

(ii) The required mass removal is calculated by summing the required mass removal for all wastewater streams combined for treatment when complying with § 60.779(g)(1)(i) or (g)(2) of this subpart.

(5) *The AMR calculation procedure for non-combustion treatment processes including closed biological treatment processes.* The AMR shall be calculated as follows:

$$AMR = (QMW_a - QMW_b) \quad (\text{Eqn WW10})$$

Where:

AMR=Actual mass removal of VOC achieved by treatment process or series of treatment processes, kilograms per hour.

QMW<sub>a</sub>=Mass flow rate of VOC in wastewater entering the treatment process or first treatment process in a series of treatment processes, kilograms per hour.

QMW<sub>b</sub>=Mass flow rate of VOC in wastewater exiting the last treatment process in a series of treatment processes, kilograms per hour.

(6) *Compare RMR to AMR.* When complying with § 60.779(f)(2)(i) or (f)(3) of this subpart, compare the RMR calculated in Equation WW9 to the AMR calculated in Equation WW10. Compliance is demonstrated if the AMR is greater than or equal to the RMR. When complying with § 60.779(g)(1)(i) or (g)(2) of this subpart, compare the RMR calculated in Equation WW9a to the AMR calculated in Equation WW10. Compliance is demonstrated if the AMR is greater than or equal to 95-percent mass removal.

(f) *Open or closed aerobic biological treatment processes: Required mass removal (RMR) option.* This paragraph (f) applies to the use of performance tests that are conducted for open or closed aerobic biological treatment processes to demonstrate compliance with the mass removal provisions for VOC. These compliance options are specified in § 60.779(f)(2)(i) and (f)(2)(ii) of this subpart. The owner or operator shall comply with the requirements specified in paragraphs (f)(1) through (f)(6) of this section. Some compounds may not require a performance test. Refer to paragraph (h) of this section and Table 14 of this subpart to determine which compounds may be exempt from the requirements of this paragraph (f).

(1) *Concentration in wastewater stream.* The concentration of VOC shall be determined as provided in this paragraph (f)(1). Concentration measurements to determine RMR shall be taken at the point of determination or

downstream of the point of determination with adjustment for concentration change made according to § 60.782(b)(6) of this subpart.

Concentration measurements to determine AMR shall be taken at the inlet and outlet to the treatment process and as provided in paragraph (a)(7) of this section for a series of treatment processes. Wastewater samples shall be collected using sampling procedures which minimize loss of organic compounds during sample collection and analysis and maintain sample integrity per § 60.782(b)(5)(ii) of this subpart. The method shall be an analytical method for wastewater which has that compound as a target analyte. Samples may be grab samples or composite samples. Samples shall be taken at approximately equally spaced time intervals over a 1-hour period. Each 1-hour period constitutes a run, and the performance test shall consist of a minimum of 3 runs.

(2) *Flow rate.* Flow rate measurements to determine RMR shall be taken at the point of determination or downstream of the point of determination with adjustment for flow rate change made according to § 60.782(c)(4) of this subpart. Flow rate measurements to determine AMR shall be taken at the inlet and outlet to the treatment process and as provided in paragraph (a)(7) of this section for a series of treatment processes. Flow rate shall be determined using inlet and outlet flow measurement devices. Where the outlet flow is not greater than the inlet flow, a flow measurement device shall be used, and may be used at either the inlet or outlet. Flow rate measurements shall be taken at the same time as the concentration measurements.

(3) *Calculation of RMR for open or closed aerobic biological treatment processes.* The required mass removal of VOC for each Group 1 wastewater stream shall be calculated using the following equation:

$$RMR = \frac{\rho}{10^9} Q \sum_{i=1}^n (C_i * Fr_i) \quad (\text{Eqn WW11})$$

Where:

RMR=Required mass removal for treatment process or series of treatment processes, kilograms per hour.

ρ=Density of the Group 1 wastewater stream, kilograms per cubic meter.

Q=Volumetric flow rate of wastewater stream at the point of determination, liters per hour.

i=Identifier for a compound.

n=Number of VOC in stream.

C<sub>i</sub>=Concentration of VOC at the point of determination, parts per million by weight.

Fr<sub>i</sub>=Fraction removal value of a VOC. Follow the procedures in § 60.778 of this subpart to develop a stream-specific list of VOC. Follow the procedures in appendix J of this part to determine Fr values.

10<sup>9</sup>=Conversion factor, mg/kg \* l/m<sup>3</sup>.

(4) The required mass removal is calculated by adding together the required mass removal for each Group 1 wastewater stream to be combined for treatment.

(5) *Actual mass removal calculation procedure for open or closed aerobic biological treatment processes.* The actual mass removal (AMR) shall be calculated using Equation WW12 as specified in paragraph (f)(5)(i) of this section when the performance test is performed across the open or closed aerobic biological treatment process only. If compliance is being demonstrated in accordance with paragraph (a)(7)(i) of this section, the AMR for the series shall be calculated using Equation WW13 in paragraph (f)(5)(ii) of this section. (This equation is for situations where treatment is performed in a series of treatment processes connected by hard-piping.) If compliance is being demonstrated in accordance with paragraph (a)(7)(ii) of this section, the AMR for the biological treatment process shall be calculated using Equation WW12 in paragraph (f)(5)(i) of this section. The AMR for the biological treatment process used in a series of treatment processes calculated using Equation WW12 shall be added to the AMR determined for each of the other individual treatment processes in the series of treatment processes.

(i) Calculate AMR for the open or closed aerobic biological treatment process as follows:

$$AMR = QMW_a * F_{bio} \quad (\text{Eqn WW12})$$

Where:

AMR=Actual mass removal of VOC achieved by open or closed biological treatment process, kilograms per hour.

QMW<sub>a</sub>=Mass flow rate of VOC in wastewater entering the treatment process, kilograms per hour.

F<sub>bio</sub>=Site-specific fraction of VOC biodegraded. F<sub>bio</sub> shall be determined as specified in paragraph (h) of this section and 40 CFR part 63, appendix C. Follow the procedures in § 60.778 of this subpart to develop a stream-specific list of VOC.

(ii) Calculate AMR across a series of treatment units where the last treatment

unit is an open or closed aerobic biological treatment process as follows:

$$AMR = QMW_a - (QMW_b)(1 - F_{bio}) \quad (\text{Eqn WW13})$$

Where:

AMR=Actual mass removal of VOC achieved by a series of treatment processes, kilograms per hour.

QMW<sub>a</sub>=Mass flow rate of VOC in wastewater entering the first treatment process in a series of treatment processes, kilograms per hour.

QMW<sub>b</sub>=Mass flow rate of VOC in wastewater exiting the last treatment process in a series of treatment processes prior to the biological treatment process, kilograms per hour.

F<sub>bio</sub>=Site-specific fraction of VOC biodegraded. F<sub>bio</sub> shall be determined as specified in paragraph (h) of this section and 40 CFR part 63, appendix C. Follow the procedures in § 60.778 of this subpart to develop a stream-specific list of VOC.

(6) *Compare RMR to AMR.* Compare the RMR calculated in Equation WW11 to the AMR calculated in either Equation WW12 or WW13, as applicable. Compliance is demonstrated if the AMR is greater than or equal to the RMR.

(g) *Open or closed aerobic biological treatment processes: 95-percent mass removal option.* This paragraph (g) applies to performance tests that are conducted for open or closed aerobic biological treatment processes to demonstrate compliance with the 95-percent mass removal provisions for VOC. This compliance option is specified in § 60.779(g) of this subpart. The RMR for this option is 95-percent mass removal. The owner or operator shall comply with the requirements specified in paragraphs (g)(1) of this section to determine AMR, paragraphs (e)(3)(ii) and (e)(4)(ii) of this section to determine RMR, and paragraph (g)(2) of this section to determine whether compliance has been demonstrated. Some compounds may not require a performance test. Refer to paragraph (h) of this section and Table 14 of this subpart to determine which compounds may be exempt from the requirements of this paragraph (g).

(1) The owner or operator shall comply with the requirements specified in paragraphs (f)(1), (f)(2), and (f)(5) of this section to determine AMR. References to Group 1 wastewater streams shall be deemed all wastewater

streams combined for treatment for the purposes of this paragraph (g)(1).

(2) *Compare RMR to AMR.* Compliance is demonstrated if the AMR is greater than or equal to RMR.

(h) *Site-specific fraction biodegraded (F<sub>bio</sub>).* The VOC are divided into two sets for the purposes of determining whether F<sub>bio</sub> must be determined, and if F<sub>bio</sub> must be determined, which procedures may be used to determine compound-specific kinetic parameters. These sets are VOC in Table 14 of this subpart, and all other VOC.

(1) *Performance test exemption.* If a biological treatment process meets the requirements specified in paragraphs (h)(1)(i) and (h)(1)(ii) of this section, the owner or operator is not required to determine F<sub>bio</sub> and is exempt from the applicable performance test requirements specified in § 60.779 of this subpart.

(i) The biological treatment process meets the definition of "enhanced biological treatment process" in § 60.771 of this subpart.

(ii) At least 99 percent by weight of all VOC that are present in the aggregate of all wastewater streams using the biological treatment process to comply with § 60.779 of this subpart are compounds on Table 14 of this subpart.

(2) *F<sub>bio</sub> determination.* If a biological treatment process does not meet the requirement specified in paragraph (h)(1)(i) of this section, the owner or operator shall determine F<sub>bio</sub> for the biological treatment process using the procedures in 40 CFR part 63, appendix C, and paragraph (h)(2)(ii) of this section. If a biological treatment process meets the requirements of paragraph (h)(1)(i) of this section but does not meet the requirement specified in paragraph (h)(1)(ii) of this section, the owner or operator shall determine F<sub>bio</sub> for the biological treatment process using the procedures in 40 CFR part 63, appendix C, and paragraph (h)(2)(i) of this section.

(i) *Enhanced biological treatment processes.* If the biological treatment process meets the definition of "enhanced biological treatment process" in § 60.771 of this subpart and the wastewater streams include one or more compounds not on Table 14 of this subpart that do not meet the criteria in paragraph (h)(1)(ii) of this section, the owner or operator shall determine F<sub>bio</sub> for VOC not on Table 14 of this subpart using any of the procedures specified in

40 CFR part 63, appendix C. (stream-specific list) (The symbol F<sub>bio</sub> represents the site specific fraction of an individual VOC that is biodegraded.) The owner or operator shall calculate F<sub>bio</sub> for the VOC on Table 14 of this subpart using the defaults provided for first order biodegradation rate constants (K1) of this subpart and follow the procedure explained in Form III of 40 CFR part 63, appendix C, or any of the procedures specified in 40 CFR part 63, appendix C.

(ii) *Biological treatment processes that are not enhanced biological treatment processes.* For biological treatment processes that do not meet the definition for "enhanced biological treatment process" in § 60.771 of this subpart, the owner or operator shall determine the F<sub>bio</sub> for VOC on Table 14 of this subpart and all other VOC using any of the procedures in 40 CFR part 63, appendix C, except procedure 3 (inlet and outlet concentration measurements).

(i) *Performance tests for control devices other than flares.* This paragraph (i) applies to performance tests that are conducted to demonstrate compliance of a control device with the efficiency limits specified in § 60.780(c) of this subpart. If complying with the 95-percent reduction efficiency requirement, comply with the requirements specified in paragraphs (i)(1) through (i)(9) of this section. If complying with the 20 ppm by volume requirement, comply with the requirements specified in paragraphs (i)(1) through (i)(6) and (i)(9) of this section. The 20 ppm by volume limit or 95 percent reduction efficiency requirement shall be measured as either total VOC or as TOC minus methane and ethane.

(1) *Sampling sites.* Sampling sites shall be selected using Method 1 or 1A of appendix A of this part, as appropriate. For determination of compliance with the 95 percent reduction requirement, sampling sites shall be located at the inlet and the outlet of the control device. For determination of compliance with the 20 parts per million by volume limit, the sampling site shall be located at the outlet of the control device.

(2) *Concentration in gas stream entering or exiting the control device.* The concentration of total VOC or TOC

in a gas stream shall be determined as provided in this paragraph (i)(2). Samples may be grab samples or composite samples (i.e., integrated samples). Samples shall be taken at approximately equally spaced time intervals over a 1-hour period. Each 1-hour period constitutes a run, and the performance test shall consist of a minimum of 3 runs. Concentration measurements shall be determined using Method 18, 40 CFR part 60, appendix A. Alternatively, any other test method validated according to the procedures in Method 301, 40 CFR part 60, appendix A may be used.

(3) *Volumetric flow rate of gas stream entering or exiting the control device.* The volumetric flow rate of the gas stream shall be determined using Method 2, 2A, 2C, or 2D, 40 CFR part 60, appendix A, as appropriate. Volumetric flow rate measurements shall be taken at the same time as the concentration measurements.

