ & 30 \text { (annual AT) } \\ & 1 \text { (1E-5 cancer) } \\ & \hline \end{aligned}$ | $\begin{aligned} & \hline 1000(1 \mathrm{hr} \mathrm{AT}) \\ & 30 \text { (chronic) } \\ & 1.3 \text { (1E-5 cancer) } \end{aligned}$ |  | $\begin{aligned} & 0.13(1 \mathrm{E}-6 \\ & \text { cancer) } \end{aligned}$ |
| Benzo(a)pyrene \#50-32-8 | $\begin{aligned} & \text { 5E-3 (1E-5 } \\ & \text { cancer) } \\ & \hline \end{aligned}$ | $\begin{aligned} & 9.1 \mathrm{E}-3(1 \mathrm{E}-5 \\ & \text { cancer) } \end{aligned}$ |  | $\begin{array}{\|l} \hline 9.1 \mathrm{E}-4(1 \mathrm{E}-6 \\ \text { cancer) } \end{array}$ |
| $\begin{aligned} & \text { Cadmium \#7440- } \\ & 43-9 \end{aligned}$ | $\begin{aligned} & 6 \mathrm{E}-3(1 \mathrm{E}-5 \\ & \text { cancer) } \end{aligned}$ | 0.02 (chronic) 5.6E-3 (1E-5 cancer) |  | $\begin{aligned} & 5.6 \mathrm{E}-4(1 \mathrm{E}-6 \\ & \text { cancer) } \end{aligned}$ |
| Chlorine \#7782-50-5 | $\begin{aligned} & \hline 500 \text { (8 hr AT) } \\ & 0.3 \text { (annual AT) } \\ & \hline \end{aligned}$ | $\begin{array}{\|l\|} \hline 290 \text { ( } 1 \mathrm{hr} \mathrm{AT} \text { ) } \\ 0.2 \text { (chronic) } \\ \hline \end{array}$ |  | 34.8 (24 hr AT) |
| Diethylene glycol monobutyl ether | 20 (24 hr AT) | 0.1 (chronic) |  | $\begin{aligned} & 2320 \text { (24 hr AT) } \\ & 13000 \text { (annual } \\ & \hline \end{aligned}$ |