(4) *Calculation of TOC concentration.* The TOC concentration (CGT) is the sum of the concentrations of the

individual components. If compliance is being determined based on TOC, the owner or operator shall compute TOC for each run using the following equation:

$$CG_T = \frac{1}{m} \sum_{j=1}^m \left( \sum_{i=1}^n CGS_{i,j} \right) \quad (\text{Eqn WW14})$$

Where:

CG<sub>T</sub>=Total concentration of TOC (minus methane and ethane) in vented gas stream, average of samples, dry basis, parts per million by volume.

CGS<sub>i,j</sub>=Concentration of sample components in vented gas stream for sample j, dry basis, parts per million by volume.

i=Identifier for a compound.

n=Number of components in the sample.

j=Identifier for a sample.

m=Number of samples in the sample run.

(5) *Calculation of total VOC concentration.* The owner or operator determining compliance based on total VOC concentration (CVOC) shall

compute C VOC according to the Equation WW14.

(6) *Percent oxygen correction for combustion control devices.* If the control device is a combustion device, comply with the requirements specified in paragraph (i)(6)(i) of this section to determine oxygen concentration, and in paragraph (i)(6)(ii) of this section to calculate the percent oxygen correction.

(i) *Oxygen concentration.* The concentration of TOC or total VOC shall be corrected to 3 percent oxygen if the control device is a combustion device. The emission rate correction factor for excess air, composite sampling (i.e., integrated sampling) and analysis procedures of Method 3B, 40 CFR part 60, appendix A shall be used to determine the actual oxygen concentration (%O<sub>2d</sub>). The samples shall be taken during the same time that the TOC (minus methane or ethane) or total VOC samples are taken.

(ii) *3 percent oxygen calculation.* The concentration corrected to 3 percent oxygen (CGc), when required, shall be computed using the following equation:

$$CG_c = CG_T \left( \frac{17.9}{20.9 - \%O_{2d}} \right) \quad (\text{Eqn WW15})$$

Where:

CG<sub>c</sub>=Concentration of TOC or VOC corrected to 3 percent oxygen, dry basis, parts per million by volume.

CG<sub>T</sub>=Total concentration of TOC (minus methane and ethane) in vented gas

stream, average of samples, dry basis, parts per million by volume.

%O<sub>2d</sub>=Concentration of oxygen measured in vented gas stream, dry basis, percent by volume.

(7) *Mass rate calculation.* The mass rate of either TOC (minus methane and

ethane) or total VOC shall be calculated using the following equations. Where the mass rate of TOC is being calculated, all organic compounds (minus methane and ethane) measured by methods specified in paragraph (i)(2) of this section are summed using Equations WW16 and WW17.

$$OMG_a = K_2 \left( \sum_{i=1}^n CG_{a,i} MW_i \right) QG_a \quad (\text{Eqn WW16})$$

$$OMG_b = K_2 \left( \sum_{i=1}^n CG_{b,i} MW_i \right) QG_b \quad (\text{Eqn WW17})$$

Where:

CG<sub>a,i</sub>, CG<sub>b,i</sub>=Concentration of TOC (minus methane and ethane) or total VOC, in vented gas stream, entering (CG<sub>a,i</sub>) and exiting (CG<sub>b,i</sub>) the control device, dry basis, parts per million by volume.

QMG<sub>a</sub>, QMG<sub>b</sub>=Mass rate of TOC (minus methane and ethane) or total VOC, in vented gas stream, entering (QMG<sub>a</sub>) and exiting (QMG<sub>b</sub>) the

control device, dry basis, kilograms per hour.

Mw<sub>i</sub>=Molecular weight of a component, kilogram/kilogram-mole.

QG<sub>a</sub>, QG<sub>b</sub>=Flow rate of gas stream entering (QG<sub>a</sub>) and exiting (QG<sub>b</sub>) the control device, dry standard cubic meters per hour.

K<sub>2</sub>=Constant, 41.57 x 10<sup>-9</sup> (parts per million)<sup>-1</sup> (gram-mole per standard cubic meter) (kilogram/gram), where standard temperature (gram-

mole per standard cubic meter) is 20° Celsius.

i=Identifier for a compound.

n=Number of components in the sample.

(8) *Percent reduction calculation.* The percent reduction in TOC (minus methane and ethane) or total VOC shall be calculated as follows:

$$E = \frac{QMG_a - QMG_b}{QMG_a} (100\%) \quad (\text{Eqn WW18})$$

Where:

E=Destruction efficiency of control device, percent.

$QMG_a, QMG_b$ =Mass rate of TOC (minus methane and ethane) or total VOC, in vented gas stream entering and exiting ( $QMG_b$ ) the control device, dry basis, kilograms per hour.

(9) *Compare mass destruction efficiency to required efficiency.* If complying with the 95 percent reduction efficiency requirement, compliance is demonstrated if the mass destruction efficiency (calculated in Equation WW18) is 95 percent or greater. If complying with the 20 parts per million by volume limit in § 60.780(c) of this subpart, compliance is demonstrated if the outlet total organic compound concentration, less methane and ethane, or total VOC concentration is 20 parts per million by volume, or less. For combustion control devices, the concentration shall be calculated on a dry basis, corrected to 3 percent oxygen.

(j) *Compliance demonstration for flares.* When a flare is used to comply with § 60.780(c) of this subpart, the owner or operator shall comply with the flare provisions in 40 CFR 63.11(b) and table 2A of this subpart, and with paragraphs (j)(1), (j)(2), and (j)(3) of this section. An owner or operator is not required to conduct a performance test to determine percent emission reduction or outlet VOC or TOC concentration when a flare is used. If a compliance demonstration has been conducted previously for a flare, using the techniques specified in paragraphs (h)(1) through (h)(3) of this section, that compliance demonstration may be used to satisfy the requirements of this paragraph (j) if either no deliberate process changes have been made since the compliance demonstration, or the results of the compliance demonstration reliably demonstrate compliance despite process changes.

(1) The compliance determination shall be conducted as specified in 40 CFR 63.11(b)(4) and table 2A of this subpart, to determine visible emissions.

(2) Determine the net heating value of the gas being combusted, using the techniques specified in 40 CFR 63.11(b)(6) and table 2A of this subpart; and

(3) Determine the exit velocity using the techniques specified in either 40 CFR 63.11(b)(7)(i) (and 40 CFR 63.11(b)(7)(iii), where applicable) or 40

CFR 63.11(b)(8), and table 2A of this subpart, as appropriate.

#### § 60.784 Reporting requirements.

(a) Owners or operators requesting approval to use alternative monitoring, recordkeeping, or reporting shall comply with the provisions in paragraph (b) of this section. Each owner or operator shall submit the reports specified in paragraphs (a)(1) through (a)(4) of this section, as applicable:

(1) Reports required by subpart A of part 60 of this part, as specified in table 2 of this subpart,

(2) Reports of certain subpart A provisions of 40 CFR part 63, as required by table 2A of this subpart,

(3) Reports required in paragraphs (c) through (g) of this section, and

(4) Start-up, shutdown, and malfunction reports specified in § 60.787 of this subpart.

(b) *Alternative monitoring and recordkeeping.* An owner or operator may request approval to use alternatives to the continuous operating parameter monitoring and recordkeeping provisions of this subpart.

(1) Requests for approval to use alternatives to the continuous monitoring and recordkeeping provisions shall be submitted prior to the implementation of the alternative monitoring system for which approval is being requested if not already included in the operating permit application. The request shall contain the information specified in paragraphs (b)(3) and (b)(4) of this section, as applicable.

(2) [Reserved]

(3) An owner or operator of an affected facility that does not have an automated monitoring and recording system capable of measuring parameter values at least once every 15 minutes and generating continuous records may request approval to use a non-automated system with less frequent monitoring.

(i) The requested system shall include manual reading and recording of the value of the relevant operating parameter no less frequently than once per hour. Daily average values shall be calculated from these hourly values and recorded.

(ii) The request shall contain:

(A) A description of the planned monitoring and recordkeeping system;

(B) Documentation that the affected facility does not have an automated monitoring and recording system;

(C) Justification for requesting an alternative monitoring and recordkeeping system; and

(D) Demonstration to the Administrator's satisfaction that the proposed monitoring frequency is sufficient to represent control device operating conditions considering typical variability of the specific process and control device operating parameter being monitored.

(4) An owner or operator may request approval to use an automated data compression recording system that does not record monitored operating parameter values at a set frequency (for example once every 15 minutes) but records all values that meet set criteria for variation from previously recorded values.

(i) The requested system shall be designed to:

(A) Measure the operating parameter value at least once every 15 minutes.

(B) Record at least four values each hour during periods of operation.

(C) Record the date and time when monitors are turned off or on.

(D) Recognize unchanging data that may indicate the monitor is not functioning properly, alert the operator, and record the incident.

(E) Compute daily average values of the monitored operating parameter based on recorded data.

(F) If the daily average is not an excursion, as defined in paragraphs (d)(3)(i) through (d)(3)(iii) of this section, the data for that operating day may be converted to hourly average values and the four or more individual records for each hour in the operating day may be discarded.

(ii) The request shall contain a description of the monitoring system and data compression recording system, including the criteria used to determine which monitored values are recorded and retained, the method for calculating daily averages, and a demonstration that the system meets all criteria in paragraph (b)(4)(i) of this section.

(5) [Reserved]

(6) For each waste management unit, treatment process, or control device used to comply with §§ 60.774 through 60.775 of this subpart for which the owner or operator seeks to monitor a parameter other than those specified in Table 5, Table 7, and Table 8 of this subpart, the owner or operator shall submit a request for approval to monitor alternative parameters. The owner or operator who requests approval to

monitor a different parameter than those listed in Table 5, Table 7, and Table 8 of this subpart shall submit the information specified in paragraphs (b)(6)(i), (ii), and (iii) of this section.

(i) A description of the parameter(s) to be monitored to ensure the waste management unit, treatment process, or control device measure is operated in conformance with its design and achieves the specified emission limit, percent reduction, or nominal efficiency, and an explanation of the criteria used to select the parameter(s).

(ii) A description of the methods and procedures that will be used to demonstrate that the parameter indicates proper operation of the waste management unit, treatment process, or control device, and the schedule for this demonstration, and a statement that the owner or operator will establish, as part of the demonstration, an operating parameter value for the monitored parameter that indicates proper operation and maintenance of the unit, process, or device.

(iii) The frequency and content of monitoring, recording, and reporting if monitoring and recording is not continuous, or if semiannual reports required under paragraph (d) of this section will not include reports of daily average values when the monitored operating parameter is not above or below (as appropriate) the operating parameter value established in paragraph (c)(7)(ii) of this section. The rationale for the proposed monitoring, recording, and reporting system shall be included.

(c) *Notification of Compliance Status.* Each owner or operator subject to this subpart shall submit a Notification of Compliance Status within 150 days after the compliance dates specified in § 60.770(a) of this subpart. The Notification of Compliance Status shall include the results of any emission point group determinations, performance tests, inspections, continuous monitoring system performance evaluations, values of monitored parameters established during performance tests, and any other information specified in paragraphs (c)(1) through (c)(14) of this section used to demonstrate compliance or required to be included in the Notification of Compliance Status.

(1) The owner or operator shall identify each designated CPU and list the components in the designated CPU. The owner or operator shall identify each affected facility and describe the process wastewater, maintenance wastewater, and aqueous in-process streams generated by the affected facility. The information shall clearly

link all applicable CPU, designated CPU, and affected facilities and demonstrate that all components of a CPU were assigned to a designated CPU.

(2) For each affected facility, the owner or operator shall submit the information specified in Table 9 of this subpart for each wastewater stream generated.

(3) For each treatment process identified in Table 9 of this subpart that receives, manages, or treats a wastewater stream (i.e., Group 1 wastewater stream or Group 2 wastewater stream selected by the owner or operator for control) or residual removed from a wastewater stream, the owner or operator shall submit the information specified in Table 10 of this subpart.

(4) For each waste management unit identified in Table 9 of this subpart that receives or manages a wastewater stream (i.e., Group 1 wastewater stream or Group 2 wastewater stream selected by the owner or operator for control) or residual removed from a wastewater stream, the owner or operator shall submit the information specified in Table 11 of this subpart.

(5) For each waste management unit identified in table 9 of this subpart, the owner or operator shall include in the Notification of Compliance Status the compliance option that will be used to comply with § 60.774 of this subpart, and the applicable provisions of other subparts that the owner or operator will use to comply with the compliance option, as allowed in § 60.774 of this subpart.

(6) For each residual removed from a wastewater stream (i.e., Group 1 wastewater stream or Group 2 wastewater stream selected by the owner or operator for control), the owner or operator shall submit the information specified in Table 12 of this subpart.

(7) For each control device used to comply with §§ 60.774, 60.775, and 60.779 of this subpart, the owner or operator shall submit the information specified in paragraphs (c)(7)(i) and (c)(7)(ii) of this section.

(i) For each flare, the owner or operator shall submit the information specified in paragraphs (c)(7)(i)(A) through (c)(7)(i)(C) of this section.

(A) Flare design (i.e., steam-assisted, air-assisted, or non-assisted);

(B) All visible emission readings, heat content determinations, flow rate measurements, and exit velocity determinations made during the compliance determination as specified by § 60.780(c)(3) of this subpart; and

(C) Reports of the times and durations of all periods during the compliance

determination when the pilot flame is absent or the monitor is not operating.

(ii) For each control device other than a flare, the owner or operator shall submit the information specified in paragraph (c)(7)(ii)(A) of this section and in either paragraph (c)(7)(ii)(B) or (c)(7)(ii)(C) of this section.

(A) The information in paragraphs (c)(7)(ii)(A)(1), (2), and (3) of this section on operating parameter values required to be established under § 60.781(f) of this subpart for the applicable parameters specified in Table 8 of this subpart, unless the operating parameter value has already been established in the operating permit.

(1) The specific operating parameter value of the monitored parameter(s) for each emission point;

(2) The rationale for the specific operating parameter value for each parameter for each emission point, including any data and calculations used to develop the value and a description of why the value indicates proper operation of the control device.

(i) If a performance test is conducted for a control device, the operating parameter value shall be based on the parameter values measured during the performance test supplemented by engineering analyses and/or manufacturer's recommendations. Performance testing is not required to be conducted over the entire range of permitted parameter values.

(ii) If a performance test is not conducted for a control device, the operating parameter value may be based solely on engineering analyses and/or manufacturer's recommendations.

(3) A definition of the affected facility's operating day for purposes of determining daily average values of monitored parameters. The definition shall specify the times at which an operating day begins and ends.

(B) The design evaluation specified in § 60.780(d)(2) of this subpart; or

(C) Results of the performance test specified in § 60.780(d)(1) of this subpart. Performance test results shall include operating ranges of key process and control parameters during the performance test; the value, averaged over the period of the performance test, of each parameter identified in the operating permit as being monitored in accordance with § 60.781 of this subpart; and applicable supporting calculations.

(8) For each treatment process used to comply with this subpart, the owner or operator shall submit the information specified in paragraphs (c)(8)(i) and (c)(8)(ii) of this section.

(i) For Items 1 and 2 in Table 7 of this subpart, the owner or operator shall

submit the information specified in paragraphs (c)(8)(i)(A) and (c)(8)(i)(B) of this section.

(A) The information specified in paragraph (c)(6)(ii)(A) of this section for the operating parameter value required to be established under § 60.781(f) of this subpart for the monitoring parameters approved by the Administrator, unless the operating parameter value has already been established in the operating permit.

(B) Results of the initial measurements of the parameters approved by the Administrator and any applicable supporting calculations.

(ii) For Item 3 in Table 7 of this subpart, the owner or operator shall submit the information specified in paragraph (c)(7)(ii)(A) of this section for the monitored operating parameter values required to be established under § 60.781(f) of this subpart, unless the operating parameter value has already been established in the operating permit.

(9) Except as provided in paragraph (c)(9)(iii) of this section, for each waste management unit or treatment process used to comply with this subpart, the owner or operator shall submit the information specified in either paragraph (c)(9)(i) or (c)(9)(ii) of this section.

(i) The design evaluation and supporting documentation specified in § 60.779(j)(1) of this subpart.

(ii) Results of the performance test specified in § 60.779(j)(2) of this subpart. Performance test results shall include operating ranges of key process and control parameters during the performance test; the value, averaged over the period of the performance test, of each parameter identified in the operating permit as being monitored in accordance with § 60.781(f) of this subpart; and applicable supporting calculations.

(iii) If the owner or operator elects to use one of the options for treatment in a RCRA unit specified in § 60.779(h) of this subpart, the owner or operator is exempt from the requirements specified in paragraphs (c)(9)(i) and (c)(9)(ii) of this section.

(10) For performance tests and group determinations that are based on measurements, and for estimates of VOC emissions, the Notification of Compliance Status shall include one complete test report for each test method used for a particular kind of emission point. For additional tests performed for the same kind of emission point using the same method, the results and any other information required shall be submitted, but a complete test report is not required. A complete test

report shall include a brief process description, sampling site description, description of sampling and analysis procedures and any modifications to standard procedures, quality assurance procedures, record of operating conditions during the test, record of preparation of standards, record of calibrations, raw data sheets for field sampling, raw data sheets for field and laboratory analyses, documentation of calculations, and any other information required by the test method.

(11) An owner or operator who transfers a Group 1 wastewater stream or residual removed from a Group 1 wastewater stream for treatment pursuant to § 60.773(e) shall include in the Notification of Compliance Status the name and location of the transferee and a description of the Group 1 wastewater stream or residual removed from a Group 1 wastewater stream sent to the treatment facility.

(12) The owner or operator who chooses to comply with the provisions in § 60.789 of this subpart shall include in the Notification of Compliance Status a statement specifying which regulation(s) is being used to comply with this subpart.

(13) Notification that the owner or operator has elected to comply with the reduced recordkeeping program in 60.785(j) of this subpart.

(14) Notification of the waste management unit compliance option used to comply with the provisions of this subpart, as specified in § 60.774 of this subpart, shall be submitted in the Notification of Compliance Status. If the owner or operator is complying with the recordkeeping and reporting provisions of a rule other than this subpart, as specified in § 60.774 of this subpart, a statement containing this information shall be submitted.

(d) *Semiannual reports.* Each owner or operator subject to the provisions of this subpart shall submit to the Administrator semiannual reports. The reports shall be submitted semiannually no later than 60 calendar days after the end of each 6-month period. The first report shall be submitted no later than 8 months after the due date of the notification of initial start-up required by § 60.7(a)(3) of this part and shall cover the 6-month period beginning on the due date of the notification of initial start-up.

(1) [Reserved]

(2) The semiannual report shall include reports of all excursions and all periods when monitoring parameters are above the maximum or below the minimum established value.

(3) The semiannual report shall include the daily average values of

monitored parameters for all excursions, as defined by paragraphs (d)(3)(i), (d)(3)(ii), or (d)(3)(iii) of this section. For excursions caused by lack of monitoring data, the duration of periods when monitoring data were not collected shall be reported. For a control device where multiple parameters are monitored, if one or more of the parameters meets the excursion criteria in paragraphs (d)(3)(i), (d)(3)(ii), or (d)(3)(iii) of this section, this is considered a single excursion for the control device.

(i) When the daily average value of one or more monitored parameters is above the maximum or below the minimum (as appropriate) established operating parameter value.

(ii) When the period of control device operation is 4 hours or greater in an operating day and monitoring data are insufficient to constitute a valid hour of data for at least 75 percent of the operating hours.

(iii) When the period of control device operation is less than 4 hours in an operating day and more than one of the hours during the period of operation does not constitute a valid hour of data due to insufficient monitoring data.

(iv) Monitoring data are insufficient to constitute a valid hour of data, as used in paragraphs (d)(3)(ii) and (d)(3)(iii) of this section, if measured values are unavailable for any of the 15-minute periods within the hour. For data compression systems approved under paragraph (b)(4) of this section, monitoring data are insufficient to calculate a valid hour of data if there are less than 4 data values recorded during the hour.

(4) Each control device is allowed one excused excursion per semiannual period. The first semiannual period is the 6-month period covered by the first semiannual report.

(5)(i) Paragraphs (d)(5)(i)(A) through (d)(5)(i)(D) of this section specify when an excursion is not a violation. In cases where continuous monitoring is required, the excursion does not count toward the number of excused excursions for determining compliance.

(A) If a monitored parameter is below the minimum established value and the affected facility is operated during such period in accordance with the affected facility's start-up, shutdown, and malfunction plan,

(B) If a monitored parameter is above the maximum established value and the affected facility is operated during such period in accordance with the affected facility's start-up, shutdown, and malfunction plan,

(C) If monitoring data are not collected during periods of start-up, shutdown, or malfunction and the

affected facility is operated during such period in accordance with the affected facility's start-up, shutdown, and malfunction plan, or

(D) If cessation of the emissions to which the monitoring applies occurs during periods of non-operation of the chemical process unit or portion thereof.

(ii) Nothing in paragraphs (d)(3) through (d)(5) of this section shall be construed to allow or excuse a monitoring parameter excursion caused by any activity that violates other applicable provisions of this subpart.

(iii) Paragraphs (d)(3) through (d)(5) of this section, except paragraph (d)(5)(i) of this section, shall apply only to emission points and control devices for which continuous monitoring is required by this subpart.

(6) The semiannual report shall include results of any performance tests conducted during the reporting period including one complete report for each test method used for a particular kind of emission point tested. For additional tests performed for a similar emission point using the same method, results and any other information required shall be submitted, but a complete test report is not required. A complete test report shall contain a brief process description, sampling site data, description of sampling and analysis procedures and any modifications to standard procedures, quality assurance procedures, record of operating conditions during the test, record of preparation of standards, record of calibrations, raw data sheets for field sampling, raw data sheets for field and laboratory analyses, documentation of calculations, and any other information required by the test method.

(7) The semiannual report shall include notification that the owner or operator has elected to comply with the reduced recordkeeping program in § 60.785(j) of this subpart.

(8) The semiannual report shall include notification that the owner or operator has elected not to retain the daily average values, as specified in § 60.785(j)(2)(i) of this subpart.

(9) The semiannual report shall include periods recorded under § 60.785(f)(10) of this subpart when the vent is diverted from the control device through a bypass line, with the next semiannual report.

(10) The semiannual report shall include notification of all occurrences recorded under § 60.785(f)(11) of this subpart in which the seal mechanism is broken, the bypass line damper or valve position has changed, or the key to unlock the bypass line damper or valve

was checked out, with the next semiannual report.

(11) The semiannual report shall include notification that semiannual report information for waste management units will be submitted with semiannual reports required by another rule that is one of the compliance options for waste management units as specified in § 60.784 of this subpart.

(12) The semiannual report shall include notification of each affected facility that ceases to generate at least one process wastewater stream or aqueous in-process stream or no longer produces a primary product that is a SOCOMI product.

(e) *Semiannual reporting for treatment processes.* Except as provided in paragraph (g) of this section, for each treatment process used to comply with this subpart, the owner or operator shall submit as part of the next semiannual report required by paragraph (d) of this section the information specified in paragraphs (e)(1) and (e)(2) of this section.

(1) For Item 1 in Table 7 of this subpart, the owner or operator shall submit the results of measurements that indicate that the biological treatment unit is outside the parameters established in the Notification of Compliance Status or operating permit.

(2) For Item 2 in Table 7 of this subpart, the owner or operator shall submit the monitoring results for each operating day during which the daily average value of any monitored parameter was above the maximum or below the minimum operating parameter value established in the Notification of Compliance Status or operating permit.

(3) For Item 3 in Table 7 of this subpart, the owner or operator shall submit the monitoring results for each operating day during which the daily average value of any monitored parameter specified in Item 3 of Table 7 of this subpart was above the maximum or below the minimum (as appropriate) operating parameter value established in the Notification of Compliance Status or operating permit.

(f) *Semiannual reporting for control devices.* Except as provided in paragraph (g) of this section, for each control device used to comply with §§ 60.774 through 60.780 of this subpart, the owner or operator shall submit as part of the next semiannual report required by paragraph (d) of this section the information specified in either paragraph (f)(1) or (f)(2) of this section.

(1) The information specified in Table 13 of this subpart, or

(2) If the owner or operator elects to comply with § 60.781(e)(2) of this subpart, i.e., an organic monitoring device installed at the outlet of the control device, the owner or operator shall submit the monitoring results for each operating day during which the daily average concentration level or reading is above the maximum or below the minimum (as appropriate) operating parameter value established as a requirement of § 60.781(f) of this subpart or established in the facility's operating permit.

(g) Where the owner or operator obtains approval to use a treatment process or control device other than one for which monitoring requirements are specified in § 60.781 of this subpart, or to monitor parameters other than those specified in Table 7 or 8 of this subpart, the owner or operator shall comply with the appropriate reporting requirements established by the Administrator.

#### § 60.785 Recordkeeping requirements.

(a) Data retention requirements are specified in paragraph (b) of this section. Each owner or operator shall keep the records specified in paragraphs (a)(1) through (a)(4) of this section, as applicable:

(1) Records required by subpart A of part 60 of this part, as specified in table 2 of this subpart,

(2) Records of certain subpart A provisions of 40 CFR part 63, as required by table 2A of this subpart,

(3) Records required in paragraphs (c) through (j) of this section, and

(4) Start-up, shutdown, and malfunction records specified in § 60.787 of this subpart.

(b) *Data retention.* Unless otherwise specified in this subpart, each owner or operator of an affected facility shall keep copies of all applicable records and reports required by this subpart for at least 5 years. All applicable records shall be maintained in such a manner that they can be readily accessed. Records of the most recent 2 years shall be retained onsite or shall be accessible to an inspector while onsite. The records of the remaining 3 years may be retained offsite. Records may be maintained in hard copy or computer-readable form including, but not limited to, on paper, microfilm, computer, floppy disk, magnetic tape, or microfiche.