| (butyl cellosolve) \#112-34-5 |  |  |  | AT) |
| :---: | :---: | :---: | :---: | :---: |
| Epichlorohydrin \#106-89-8 | $\begin{aligned} & 1 \text { ( } 24 \mathrm{hr} \text { AT) } \\ & 8 \text { (1E-5 cancer) } \end{aligned}$ | $\begin{aligned} & 1300(1 \mathrm{hr} \mathrm{AT}) \\ & 1 \text { (chronic) } \\ & 8.3 \text { (1E-5 cancer) } \end{aligned}$ |  | $\begin{array}{\|l} \hline 45.4(24 \mathrm{hr} \text { AT }) \\ 0.83(1 \mathrm{E}-6 \\ \text { cancer) } \\ \hline \end{array}$ |
| Ethylene glycol \#107-21-1 | 1000 (1 hr AT) | 400 (chronic) |  | N/A |
| Ethylene oxide \#75-21-8 | $\begin{aligned} & 0.3 \text { (1E-5 } \\ & \text { cancer) } \end{aligned}$ | $\begin{aligned} & \hline 30 \text { (chronic) } \\ & 0.11 \text { (1E-5 } \\ & \text { cancer) } \\ & \hline \end{aligned}$ |  | $\begin{aligned} & \text { 1.1E-2 (1E-6 } \\ & \text { cancer) } \end{aligned}$ |
| Table 3, continued... |  |  |  |  |
| Chemical | MDEQ-AQD | MPCA | Ohio EPA ${ }^{1}$ | WDNR |
| Formaldehyde \#50-00-0 | $\begin{aligned} & 9(8 \mathrm{hr} \mathrm{AT}) \\ & 0.8(1 \mathrm{E}-5 \\ & \text { cancer }) \end{aligned}$ | $\begin{aligned} & 94 \text { (1 hr AT) } \\ & 9 \text { (chronic) } \\ & 2 \text { (1E-5 cancer) } \end{aligned}$ |  | 7.7E-2 (1E-6 cancer) |
| Hexane \#110-54-3 | 700 (24 hr AT) | 2000 (chronic) |  | $\begin{aligned} & 4320(24 \mathrm{hr} \text { AT }) \\ & 200 \text { (annual } \\ & \text { AT) } \\ & \hline \end{aligned}$ |
| Hydrogen chloride \#7647-01-0 | $\begin{aligned} & 2100(1 \mathrm{hr} \text { AT) } \\ & 20 \text { (annual AT) } \end{aligned}$ | $\begin{aligned} & 2700(1 \mathrm{hr} \text { AT) } \\ & 20 \text { (chronic) } \end{aligned}$ |  | $\begin{array}{\|l\|} \hline 746(1 \mathrm{hr} \mathrm{AT}) \\ 20 \text { (annual AT) } \end{array}$ |
| Hydrogen sulfide \# 7783-06-4 | $\begin{aligned} & 100 \text { (24 hr AT) } \\ & 2 \text { (annual AT) } \\ & \hline \end{aligned}$ | $\begin{aligned} & \hline 42 \text { ( } 1 \mathrm{hr} \text { AT) } \\ & 2 \text { (chronic) } \\ & \hline \end{aligned}$ |  | 335 (24 hr AT) |
| Manganese | $\begin{aligned} & 0.05 \text { (annual } \\ & \text { AT) } \end{aligned}$ | 0.2 (chronic) |  | 4.8 (24 hr AT) |
| Mercury \#7439-97-6 | $\begin{aligned} & \text { (no ITSL; } \\ & \text { inhalation-only } \\ & \text { RfC }=0.3 \\ & \mu \mathrm{~g} / \mathrm{m}^{3} \text { ) } \end{aligned}$ | $\begin{aligned} & \hline 0.6 \text { (1 hr AT) } \\ & 0.3 \text { (chronic) } \end{aligned}$ |  | Inorganic: 0.6 ( 24 hr AT); 0.3 (annual AT). <br> Alkyl cpds: 0.24 (24 hr AT) |
| Methyl bromide \#74-83-9 | 5 (24 hr AT) | $\begin{aligned} & 2000(1 \mathrm{hr} \text { AT) } \\ & 5 \text { (chronic) } \\ & \hline \end{aligned}$ |  | 93.2 (24 hr AT) <br> 5 (annual AT) |
| Naphthalene \#91-20-3 | $\begin{aligned} & 3(24 \mathrm{hr} \mathrm{AT}) \\ & 0.8(1 \mathrm{E}-5 \\ & \text { cancer }) \end{aligned}$ | $\begin{aligned} & 200(1 \mathrm{hr} \mathrm{AT}) \\ & 9 \text { (chronic) } \\ & 0.29(1 \mathrm{E}-5 \\ & \text { cancer) } \\ & \hline \end{aligned}$ |  | 1258 (24 hr AT) |
| $\begin{aligned} & \text { Nickel \#7440-02- } \\ & 0 \end{aligned}$ | $\begin{aligned} & \text { 4.2E-2 (1E-5 } \\ & \text { cancer) } \end{aligned}$ | $\begin{aligned} & \hline 11 \text { (1 hr AT) } \\ & 0.014 \text { (chronic) } \\ & 2.1 \mathrm{E}-2 \text { (1E-5 } \\ & \text { cancer) } \end{aligned}$ |  | $\begin{aligned} & 3.8 \mathrm{E}-3(1 \mathrm{E}-6 \\ & \text { cancer) } \end{aligned}$ |
| ```Phenol #108-95- 2``` | 190 (8 hr AT) | $\begin{aligned} & 5800 \text { (1 hr AT) } \\ & 200 \text { (chronic) } \\ & \hline \end{aligned}$ |  | 462 (24 hr AT) |
| $\begin{aligned} & \text { Styrene \#100- } \\ & \text { 42-5 } \end{aligned}$ | $\begin{aligned} & 1000(24 \mathrm{hr} \mathrm{AT}) \\ & 17(1 \mathrm{E}-5 \\ & \text { cancer }) \\ & \hline \end{aligned}$ | $\begin{aligned} & 21000 \text { (1 hr AT) } \\ & 1000 \text { (chronic) } \end{aligned}$ |  | $\begin{aligned} & \hline 2045 \text { (24 hr AT) } \\ & 1000 \text { (annual } \\ & \text { AT) } \\ & \hline \end{aligned}$ |
| Toluene \#108- | 5000 (24 hr AT) | 37000 (1 hr AT) |  | 4522 (24 hr AT) |