(c) *Miscellaneous records.* The owner or operator shall keep the records specified in paragraphs (c)(1) through (c)(8) of this section.

(1) A record that each waste management unit inspection required by § 60.774 of this subpart was performed.



(2) A record that each inspection for control devices required by § 60.780(f) of this subpart was performed.

(3) For Item 1 and Item 2 of Table 7 of this subpart, the owner or operator shall keep the records approved by the Administrator.

(4) Except as provided in paragraph (c)(5) of this section, continuous records of the monitored parameters specified in Item 3 of Table 7, in Table 8, or in § 60.781(e)(2) of this subpart, as appropriate.

(5) Where the owner or operator obtains approval to use a treatment process or control device other than one for which monitoring requirements are specified in § 60.781 of this subpart, or to monitor parameters other than those specified in Table 7 or Table 8 of this subpart, the owner or operator shall comply with the recordkeeping requirements established by the Administrator as part of the review of the permit application or other appropriate means.

(6) The owner or operator who is complying with the provisions in § 60.789(c)(1) of this subpart shall keep a record of the information used to determine which control, testing, monitoring, recordkeeping, and reporting requirements are the most stringent.

(7) Documentation of a decision to use a delay of repair due to unavailability of parts, as specified in § 60.777(c) of this subpart, shall include a description of the failure, the reason additional time was necessary (including a statement of why replacement parts were not kept on site and when the manufacturer promised delivery), the date when repair would have been completed if parts had been available, and the date when repair was completed.

(8) The owner or operator shall keep a record of each affected facility that ceases to generate at least one process wastewater stream or aqueous in-process stream or no longer produces a primary product that is a SOCMI product.

(d) *Record of notice sent to treatment operator.* The owner or operator transferring a Group 1 wastewater stream or residual removed from a Group 1 wastewater stream in accordance with § 60.773(e) of this subpart shall keep a record of the notice sent to the treatment operator stating that the wastewater stream or residual contains VOC which are required to be managed and treated in accordance with the provisions of this subpart.

(e) *Control device records.* For each control device used to comply with this subpart, the owner or operator shall keep a record of the information

specified in paragraphs (e)(1) through (e)(3) of this section.

(1) Identification of all parts of the control device that are designated as unsafe to inspect, as specified in § 60.786(g) of this subpart, an explanation stating why the equipment is unsafe to inspect, and the plan for inspecting the equipment.

(2) Identification of all parts of the control device that are designated as difficult to inspect, as specified in § 60.786(h) of this subpart, an explanation stating why the equipment is difficult to inspect, and the plan for inspecting the equipment.

(3) For each boiler or process heater used to comply with this subpart, the owner or operator shall keep a record of any changes in the location at which the vent stream is introduced into the flame zone.

(f) *Continuous records.* Owners or operators required to keep continuous records by any section of this subpart shall keep records as specified in paragraphs (f)(1) through (f)(11) of this section, unless an alternative recordkeeping system has been requested and approved under § 60.784(b) of this subpart, except as provided in § 60.784(d)(5)(i) of this subpart.

(1) The monitoring system shall measure data values at least once every 15 minutes.

(2) The owner or operator shall record either:

(i) Each measured data value; or  
(ii) Block average values for 15-minute or shorter periods calculated from all measured data values during each period or at least one measured data value per minute if measured more frequently than once per minute.

(3) If the daily average value of a monitored parameter for a given operating day is below the maximum or above the minimum established value in the report required by § 60.784(c) of this subpart or the operating permit, the owner or operator shall either:

(i) Retain block hourly average values for that operating day for 5 years and discard, at or after the end of that operating day, the 15-minute or more frequent average values and readings recorded under paragraph (f)(2) of this section; or

(ii) Retain the data recorded in paragraph (f)(2) of this section for 5 years.

(4) If the daily average value of a monitored parameter for a given operating day is above the maximum or below the minimum established value in the report required by § 60.784(c) of this subpart or operating permit, the owner or operator shall retain the data

recorded that operating day under paragraph (f)(2) of this section for 5 years.

(5) Daily average values of each continuously monitored parameter shall be calculated for each operating day, and retained for 5 years, except as specified in paragraphs (f)(6) and (f)(7) of this section.

(i) The daily average shall be calculated as the average of all values for a monitored parameter recorded during the operating day. The average shall cover a 24-hour period if operation is continuous, or the number of hours of operation per operating day if operation is not continuous.

(ii) The operating day shall be the period defined in the operating permit or the report required by § 60.784(c) of this subpart. It may be from midnight to midnight or another daily period.

(6) If all recorded values for a monitored parameter during an operating day are below the maximum or above the minimum established value in the report required by § 60.784(c) of this subpart or operating permit, the owner or operator may record this fact and retain this record for 5 years rather than calculating and recording a daily average for that operating day. For these operating days, the records required in paragraph (f)(3) of this section shall also be retained for 5 years.

(7) Monitoring data recorded during periods identified in paragraphs (f)(7)(i) through (f)(7)(v) of this section shall not be included in any average computed under this subpart. Records shall be kept of the times and durations of all such periods and any other periods during process or control device operation when monitors are not operating.

(i) Monitoring system breakdowns, repairs, calibration checks, and zero (low-level) and high-level adjustments:

(ii) Start-ups;  
(iii) Shutdowns;  
(iv) Malfunctions;

(v) Periods of non-operation of the chemical process unit (or portion thereof), resulting in cessation of the emissions to which the monitoring applies.

(8) For flares, records of the times and duration of all periods during which all pilot flames are simultaneously absent shall be kept rather than daily averages.

(9) For carbon adsorbers, the owner or operator shall keep the records specified in paragraphs (e)(9)(i) and (e)(9)(ii) of this section instead of daily averages.

(i) Records of the total regeneration stream mass flow for each carbon bed regeneration cycle.



(ii) Records of the temperature of the carbon bed after each regeneration cycle.

(10) Hourly records of whether the flow indicator for bypass lines specified in § 60.786(f)(1) of this subpart was operating and whether a diversion was detected at any time during the hour. Also, records of the times of all periods when the vent is diverted from the control device or the flow indicator specified in § 60.786(f)(1) of this subpart is not operating.

(11) Where a seal or closure mechanism is used to comply with § 60.786(f)(2) of this subpart, hourly records of whether a diversion was detected at any time are not required. The owner or operator shall record whether the monthly visual inspection of the seals or closure mechanisms has been done, and shall record the occurrence of all periods when the seal mechanism is broken, the bypass line damper or valve position has changed, or the key for a lock-and-key type configuration has been checked out, and records of any car-seal that has broken.

(g) *Process knowledge records.* If the owner or operator determines that a wastewater stream is not a Group 1 wastewater stream by using process knowledge to determine the annual average concentration of a wastewater stream as specified in § 60.782(b)(3) of this subpart and/or uses process knowledge to determine the annual average flow rate as specified in § 60.782(c)(1) of this subpart, the owner or operator shall keep the documentation of how process knowledge was used to determine the annual average concentration and/or the annual average flow rate of the wastewater stream as specified in § 60.782(b)(3) or (c)(1) of this subpart, as appropriate.

(h) *Continuous monitoring system records.* For continuous monitoring systems used to comply with this subpart, records documenting the completion of calibration checks, and records documenting the maintenance of continuous monitoring systems that are specified in the manufacturer's instructions or that are specified in other written procedures that provide adequate assurance that the equipment would reasonably be expected to monitor accurately.

(i) [Reserved]

(j) *Reduced recordkeeping program.* For any parameter with respect to any item of equipment, the owner or operator may implement the recordkeeping requirements specified in paragraph (j)(1) or (j)(2) of this section as alternatives to the continuous operating parameter monitoring and

recordkeeping provisions specified in this subpart. The owner or operator shall retain for a period of 5 years each record required by paragraph (j)(1) or (j)(2) of this section.

(1) The owner or operator may retain only the daily average value, and is not required to retain more frequent monitored operating parameter values, for a monitored parameter with respect to an item of equipment, if the requirements of paragraphs (j)(1)(i) through (j)(1)(vi) of this section are met. An owner or operator electing to comply with the requirements of paragraph (j)(1) of this section shall notify the Administrator in the Notification of Compliance Status as specified in § 60.784(c)(13) of this subpart or, if the Notification of Compliance Status has already been submitted, in the semiannual report immediately preceding implementation of the requirements of paragraph (j)(1) of this section as specified in § 60.784(d)(7) of this subpart.

(i) The monitoring system is capable of detecting unrealistic or impossible data during periods of operation other than start-ups, shutdowns, or malfunctions (e.g., a temperature reading of  $-200^{\circ}\text{C}$  on a boiler), and will alert the operator by alarm or other means. The owner or operator shall record the occurrence. All instances of the alarm or other alert in an operating day constitute a single occurrence.

(ii) The monitoring system generates, updated at least hourly throughout each operating day, a running average of the monitoring values that have been obtained during that operating day, and the capability to observe this running average is readily available to the Administrator on-site during the operating day. The owner or operator shall record the occurrence of any period meeting the criteria in paragraphs (j)(1)(ii)(A) through (j)(1)(ii)(C) of this section. All instances in an operating day constitute a single occurrence.

(A) The running average is above the maximum or below the minimum established limits;

(B) The running average is based on at least six 1-hour average values; and

(C) The running average reflects a period of operation other than a start-up, shutdown, or malfunction.

(iii) The monitoring system is capable of detecting unchanging data during periods of operation other than start-ups, shutdowns, or malfunctions, except in circumstances where the presence of unchanging data is the expected operating condition based on past experience (e.g., pH in some scrubbers), and will alert the operator by alarm or

other means. The owner or operator shall record the occurrence. All instances of the alarm or other alert in an operating day constitute a single occurrence.

(iv) The monitoring system will alert the owner or operator by an alarm or other means, if the running average parameter value calculated under paragraph (j)(1)(ii) of this section reaches a set point that is appropriately related to the established limit for the parameter that is being monitored.

(v) The owner or operator shall verify the proper functioning of the monitoring system, including its ability to comply with the requirements of paragraph (j)(1) of this section, at the times specified in paragraphs (j)(1)(v)(A) through (j)(1)(v)(C). The owner or operator shall document that the required verifications occurred.

(A) Upon initial installation.

(B) Annually after initial installation.

(C) After any change to the programming or equipment constituting the monitoring system, which might reasonably be expected to alter the monitoring system's ability to comply with the requirements of this section.

(vi) The owner or operator shall retain the records identified in paragraphs (j)(1)(vi)(A) through (j)(1)(vi)(D) of this section.

(A) Identification of each parameter, for each item of equipment, for which the owner or operator has elected to comply with the requirements of paragraph (j) of this section.

(B) A description of the applicable monitoring system(s), and of how compliance will be achieved with each requirement of paragraphs (j)(1)(i) through (j)(1)(v) of this section. The description shall identify the location and format (e.g., on-line storage, log entries) for each required record. If the description changes, the owner or operator shall retain both the current and the most recent superseded description, as provided in paragraph (a) of this section, except as provided in paragraph (j)(1)(vi)(D) of this section.

(C) A description, and the date, of any change to the monitoring system that would reasonably be expected to impair its ability to comply with the requirements of paragraph (j)(1) of this section.

(D) Owners and operators subject to paragraph (j)(1)(vi)(B) of this section shall retain the current description of the monitoring system as long as the description is current, but not less than 5 years from the date of its creation. The current description shall, at all times, be retained on-site or be accessible from a central location by computer or other means that provides access within 2

hours after a request. The owner or operator shall retain all superseded descriptions for at least 5 years after the date of their creation. Superseded descriptions shall be retained on-site (or accessible from a central location by computer or other means that provides access within 2 hours after a request) for at least 6 months after their creation. Thereafter, superseded descriptions may be stored off-site.

(2) If an owner or operator has elected to implement the requirements of paragraph (j)(1) of this section for a monitored parameter with respect to an item of equipment and a period of 6 consecutive months has passed without an excursion as defined in paragraph (j)(2)(iv) of this section, the owner or operator is no longer required to record the daily average value for any operating day when the daily average value is less than the maximum or greater than the minimum established limit. With approval by the Administrator, monitoring data generated prior to the compliance date of this subpart shall be credited toward the period of 6 consecutive months, if the parameter limit and the monitoring accomplished during the period prior to the compliance date was required and/or approved by the Administrator.

(i) If the owner or operator elects not to retain the daily average values, the owner or operator shall notify the Administrator in the next semiannual report as specified in § 60.784(d)(8) of this subpart. The notification shall identify the parameter and unit of equipment.

(ii) If, on any operating day after the owner or operator has ceased recording daily average values as provided in paragraph (j)(2) of this section, there is an excursion as defined in paragraph (j)(2)(iv) of this section, the owner or operator shall immediately resume retaining the daily average value for each operating day and shall notify the Administrator in the next semiannual report. The owner or operator shall continue to retain each daily average value until another period of 6 consecutive months has passed without an excursion as defined in paragraph (j)(2)(iv) of this section.

(iii) The owner or operator shall retain the records specified in paragraphs (j)(1)(i) through (j)(1)(iv) of this section, for the duration specified in paragraph (j) of this section. For any calendar week, if compliance with paragraphs (j)(1)(i) through (j)(1)(iv) of this section does not result in retention of a record of at least one occurrence or measured parameter value, the owner or operator shall record and retain at least one parameter value during a period of

operation other than a start-up, shutdown, or malfunction.

(iv) For purposes of paragraph (j) of this section, an excursion means that the daily average value of monitoring data for a parameter is greater than the maximum, or less than the minimum established value, except that the daily average value during any start-up, shutdown, or malfunction shall not be considered an excursion for purposes of paragraph (j)(2) of this section, if the owner or operator follows the applicable provisions of the start-up, shutdown, and malfunction plan required by § 60.787 of this subpart. An excused excursion, as described in § 60.784(d)(4) of this subpart, shall not be considered an excursion for purposes of this paragraph (j)(2).

#### § 60.786 Leak inspection provisions.

(a) For each vapor collection system, closed-vent system, fixed roof, cover, or enclosure required to comply with this section, the owner or operator shall comply with the requirements of paragraphs (b) through (i) of this section, unless otherwise specified in this subpart.

(b) Except as provided in paragraphs (g) and (h) of this section, each vapor collection system and closed-vent system shall be inspected according to the procedures and schedule specified in paragraphs (b)(1) and (b)(2) of this section and each fixed roof, cover, and enclosure shall be inspected according to the procedures and schedule specified in paragraph (b)(3) of this section.

(1) If the vapor collection system or closed vent system is constructed of hard-piping, the owner or operator shall:

- (i) Conduct an initial inspection according to the procedures in paragraph (c) of this section; and
- (ii) Conduct annual visual inspections for visible, audible, or olfactory indications of leaks.

(2) If the vapor collection system or closed vent system is constructed of duct work, the owner or operator shall:

- (i) Conduct an initial inspection according to the procedures in paragraph (c) of this section, and
- (ii) Conduct annual visual inspections for visible, audible, or olfactory indications of leaks.

(2) If the vapor collection system or closed vent system is constructed of duct work, the owner or operator shall:

- (i) Conduct an initial inspection according to the procedures in paragraph (c) of this section; and
- (ii) Conduct annual inspections according to the procedures in paragraph (c) of this section.

(iii) Conduct annual visual inspections for visible, audible, or olfactory indications of leaks.

(3) For each fixed roof, cover, and enclosure, the owner or operator shall:

(i) Conduct an initial inspection according to the procedures in paragraph (c) of this section; and

(ii) Conduct semi-annual visual inspections for visible, audible, or olfactory indications of leaks.

(c) Each vapor collection system, closed vent system, fixed roof, cover, and enclosure shall be inspected according to the procedures specified in paragraphs (c)(1) through (c)(6) of this section.

(1) Inspections shall be conducted in accordance with Method 21, 40 CFR part 60, appendix A, and with the exceptions and modifications specified in this subpart.

(2) The detection instrument shall meet the performance criteria of Method 21, 40 CFR part 60, appendix A except the instrument response factor criteria in Section 3.1.2(a) of Method 21 shall be for the average composition of the process fluid not each individual VOC in the stream.

(i) Except as provided in paragraph (c)(2)(ii) of this section, the detection instrument shall meet the performance criteria of Method 21, 40 CFR part 60, appendix A, except the instrument response factor criteria in section 3.1.2(a) of Method 21 shall be for the average composition of the process fluid not each individual VOC in the stream. For process streams that contain nitrogen, air, or other inerts which are not organic hazardous air pollutants or VOCs, the average stream response factor shall be calculated on an inert-free basis.

(ii) If no instrument is available at the plant site that will meet the performance criteria specified in paragraph (c)(2)(i) of this section, the instrument readings may be adjusted by multiplying by the average response factor of the process fluid, calculated on an inert-free basis as described in paragraph (c)(2)(i) of this section.

(3) The detection instrument shall be calibrated before use on each day of its use by the procedures specified in Method 21, 40 CFR part 60, appendix A.

(4) Calibration gases shall be as follows:

- (i) Zero air (less than 10 parts per million hydrocarbon in air); and
- (ii) Mixtures of methane in air at a concentration less than 10,000 parts per million. A calibration gas other than methane in air may be used if the instrument does not respond to methane or if the instrument does not meet the performance criteria specified in

paragraph (c)(2)(i) of this section. In such cases, the calibration gas may be a mixture of one or more of the compounds to be measured in the air.

(5) An owner or operator may elect to adjust or not adjust instrument readings for background. If an owner or operator elects to not adjust readings for background, all such instrument readings shall be compared directly to the applicable leak definition to determine whether there is a leak. If an owner or operator elects to adjust instrument readings for background, the owner or operator shall measure background concentration using the procedures in 40 CFR 63.180(b) and (c). The owner or operator shall subtract background reading from the maximum concentration indicated by the instrument.

(6) The arithmetic difference between the maximum concentration indicated by the instrument and the background level shall be compared with 500 parts per million for determining compliance.

(d) Leaks, as indicated by an instrument reading greater than 500 parts per million above background or by visual inspections, shall be repaired as soon as practical, except as provided in paragraph (e) of this section.

(1) A first attempt at repair shall be made no later than 5 calendar days after the leak is detected.

(2) Repair shall be completed no later than 15 calendar days after the leak is detected.

(e) Delay of repair of a vapor collection system, closed vent system, fixed roof, cover, or enclosure for which leaks have been detected is allowed if the repair is technically infeasible without a shutdown or if the owner or operator determines that emissions resulting from immediate repair would be greater than the fugitive emissions likely to result from delay of repair. Repair of such equipment shall be complete by the end of the next shutdown.

(f) For each vapor collection system or closed vent system that contains bypass lines that could divert emissions away from a control device, the owner or operator shall comply with the provisions of either paragraph (f)(1) or (f)(2) of this section. Equipment such as low leg drains, high point bleeds, analyzer vents, open-ended valves or lines, and pressure relief valves needed for safety purposes are not subject to this paragraph (f).

(1) Properly install, maintain, and operate a flow indicator that takes a reading at least once every 15 minutes. Records shall be generated as specified in § 60.785(f)(10) of this subpart. The flow indicator shall be installed at the

entrance to any bypass line that could divert emissions away from the control device and to the atmosphere; or

(2) Secure the bypass line damper or valve in the non-diverting position with a car-seal or a lock-and-key type configuration. A visual inspection of the seal or closure mechanism shall be performed at least once every month to ensure that the damper or valve is maintained in the non-diverting position and emissions are not diverted through the bypass line. Records shall be generated as specified in § 60.785(e)(11) of this subpart.

(g) Any parts of the vapor collection system, closed vent system, fixed roof, cover, or enclosure that are designated, as described in paragraph (i)(1) of this section, as unsafe to inspect are exempt from the inspection requirements of paragraphs (b)(1), (b)(2), and (b)(3)(i) of this section if:

(1) The owner or operator determines that the equipment is unsafe to inspect because inspecting personnel would be exposed to an imminent or potential danger as a consequence of complying with paragraphs (b)(1), (b)(2), or (b)(3)(i) of this section; and

(2) The owner or operator has a written plan that requires inspection of the equipment as frequently as practicable during safe-to-inspect times.

(h) Any parts of the vapor collection system, closed vent system, fixed roof, cover, or enclosure that are designated, as described in paragraph (i)(2) of this section, as difficult to inspect are exempt from the inspection requirements of paragraphs (b)(1), (b)(2), and (b)(3)(i) of this section if:

(1) The owner or operator determines that the equipment cannot be inspected without elevating the inspecting personnel more than 2 meters above a support surface; and

(2) The owner or operator has a written plan that requires inspection of the equipment at least once every 5 years.

(i) The owner or operator shall record the information specified in paragraphs (i)(1) through (i)(5) of this section.

(1) Identification of all parts of the vapor collection system, closed vent system, fixed roof, cover, or enclosure that are designated as unsafe to inspect, an explanation of why the equipment is unsafe to inspect, and the plan for inspecting the equipment.

(2) Identification of all parts of the vapor collection system, closed vent system, fixed roof, cover, or enclosure that are designated as difficult to inspect, an explanation of why the equipment is difficult to inspect, and the plan for inspecting the equipment.

(3) For each vapor collection system or closed vent system that contains bypass lines that could divert a vent stream away from the control device and to the atmosphere, the owner or operator shall keep a record of the information specified in either paragraph (i)(3)(i) or (i)(3)(ii) of this section.

(i) Hourly records of whether the flow indicator specified under paragraph (f)(1) of this section was operating and whether a diversion was detected at any time during the hour, as well as records of the times of all periods when the vent stream is diverted from the control device or the monitor is not operating.

(ii) Where a seal mechanism is used to comply with paragraph (f)(2) of this section, hourly records of flow are not required. In such cases, the owner or operator shall record whether the monthly visual inspection of the seals or closure mechanisms has been done, and shall record the occurrence of all periods when the seal mechanism is broken, the bypass line valve position has changed, or the key for a lock-and-key type configuration has been checked out, and records of any car-seal that has broken.

(4) For each inspection during which a leak is detected, a record of the information specified in paragraphs (i)(4)(i) through (i)(4)(viii) of this section.

(i) The instrument identification numbers; the name or initials of the person conducting the inspection; and identification of the equipment.

(ii) The date the leak was detected and the date of the first attempt to repair the leak.

(iii) Maximum instrument reading measured by the method specified in paragraph (d) of this section after the leak is successfully repaired or determined to be nonreparable.

(iv) "Repair delayed" and the reason for the delay if a leak is not repaired within 15 calendar days after discovery of the leak.

(v) The name, initials, or other form of identification of the owner or operator (or designee) whose decision it was that repair could not be effected without a shutdown.

(vi) The expected date of successful repair of the leak if a leak is not repaired within 15 calendar days.

(vii) Dates of shutdowns that occur while the equipment is unrepaired.

(viii) The date of successful repair of the leak.

(5) For each inspection conducted in accordance with paragraph (c) of this section during which no leaks are detected, a record that the inspection was performed, the date of the

inspection, and a statement that no leaks were detected.

(6) For each visual inspection conducted in accordance with paragraph (b)(1)(ii) or (b)(3)(ii) of this section during which no leaks are detected, a record that the inspection was performed, the date of the inspection, and a statement that no leaks were detected.

**§ 60.787 Additional Requirements—Start-up, Shutdown, Malfunction, or Nonoperation; Alternative Means of Emission Limitation; and Permits**

(a) *Applicability of this subpart during periods of start-up, shutdown, malfunction, or non-operation.*

Paragraphs (a)(1) through (a)(4) of this section shall be followed during periods of start-up, shutdown, malfunction, or non-operation of the affected facility or any part thereof.

(1) The emission limitations set forth in this subpart and the emission limitations referred to in this subpart shall apply at all times except during periods of non-operation of the affected facility (or specific portion thereof) resulting in cessation of the emissions to which this subpart applies. The emission limitations of this subpart and the emission limitations referred to in this subpart shall not apply during periods of start-up, shutdown, or malfunction. During periods of start-up, shutdown, or malfunction, the owner or operator shall follow the applicable provisions of the start-up, shutdown, and malfunction plan as specified in 40 CFR 63.6(e)(3) and table 2A of this subpart. However, if a start-up, shutdown, malfunction, or period of non-operation of one portion of an affected facility does not affect the ability of a particular emission point to comply with the emission limitations to which it is subject, then that emission point shall still be required to comply with the applicable provisions of this subpart during the start-up, shutdown, malfunction, or period of non-operation.

(2) The owner or operator shall not shut down items of equipment that are required or utilized for compliance with this subpart during periods of start-up, shutdown, or malfunction during times when emissions, wastewater streams, or residuals are being routed to such items of equipment, if the shutdown would contravene requirements of this subpart applicable to such items of equipment. This paragraph (a)(2) does not apply if the item of equipment is malfunctioning. This paragraph (a)(2) also does not apply if the owner or operator shuts down the compliance equipment (other than monitoring systems) to avoid damage due to a

contemporaneous start-up, shutdown, or malfunction of the affected facility or portion thereof. If the owner or operator has reason to believe that monitoring equipment would be damaged due to a contemporaneous start-up, shutdown, or malfunction of the affected facility or portion thereof, the owner or operator shall provide documentation to the Administrator, as soon as possible, supporting such a claim. Once approved by the Administrator, the provision for ceasing to collect, during a start-up, shutdown, or malfunction, monitoring data that would otherwise be required by the provisions of this subpart must be incorporated into the start-up, shutdown, malfunction plan for that affected facility.