| 88-3 |  | 400 (chronic) | $\begin{aligned} & 400 \text { (annual } \\ & \text { AT) } \end{aligned}$ |
| :---: | :---: | :---: | :---: |
| Trichloroethylene \#79-01-6 | $\begin{aligned} & 10000(24 \mathrm{hr} \\ & \text { AT) } \\ & 2 \text { (annual AT) } \\ & 2 \text { (1E-5 cancer) } \end{aligned}$ | $\begin{aligned} & 2000(1 \mathrm{hr} \mathrm{AT}) \\ & 2 \text { (chronic) } \\ & 3 \text { (1E-5 cancer) } \end{aligned}$ | $\begin{aligned} & 0.5(1 \mathrm{E}-6 \\ & \text { cancer) } \end{aligned}$ |
| Vinyl chloride \#75-01-4 | $\begin{aligned} & 100(24 \mathrm{hr} \mathrm{AT}) \\ & 1.1(1 \mathrm{E}-5 \\ & \text { cancer }) \end{aligned}$ | $\begin{aligned} & 180000(1 \mathrm{hr} \mathrm{AT}) \\ & 100 \text { (chronic) } \\ & 1.1 \text { (1E-5 cancer) } \end{aligned}$ | ```100 (annual AT) 0.11 (1E-6 cancer)``` |
| $\begin{array}{\|l} \hline \text { Xylenes \#1330- } \\ 20-7 \\ \hline \end{array}$ | 100 (24 hr AT) | 43000 (1 hr AT) 100 (chronic) | $10421 \text { (annual }$ AT) |

${ }^{1}$ Ohio EPA does not publish their air toxics benchmarks; no list is available. They have Toxic Compound Data Sheets available (see link in Table 3), however, these appear to be justifications for listing with a summary of known hazards and toxicity information sources (e.g., IRIS unit risk values and RfCs; ACGIH OELs). It is unclear how permit applicants and staff permit reviewers determine if modeled impacts are approvable.

## ApPENDIX O:

## TESTING ReqUIREMENTS IN PERMITS TO INSTALL

# Testing Requirements in Permits to Install 

April 16, 2013

## ORR Recommendation A-1(8)

The AQD should stop requiring permit holders to conduct elaborate and costly stack tests to provide emissions research data, since the DEQ does not use this information for subsequent permit reviews.

## Update

There are many reasons why stack testing requirements are included in permits to install. These include compliance demonstrations where it is required via regulations (i.e. NSPS, NESHAPs, etc.); there is uncertainty in the quality of the emissions data; the proximity of the projected emissions to key regulatory thresholds; the source category in question has not tested to verify emissions. Stack testing is not a research project. Stack testing is a compliance demonstration and is a core component of the air program.

The AQD and many applicants routinely use historical stack test data in the evaluation of permit applications. If the data is representative of a similar process, an applicant may use it in quantifying their emissions. AQD may also use this data as a way to determine if emissions, as presented by the applicant, are similar to what other sources have provided and/or demonstrated.

Over the past several years, AQD required stack testing to confirm toxic air contaminant emissions from new asphalt plants. Effective June 1, after an evaluation of the test results, AQD determined that routine testing of asphalt plants was no longer warranted.

The need for stack testing will be determined on a case by case basis. This is consistent with how AQD routinely evaluates the need for stack testing of various
source categories. AQD will continue to make such evaluations in the future, thus not requiring stack testing where it is not warranted.

While all stack test data submitted to AQD is public information and available to applicants for review and use, it is not currently easily accessible. AQD will work with the regulated community to develop a standardized stack test result submittal template. Also, with input from the regulated community, AQD will explore ways to post stack test results on the internet to increase accessibility.

Based upon the above, AQD believes that this item has been completely addressed and can be listed as resolved.


[^0]:    *TACs with ITSL values that were either: (1) greater than 75th percentile OR (2) default ITSL, AND there was no other reason for including in a new list of TACs.
    **All TACs with IRSLs are included in the new screening level list.

[^1]:    APPENDIX L: BENCHMARKING SURVEY OF STATE AIR TOXICS ASSESSMENTS IN NEW SOURCE PERMITTING

[^2]:    ${ }^{1}$ MDEQ-AQD, and presumably the air toxics permitting agencies of the other EPA R5 states, utilizes a general hierarchy system that is not rigidly applied; professional judgment and consideration of the age and basis for the available benchmarks and methods are important factors in adopting health-based screening levels that are appropriate and defensible.