(3) During start-ups, shutdowns, and malfunctions when the emission limitations of this subpart do not apply pursuant to paragraphs (a)(1) and (a)(2) of this section, the owner or operator shall implement, to the extent reasonably available, measures to prevent or minimize excess emissions. For purposes of this paragraph (a)(3), the term "excess emissions" means emissions in excess of those that would have occurred if there were no start-up, shutdown, or malfunction and the owner or operator complied with the relevant provisions of this subpart. The measures to be taken shall be identified in the applicable start-up, shutdown, and malfunction plan, and may include, but are not limited to, air pollution control technologies, recovery technologies, work practices, pollution prevention, monitoring, and/or changes in the manner of operation of the affected facility. Back-up control devices are not required, but may be used if available.

(b) *Start-up, shutdown, and malfunction plan.* The owner or operator of an affected facility shall develop and implement a written start-up, shutdown, and malfunction plan as specified in 40 CFR 63.6(e)(3) and table 2A of this subpart. This plan shall describe, in detail, procedures for operating and maintaining the affected facility during periods of start-up, shutdown, and malfunction and a program for corrective action for malfunctioning process and air pollution control equipment used to comply with this subpart. A provision for ceasing to collect, during a start-up, shutdown, or malfunction, monitoring data that would otherwise be required by the provisions of this subpart may be included in the start-up, shutdown, and malfunction plan only if the owner or operator has demonstrated to the Administrator that the monitoring system would be damaged or destroyed

if it were not shut down during the start-up, shutdown, or malfunction. The affected facility shall keep the start-up, shutdown, and malfunction plan on-site.

(1) *Records of start-up, shutdown, and malfunction.* The owner or operator shall keep the records specified in paragraphs (b)(1)(i) and (b)(1)(ii) of this section.

(i) Records of the occurrence and duration of each start-up, shutdown, and malfunction of operation of process equipment or control devices or recovery devices or continuous monitoring systems used to comply with this subpart during which excess emissions (as defined in paragraph (a)(3) of this section) occur.

(ii) For each start-up, shutdown, or malfunction during which excess emissions (as defined in paragraph (a)(3) of this section) occur, records reflecting whether the procedures specified in the affected facility's start-up, shutdown, and malfunction plan were followed, and documentation of actions taken that are not consistent with the plan. For example, if a start-up, shutdown, and malfunction plan includes procedures for routing a control device to a backup control device, records shall be kept of whether the plan was followed. These records may take the form of a "checklist," or other form of recordkeeping that confirms conformance with the start-up, shutdown, and malfunction plan for the event.

(2) *Reports of start-up, shutdown, and malfunction.* For the purposes of this subpart, the start-up, shutdown, and malfunction reports shall be submitted on the same schedule as the semiannual reports required under § 60.784(d) of this subpart. Said reports shall include the information specified in paragraphs (b)(1)(i) and (b)(1)(ii) of this section and shall contain the name, title, and signature of the owner or operator or other responsible official who is certifying its accuracy.

(b) *Alternative means of emission limitation.* If, in the judgment of the Administrator, an alternative means of emission limitation will achieve a reduction in VOC emissions at least equivalent to the reduction in VOC achieved under any design, equipment, work practice, or operational standards in this subpart, the Administrator will publish a notice permitting the use of the alternative means for purposes of compliance with that requirement.

(1) The notice may condition the permission on requirements related to the operation and maintenance of the alternative means.

(2) Any notice under paragraph (b) of this section shall be published only after public notice and an opportunity for a hearing.

(3) Any person seeking permission to use an alternative means of compliance under this section shall collect, verify, and submit to the Administrator information showing that the alternative means achieves equivalent emission reductions.

(c) *Permit.* Each owner or operator of an affected facility subject to this subpart shall obtain a permit under 40 CFR part 70 or part 71 from the appropriate permitting authority. If EPA has approved a State operating permit program under 40 CFR part 71, the permit shall be obtained from the State authority. If the State operating permit program has not been approved, the owner or operator shall apply to the EPA regional office pursuant to 40 CFR part 70.

**§ 60.788 [Reserved]**

**§ 60.789 Relationship to other regulations.**

(a) The owner or operator who is subject to the provisions of this section shall include in the Notification of Compliance Status a statement specifying the options being used to comply with the provisions of this section.

(b) *Relationship to benzene waste.* After the compliance dates specified in § 60.770 of this subpart, the owner or operator of a Group 1 or Group 2 wastewater stream that is also subject to the provisions of 40 CFR part 61, subpart FF is required to comply with the provisions of both this subpart and 40 CFR part 61, subpart FF. Alternatively, the owner or operator may elect to comply with the provisions of paragraphs (b)(1) and (b)(2) of this section, which shall constitute compliance with the provisions of 40 CFR part 61, subpart FF.

(1) Comply with the provisions of this subpart; and

(2) For any Group 2 wastewater stream or organic stream whose benzene emissions are subject to control through the use of one or more treatment processes or waste management units under the provisions of 40 CFR part 61, subpart FF on or after September 12, 1994, comply with the requirements of

this subpart for Group 1 wastewater streams.

(c) *Relationship to RCRA.* After the compliance dates specified in § 60.770 of this subpart, the owner or operator of any Group 1 or Group 2 wastewater stream that is also subject to provisions in 40 CFR parts 260 and 272 shall comply with the requirements of either paragraph (c)(1) or (c)(2) of this section.

(1) For each Group 1 or Group 2 wastewater stream, the owner or operator shall comply with the more stringent control requirements (e.g., waste management units, numerical treatment standards, etc.) and the more stringent testing, monitoring, recordkeeping, and reporting requirements that overlap between the provisions of this subpart and the provisions of 40 CFR parts 260 through 272. The owner or operator shall keep a record of the information used to determine which requirements were the most stringent and shall submit this information if requested by the Administrator; or

(2) The owner or operator shall submit, no later than four months before the applicable compliance date specified in § 60.770 of this subpart, a request for a case-by-case determination of requirements. The request shall include the information specified in paragraphs (c)(2)(i) and (c)(2)(ii) of this section.

(i) Identification of the wastewater streams that are subject to this subpart and to the provisions in 40 CFR parts 260 through 272, determination of the Group 1/Group 2 status of those streams using the provisions specified in this subpart, determination of whether or not those streams are listed or exhibit a characteristic as specified in 40 CFR part 261, and determination of whether the waste management unit is subject to permitting under 40 CFR part 270.

(ii) Identification of the specific control requirements (e.g., waste management units, numerical treatment standards, etc.) and testing, monitoring, recordkeeping, and reporting requirements that overlap between the provisions of this subpart and the provisions of 40 CFR parts 260 through 272.

(d) *Overlap with the Vinyl Chloride NESHAP.* After the compliance dates specified in § 60.770 of this subpart, the

owner or operator of a Group 1 and Group 2 wastewater stream that is also subject to the provisions of 40 CFR part 61 subpart F shall comply with the provisions of either paragraph (d)(1) or (d)(2) of this section.

(1) The owner or operator shall comply with the provisions of both this subpart and 40 CFR part 61 subpart F or

(2) The owner or operator may submit, no later than four months before the applicable compliance date specified in § 60.770 of this subpart, information demonstrating how compliance with 40 CFR Part 61, subpart F, will also ensure compliance with this subpart. The information shall include a description of the testing, monitoring, reporting, and recordkeeping that will be performed.

(e) *Overlap with the HON.* After the compliance dates specified in § 60.770 of this subpart, the owner or operator of any Group 1 or Group 2 process wastewater stream that is also subject to and controlled according to the provisions in 40 CFR, subpart G shall comply with either 40 CFR, subpart G or this subpart.

(f) *Overlap with other regulations for monitoring, recordkeeping, or reporting with respect to combustion devices, recovery devices, or recapture devices.* After compliance dates specified in § 60.770 of this subpart, if any combustion device, recovery device, or recapture device subject to this subpart is also subject to monitoring, recordkeeping, and reporting requirements in 40 CFR part 264, subpart AA or CC, or is subject to monitoring and recordkeeping requirements in 40 CFR part 265, subpart AA or CC and other owner or operator complies with the periodic reporting requirements under 40 CFR part 264, subpart AA or CC that would apply to the device if the facility had final-permitted status, the owner or operator may elect to comply with the monitoring, recordkeeping, and reporting requirements of this subpart, or with the monitoring, recordkeeping, and reporting requirements in 40 CFR parts 264 and/or 265, as described in this paragraph (f), which shall constitute compliance with the monitoring, recordkeeping, and reporting requirements of this subpart.

TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS

| Chemical name <sup>a</sup>                        | CAS No. <sup>b</sup> |
|---|----------------------|
| (1,1,2-) Trichloro (1,2,2-) trifluoroethane ..... | 76131                |
| (2-Ethylhexyl) amine .....                        | 104756               |
| 1,4-Dichlorobutene .....                          | 110576               |
| 1-Butene .....                                    | 106989               |

TABLE 1 TO SUBPART YYY—LIST OF SOCM1 CHEMICALS—Continued

| Chemical name <sup>a</sup>                                  | CAS No. <sup>b</sup> |
|---|----------------------|
| 1-Methyl-2-pyrrolidone .....                                | 872504               |
| 1-Naphthyl-N-methylcarbamate .....                          |                      |
| 1-Phenyl ethyl hydroperoxide .....                          | 3071327              |
| 2-Butene .....  | 25167673             |
| 2-Butyne-1,4-diol .....                                     | 110656               |
| 2-Chloro-1,3-butadiene (Chloroprene) .....                  | 126998               |
| 2-Chloro-4-(ethylamino)-6-(isopropylamino)-S-triazine ..... | 1912249              |
| 2-Ethylhexanol (2-ethyl-1-hexanol) .....                    | 104767               |
| 2-Hexenedinitrile .....                                     | 13042029             |
| 3,4-Dichloro-1-butene .....                                 | 64037543             |
| 3-Hexenedinitrile .....                                     | 1119853              |
| 3-Pentenenitrile .....                                      | 4635874              |
| 6-Ethyl-1,2,3,4-tetrahydro-9,10-antracenedione .....        | 15547178             |
| Acenaphthene .....  | 83329                |
| Acetal (1,1-diethoxy-ethane) .....                          | 105577               |
| Acetaldehyde .....  | 75070                |
| Acetaldol (3-hydroxy-butanal) .....                         | 107891               |
| Acetamide .....   | 60355                |
| Acetanilide .....   | 103844               |
| Acetic anhydride .....                                      | 108247               |
| Acetic acid .....   | 64197                |
| Acetoacetanilide .....                                      | 102012               |
| Acetone cyanohydrin .....                                   | 75865                |
| Acetone .....   | 67641                |
| Acetonitrile .....  | 75058                |
| Acetophenone .....  | 98862                |
| Acetyl chloride .....                                       | 75365                |
| Acetylene tetrabromide (1,1,2,2-tetrabromoethane) .....     | 79276                |
| Acetylene .....   | 74862                |
| Acrolein .....  | 107028               |
| Acrylamide .....  | 79061                |
| Acrylic acid .....  | 79107                |
| Acrylonitrile .....   | 107131               |
| Adipic acid .....   | 124049               |
| Adiponitrile .....  | 111693               |
| Alcohols, C-11 or higher, mixtures .....                    |                      |
| Alcohols, C-11 or lower, mixtures .....                     |                      |
| Alizarin .....  | 72480                |
| Alkyl naphthalenes .....                                    |                      |
| Alkyl naphthalene sulfonates .....                          |                      |
| Alkyl anthraquinones .....                                  |                      |
| Allyl cyanide .....   | 109751               |
| Allyl chloride .....  | 107051               |
| Allyl bromide .....   | 106956               |
| Allyl alcohol .....   | 107186               |
| Aluminum acetate .....                                      | 7360443              |
| Aluminum formates .....                                     |                      |
| Aminobenzoic acid (p-) .....                                | 1321115              |
| Aminoethylethanolamine .....                                | 111411               |
| Aminophenol sulfonic acid .....                             |                      |
| Aminophenol (p-) .....                                      | 123308               |
| Ammonium acetate .....                                      | 631618               |
| Ammonium thiocyanate .....                                  | 1762954              |
| Amyl acetates .....   | 628637               |
|   | 123922               |
| Amyl chloride (n-) .....                                    | 543599               |
| Amyl phenol .....   | 1322061              |
| Amyl chlorides (mixed) .....                                |                      |
| Amyl mercaptans .....                                       | 110667               |
| Amyl alcohols (mixed) .....                                 | 30899195             |
| Amyl alcohol (tert-) .....                                  | 75854                |
| Amyl alcohol (n-) (1-pentanol) .....                        | 71410                |
| Amyl ether .....  | 693652               |
| Amylamines .....  | 110587               |
| Amylene .....   | 513359               |
| Amylenes, mixed .....                                       |                      |
| Aniline .....   | 62533                |
| Aniline hydrochloride .....                                 | 142041               |
| Anisidine (p-) .....  | 29191524             |
| Anisidine (o-) .....  | 90040                |
| Anisole (methoxy benzene) .....                             | 100663               |
| Anthracene .....  | 120127               |

TABLE 1 TO SUBPART YYY—LIST OF SOCM1 CHEMICALS—Continued

| Chemical name <sup>a</sup>   | CAS No. <sup>b</sup> |
|--|----------------------|
| Anthranilic acid   | 118923               |
| Anthraquinone  | 84651                |
| ar-Methylbenzenediamine  | 25376458             |
| Azobenzene   | 103333               |
| Barium acetate   | 543806               |
| Benzaldehyde   | 100527               |
| Benzamide  | 55210                |
| Benzene  | 71432                |
| Benzenedisulfonic acid   | 98486                |
| Benzenesulfonic acid   | 98113                |
| Benzenesulfonic acid C <sub>10-16</sub> -alkyl derivatives, sodium salts | 68081812             |
| Benzidine  |                      |
| Benzil   | 134816               |
| Benzilic acid  | 76937                |
| Benzoguanamine   |                      |
| Benzoic acid   | 65850                |
| Benzoïn  | 119539               |
| Benzonitrile   | 100470               |
| Benzophenone   | 119619               |
| Benzotrìchloride   | 98077                |
| Benzoyl chloride   | 98884                |
| Benzoyl peroxide   | 94360                |
| Benzyl acetate   | 140114               |
| Benzyl chloride  | 100447               |
| Benzyl alcohol   | 100516               |
| Benzyl dichloride  | 98873                |
| Benzyl benzoate  | 120514               |
| Benzylamine  | 100469               |
| Benzylideneacetone   | 1896624              |
| Biphenyl   | 92524                |
| Bis(Chloromethyl)Ether   | 542881               |
| Bisphenol A  | 80057                |
| Brometone  |                      |
| Bromobenzene   | 108861               |
| Bromoform  | 75252                |
| Bromonaphthalene   | 27497514             |
| Butadiene and butene fractions   |                      |
| Butadiene (1,3-)   | 106990               |
| Butane   | 106978               |
| Butanediol (1,4-)  | 110634               |
| Butanes, mixed   |                      |
| Butenes, mixed   |                      |
| Butyl hydroperoxide (tert-)  | 75912                |
| Butyl acetate (sec-)   | 105464               |
| Butyl chloride (tert-)   | 507200               |
| Butyl alcohol (tert-)  | 75650                |
| Butyl benzoate   | 136607               |
| Butyl mercaptan (n-)   | 109795               |
| Butyl acrylate (n-)  | 141322               |
| Butyl mercaptan (tert-)  | 75661                |
| Butyl methacrylate (n-)  | 97881                |
| Butyl alcohol (sec-)   | 78922                |
| Butyl acetate (tert-)  | 540885               |
| Butyl acetate (n-)   | 123864               |
| Butyl methacrylate (tert-)   |                      |
| Butyl toluene (tert-)  | 98511                |
| Butyl phenol (tert-)   | 88186                |
| Butyl alcohol (n-)   | 71363                |
| Butylamine (t-)  | 75649                |
| Butylamine (s-)  | 13952846             |
| Butylamine (n-)  | 109739               |
| Butylbenzene (tert-)   | 98066                |
| Butylbenzoic acid (p-tert-)  | 98737                |
| Butylbenzyl phthalate  | 85867                |
| Butylene glycol (1,3-)   | 107880               |
| Butylenes (n-)   |                      |
| Butyraldehyde (n-)   | 123728               |
| Butyric acid (n-)  | 107926               |
| Butyric anhydride (n-)   | 106310               |
| Butyrolacetone   | 96480                |
| Butyronitrile  | 109740               |
| Calcium acetate  | 62544                |



TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS—Continued

| Chemical name <sup>a</sup>               | CAS No. <sup>b</sup> |
|--|----------------------|
| Calcium propionate .....                 | 4075814              |
| Caproic acid .....                       | 142621               |
| Caprolactam .....                        | 105602               |
| Carbaryl .....                           | 63252                |
| Carbazole .....                          | 86748                |
| Carbon tetrabromide .....                | 558134               |
| Carbon disulfide .....                   | 75150                |
| Carbon tetrachloride .....               | 56235                |
| Carbon tetrafluoride .....               | 75730                |
| Cellulose acetate .....                  | 9004357              |
| Chloral .....                            | 75876                |
| Chloranil (o-chloranil) .....            | 2435532              |
| Chloranil (p-chloranil) .....            | 118752               |
| Chloroacetic acid .....                  | 79118                |
| Chloroacetophenone (2-) .....            | 532274               |
| Chloroaniline (o-) .....                 | 95512                |
| Chloroaniline (p-) .....                 | 106478               |
| Chloroaniline (m-) .....                 | 108429               |
| Chlorobenzaldehyde (4-) .....            | 104881               |
| Chlorobenzaldehyde (2-) .....            | 89985                |
| Chlorobenzaldehyde (3-) .....            | 587042               |
| Chlorobenzene .....                      | 108907               |
| Chlorobenzoic acid .....                 | 118912               |
|  | 535808               |
|  | 74113                |
| Chlorobenzotrichloride (p-) .....        | 5216251              |
| Chlorobenzotrichloride (o-) .....        | 2136892              |
| Chlorobenzoyl chloride (p-) .....        | 122010               |
| Chlorobenzoyl chloride (o-) .....        | 609654               |
| Chlorodifluoroethane .....               | 25497294             |
| Chlorodifluoromethane .....              | 75456                |
| Chlorofluorocarbons .....                |                      |
| Chloroform .....                         | 67663                |
| Chlorohydrin .....                       |                      |
| Chloronaphthalene .....                  | 25586430             |
| Chloronitrobenzene (o-) .....            | 88733                |
| Chloronitrobenzene (m-) .....            | 121733               |
| Chloronitrobenzene (p-) .....            | 100005               |
| Chlorophenol (o-) .....                  | 95578                |
| Chlorophenol (m-) .....                  | 108430               |
| Chlorophenol (p-) .....                  | 106489               |
| Chlorosulfonic acid .....                | 7790945              |
| Chlorotoluene (m-) .....                 | 108418               |
| Chlorotoluene (o-) .....                 | 95498                |
| Chlorotoluene (p-) .....                 | 106434               |
| Chlorotrifluoroethylene .....            | 79389                |
| Chlorotrifluoromethane .....             | 75729                |
| Choline chloride .....                   | 67481                |
| Chrysene .....                           | 218019               |
| Cinnamic acid .....                      | 140103               |
| Citric acid .....                        | 77929                |
| Cobalt acetate .....                     |                      |
| Copper acetate .....                     | 142712               |
| Cresol and cresylic acid (o-) .....      | 95487                |
| Cresol and cresylic acid (p-) .....      | 106445               |
| Cresol and cresylic acid (m-) .....      | 108394               |
| Cresols and cresylic acids (mixed) ..... | 1319773              |
| Crotonaldehyde .....                     | 4170300              |
| Crotonic acid .....                      | 3724650              |
| Cumene hydroperoxide .....               | 80159                |
| Cumene .....                             | 98828                |
| Cyanamide .....                          | 420042               |
| Cyanoacetic acid .....                   | 372098               |
| Cyanofornamide .....                     |                      |
| Cyanogen chloride .....                  | 506774               |
| Cyanuric acid .....                      | 108805               |
| Cyanuric chloride .....                  | 108770               |
| Cyclohexane, oxidized .....              | 68512152             |
| Cyclohexane .....                        | 110827               |
| Cyclohexanol .....                       | 108930               |
| Cyclohexanone oxime .....                | 100641               |
| Cyclohexanone .....                      | 108941               |

TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS—Continued

| Chemical name <sup>a</sup>                        | CAS No. <sup>b</sup> |
|---|----------------------|
| Cyclohexene                                       | 110838               |
| Cyclohexylamine                                   | 108918               |
| Cyclooctadiene                                    | 29965977             |
| Cyclooctadiene (1,3-)                             | 3806595              |
| Cyclooctadiene (1,5-)                             | 111784               |
| Cyclopentadiene (1,3-)                            |                      |
| Cyclopropane                                      | 75194                |
| Decahydronaphthalene                              | 91178                |
| Decanol   | 112301               |
| Decyl alcohol (1-decanol)                         | 112301               |
| Di-o-tolylguanidine                               | 97392                |
| Di(2-methoxyethyl) phthalate                      |                      |
| Di-n-heptyl-n-nonyl undecyl phthalate             |                      |
| Diacetone alcohol                                 | 123422               |
| Diacetoxy-2-Butene (1,4-)                         |                      |
| Diallyl phthalate                                 | 131179               |
| Diallyl isophthalate                              |                      |
| Diaminobenzoic acids                              | 27576041             |
| Diaminophenol hydrochloride                       | 137097               |
| Dibromomethane                                    | 74953                |
| Dibutanized aromatic concentrate                  |                      |
| Dibutoxyethyl phthalate                           |                      |
| Dichloro-1-butene (3,4-)                          | 760236               |
| Dichloro-2-butene (1,4-)                          | 764410               |
| Dichloro-2-butenes                                |                      |
| Dichloroaniline (mixed isomers)                   | 27134276             |
| Dichlorobenzene (p-)                              | 106467               |
| Dichlorobenzene (m-)                              | 541731               |
| Dichlorobenzene (o-)                              | 95501                |
| Dichlorobenzidine (3,3'-)                         | 91941                |
| Dichlorodifluoromethane                           | 75718                |
| Dichlorodimethylsilane                            | 75785                |
| Dichloroethane (1,2-) (Ethylene dichloride) (EDC) | 107062               |
| Dichloroethyl ether (bis(2-chloroethyl)ether)     | 111444               |
| Dichloroethylene (1,2-)                           | 540590               |
| Dichlorofluoromethane                             | 75434                |
| Dichlorohydrin (a-)                               | 96231                |
| Dichloromethyl ether                              |                      |
| Dichloronitrobenzenes                             |                      |
| Dichloropentanes                                  |                      |
| Dichlorophenol (2,4-)                             | 120832               |
| Dichloropropane (1,1-)                            | 78999                |
| Dichloropropene (1,3-)                            | 542756               |
| Dichloropropene/dichloropropane (mixed)           |                      |
| Dichlorotetrafluoroethane                         | 1320372              |
| Dicyandiamide                                     | 461585               |
| Dicyclohexylamine                                 | 101837               |
| Dicyclopentadiene                                 | 77736                |
| Diethanolamine (2,2'-Iminodiethanol)              | 111422               |
| Diethyl phthalate                                 | 84662                |
| Diethyl sulfate                                   | 64675                |
| Diethylamine                                      | 109897               |
| Diethylaniline (N,N-)                             | 91667                |
| Diethylaniline (2,6-)                             | 579668               |
| Diethylbenzene                                    | 25340174             |
| Diethylene glycol monoethyl ether                 | 111900               |
| Diethylene glycol dimethyl ether                  | 111966               |
| Diethylene glycol                                 | 111466               |
| Diethylene glycol monobutyl ether acetate         | 124174               |
| Diethylene glycol monomethyl ether                | 111773               |
| Diethylene glycol diethyl ether                   | 112367               |
| Diethylene glycol monomethyl ether acetate        | 629389               |
| Diethylene glycol monoethyl ether acetate         | 112152               |
| Diethylene glycol monohexyl ether                 | 112594               |
| Diethylene glycol monobutyl ether                 | 112345               |
| Diethylene glycol dibutyl ether                   | 112732               |
| Difluoroethane (1,1-)                             | 75376                |
| Dihydroxybenzoic acid (Resorcylic acid)           | 27138574             |
| Diisobutylene                                     | 25167708             |
| Diisodecyl phthalate                              | 26761400             |
| Diisononyl phthalate                              | 28553120             |
| Diisooctyl phthalate                              | 27554263             |

TABLE 1 TO SUBPART YYY—LIST OF SOCM1 CHEMICALS—Continued

| Chemical name <sup>a</sup>                  | CAS No. <sup>b</sup> |
|---|----------------------|
| Diisopropylamine                            | 108189               |
| Diketene (4-methylene-2-oxetanone)          | 674828               |
| Dimethyl sulfate                            | 77781                |
| Dimethyl ether                              | 115106               |
| Dimethyl sulfide                            | 75183                |
| Dimethyl phthalate                          | 131113               |
| Dimethyl sulfoxide                          | 67685                |
| Dimethyl terephthalate                      | 120616               |
| Dimethylacetamide (N,N-)                    | 127195               |
| Dimethylamine                               | 124403               |
| Dimethylaminoethanol (2-)                   | 108010               |
| Dimethylaniline (N,N)                       | 121697               |
| Dimethylbenzidine (3,3'-)                   | 119937               |
| Dimethylformamide (N,N-)                    | 68122                |
| Dimethylhydrazine (1,1-)                    | 57147                |
| Dimethylphenol (2,5-) Xylenol (2, 5-)       | 95874                |
| Dimethylphenol (2,6-) Xylenol (2, 6-)       | 576261               |
| Dimethylphenol (3,5-) Xylenol (3, 5-)       | 108689               |
| Dimethylphenol (2,4-) Xylenol (2, 4-)       | 105679               |
| Dimethylphenol (2,3-) Xylenol (2, 3-)       | 526750               |
| Dimethylphenol (3,4-) Xylenol (3, 4-)       | 95658                |
| Dinitrobenzenes (NOS) <sup>c</sup>          | 25154545             |
| Dinitrobenzoic acid (3,5-)                  | 99343                |
| Dinitrophenol (2,4-)                        | 51285                |
| Dinitrotoluene (3,4-)                       | 610399               |
| Dinitrotoluene (2,6-)                       | 606202               |
| Dinitrotoluene (2,3-)                       | 602017               |
| Dinitrotoluene (2,4-)                       | 121142               |
| Diocyl phthalate                            | 117817               |
| Dioxane (1,4-) (1,4-Diethyleneoxide)        | 123911               |
| Dioxolane (1,3-)                            | 646060               |
| Diphenyl oxide                              | 101848               |
| Diphenyl thiourea (N,N'-)                   | 102089               |
| Diphenyl methane                            | 101815               |
| Diphenylamine                               | 122394               |
| Dipropylene glycol                          | 110985               |
| Dodecandedioic acid                         | 693232               |
| Dodecene (branched)                         | 112414               |
| Dodecene (n-)                               | 25378227             |
| Dodecyl phenol (branched)                   | 121158585            |
| Dodecyl benzene (branched)                  | 123013               |
| Dodecylaniline                              | 28675174             |
| Dodecylbenzene sulfonic acid                | 27176870             |
| Dodecylbenzene, nonlinear                   |                      |
| Dodecylbenzene (n-)                         | 121013               |
| Dodecylbenzene sulfonic acid, sodium salt   | 25155300             |
| Dodecylmercaptan (branched)                 | 25103586             |
| Dodecylphenol                               | 27193868             |
| Epichlorohydrin (1-chloro-2,3-epoxypropane) | 106898               |
| Ethane                                      | 74840                |
| Ethanol                                     | 64175                |
| Ethanolamine                                | 141435               |
| Ethyl ether                                 | 60297                |
| Ethyl oxalate                               | 95921                |
| Ethyl orthoformate                          | 122510               |
| Ethyl acetate                               | 141786               |
| Ethyl bromide                               | 74964                |
| Ethyl chloride (Chloroethane)               | 75003                |
| Ethyl cyanide                               | 107120               |
| Ethyl acrylate                              | 140885               |
| Ethyl sodium oxalacetate                    | 41892711             |
| Ethyl acetoacetate                          | 141979               |
| Ethyl chloroacetate                         | 105395               |
| Ethyl mercaptan (ethanethiol)               | 75081                |
| Ethylamine                                  | 75047                |
| Ethylaniline (o-)                           | 578541               |
| Ethylaniline (n-)                           | 103695               |
| Ethylbenzene                                | 100414               |
| Ethylcellulose                              | 9004573              |
| Ethylcyanoacetate                           | 105566               |
| Ethylene glycol dibutyl ether               | 112481               |
| Ethylene dibromide (Dibromoethane)          | 106934               |

TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS—Continued

| Chemical name <sup>a</sup>                         | CAS No. <sup>b</sup> |
|--|----------------------|
| Ethylene oxide                                     | 75218                |
| Ethylene glycol monoethyl ether                    | 110805               |
| Ethylene glycol monoethyl ether acetate            | 111159               |
| Ethylene glycol monomethyl ether                   | 109864               |
| Ethylene dichloride                                | 107062               |
| Ethylene glycol monobutyl ether acetate            | 112072               |
| Ethylene glycol dimethyl ether                     | 110714               |
| Ethylene   | 74851                |
| Ethylene glycol monophenyl ether                   | 122996               |
| Ethylene glycol monoacetate                        | 542596               |
| Ethylene carbonate                                 | 96491                |
| Ethylene glycol monoethyl ether                    |                      |
| Ethylene glycol diacetate                          | 111557               |
| Ethylene glycol diethyl ether (1,2-diethoxyethane) | 629141               |
| Ethylene glycol monopropyl ether                   | 2807309              |
| Ethylene glycol monohexyl ether                    | 112254               |
| Ethylene glycol monomethyl ether acetate           | 110496               |
| Ethylene glycol monobutyl ether                    | 111762               |
| Ethylene chlorohydrin                              | 107073               |
| Ethylene glycol                                    | 107211               |
| Ethylenediamine                                    | 107153               |
| Ethylenediamine tetraacetic acid                   | 60004                |
| Ethylenimine (Aziridine)                           | 151564               |
| Ethylhexanoic acid (2-)                            | 149575               |
| Ethylhexyl succinate (2-)                          |                      |
| Ethylhexyl acrylate (2-isomer)                     | 103117               |
| Ethylmethylbenzene                                 | 25550145             |
| Fluoranthene                                       | 206440               |
| Formaldehyde                                       | 50000                |
| Formamide  | 75127                |
| Formic acid  | 64186                |
| Fumaric acid                                       | 110178               |
| Furfural (2-furan carboxaldehyde)                  | 98011                |
| Glutaraldehyde                                     | 111308               |
| Glyceraldehyde                                     | 367475               |
| Glycerol dichlorohydrin                            | 26545737             |
| Glycerol tri(polyoxypropylene)ether                | 25791962             |
| Glycerol   | 56815                |
| Glycidol   | 556525               |
| Glycine  | 56406                |
| Glycol ethers                                      |                      |
| Glyoxal (ethane dial)                              | 107222               |
| Guanidine  |                      |
| Guanidine nitrate                                  | 506934               |
| Heptenes   |                      |
| Hexachlorobenzene                                  | 118741               |
| Hexachlorobutadiene                                | 87683                |
| Hexachlorocyclopentadiene                          | 77474                |
| Hexachloroethane                                   | 67721                |
| Hexadecyl chloride                                 |                      |
| Hexadecyl alcohol (1-hexadecanol)                  | 36653824             |
| Hexadiene (1,4-)                                   | 592450               |
| Hexamethylene glycol                               | 629118               |
| Hexamethylene diamine adipate                      | 3323533              |
| Hexamethylenediamine                               | 124094               |
| Hexamethylenetetramine                             | 100970               |
| Hexane   | 110543               |
| Hexanetriol (1,2,6-)                               | 106694               |
| Hexyl alcohol                                      | 111273               |
| Hexylene glycol                                    | 107415               |
| Higher glycols                                     |                      |
| Hydrogen cyanide                                   | 74908                |
| Hydroquinone                                       | 123319               |
| Hydroxyadipaldehyde                                | 141311               |
| Hydroxybenzoic acid (p-)                           | 99967                |
| Iminodiethanol (2,2-) (diethanolamine)             | 111422               |
| Isoamyl alcohol                                    | 123513               |
| Isoamyl chloride (mixed)                           |                      |
| Isoamylene   | 26760645             |
| Isobutane  | 75285                |
| Isobutanol   | 78831                |
| Isobutyl methacrylate                              | 97869                |

TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS—Continued

| Chemical name <sup>a</sup>                                    | CAS No. <sup>b</sup> |
|---|----------------------|
| Isobutyl acetate  | 110190               |
| Isobutyl acrylate   | 106638               |
| Isobutyl vinyl ether  | 109535               |
| Isobutyl alcohol  | 78831                |
| Isobutylene   | 115117               |
| Isobutyraldehyde (2-methyl-propanal)                          | 78842                |
| Isobutyric acid   | 79312                |
| Isodecanol  | 25339177             |
| Isohexyldecyl alcohol   |                      |
| Isononyl alcohol  |                      |
| Isocetyl alcohol  | 26952216             |
| Isopentane  | 78784                |
| Isophorone nitrile  |                      |
| Isophorone  | 78591                |
| Isophthalic acid  | 121915               |
| Isoprene  | 78795                |
| Isopropanol   | 67630                |
| Isopropyl acetate   | 108214               |
| Isopropyl ether   | 108203               |
| Isopropyl chloride  | 75296                |
| Isopropylamine  | 75310                |
| Isopropylphenol   | 25168063             |
| Ketene  | 463514               |
| Lactic acid   | 79334                |
| Lauryl dimethylamine oxide                                    |                      |
| Lead subacetate   | 1335326              |
| Lead phthalate  |                      |
| Lead acetate  | 6080564              |
| Linear alcohols, ethoxylated and sulfated, sodium salt, mixed |                      |
| Linear alcohols, ethoxylated, mixed                           |                      |
| Linear alkyl sulfonate  |                      |
| Linear alcohols, sulfated, sodium salt, mixed                 |                      |
| Magnesium acetate   | 142723               |
| Maleic anhydride  | 108316               |
| Maleic hydrazide  | 123331               |
| Maleic acid   | 110167               |
| Malic acid  | 6915157              |
| Manganese acetate   | 638380               |
| Melamine (1,3,5-triazine-2,4,6-triamine)                      | 108781               |
| Mercuric acetate  | 1600277              |
| Mesityl oxide   | 141797               |
| Metanilic acid  | 121471               |
| Methacrylic acid  | 79414                |
| Methacrylonitrile   | 126987               |
| Methallyl chloride  | 563473               |
| Methallyl alcohol   | 513428               |
| Methane   | 74828                |
| Methanol  | 67561                |
| Methionine  | 63683                |
| Methyl mercaptan  | 74931                |
| Methyl iodide   | 74884                |
| Methyl ethyl ketone (2-butanone)                              | 78933                |
| Methyl isobutyl carbinol                                      | 108112               |
| Methyl acetate  | 79209                |
| Methyl chloride (Chloromethane)                               | 74873                |
| Methyl salicylate   | 119368               |
| Methyl acetoacetate   | 105453               |
| Methyl bromide (Bromomethane)                                 | 74839                |
| Methyl formate  | 107313               |
| Methyl phenyl carbinol  | 98851                |
| Methyl methacrylate   | 80626                |
| Methyl tert-butyl ether                                       | 1634044              |
| Methyl isocyanate   | 624839               |
| Methyl butynol  | 37365712             |
| Methyl hydrazine  | 60344                |
| Methyl isobutyl ketone (Hexone)                               | 108101               |
| Methyl acrylate   | 96333                |
| Methyl butenols   |                      |
| Methyl anthranilate   | 134203               |
| Methylamine   | 74895                |
| Methylaniline (N-)  | 100618               |
| Methylbutanol (2-)  | 137326               |

TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS—Continued

| Chemical name <sup>a</sup>                    | CAS No. <sup>b</sup> |
|---|----------------------|
| Methylcyclohexane                             | 108872               |
| Methylcyclohexanol                            | 25639423             |
| Methylcyclohexanone                           | 1331222              |
| Methylene chloride (Dichloromethane)          | 75092                |
| Methylene dianiline (4,4')                    | 101779               |
| Methylene diphenyl diisocyanate (4,4'-) (MDI) | 101688               |
| Methylenes (a-)                               | 79696                |
| Methylnaphthalene (2-)                        | 91576                |
| Methylnaphthalene (1-)                        | 90120                |
| Methylpentane (2-)                            | 107835               |
| Methylpentynol                                | 77758                |
| Methylstyrene (a-)                            | 98839                |
| Monomethylhydrazine                           |                      |
| Morpholine                                    | 110918               |
| n-Heptane                                     | 142825               |
| n-Propanol                                    | 71238                |
| N-Vinyl-2-pyrrolidone                         |                      |
| Naphthalene sulfonic acid (a-)                | 85472                |
| Naphthalene                                   | 91203                |
| Naphthalene sulfonic acid (b-)                | 120183               |
| Naphthenic acids                              |                      |
| Naphthol (a-)                                 | 90153                |
| Naphthol (b-)                                 | 135193               |
| Naphtholsulfonic acid (1-)                    | 567180               |
| Naphthylamine sulfonic acid (1,4-)            | 84866                |
| Naphthylamine (1-)                            | 134327               |
| Naphthylamine (2-)                            | 91598                |
| Naphthylamine sulfonic acid (2,1-)            | 81163                |
| Neohexane                                     | 75832                |
| Neopentanoic acid                             | 75989                |
| Neopentyl glycol                              | 126307               |
| Nickel formate                                |                      |
| Nitriloacetic acid                            |                      |
| Nitrilotriacetic acid                         | 139139               |
| Nitroaniline (m-)                             | 99092                |
| Nitroaniline (p-)                             | 100016               |
| Nitroaniline (o-)                             | 88744                |
| Nitroanisole (p-)                             | 100174               |
| Nitroanisole (o-)                             | 91236                |
| Nitrobenzene                                  | 98953                |
| Nitrobenzoic acid (m-)                        | 121926               |
| Nitrobenzoic acid (o-)                        | 552169               |
| Nitrobenzoic acid (p-)                        | 62237                |
| Nitrobenzoyl chloride (p-)                    |                      |
| Nitroethane                                   | 79243                |
| Nitroguanidine                                | 556887               |
| Nitromethane                                  | 75525                |
| Nitronaphthalene (1-)                         | 86577                |
| Nitrophenol (p-)                              | 100027               |
| Nitrophenol (o-)                              | 88755                |
| Nitropropane (1-)                             | 25322014             |
| Nitropropane (2-)                             | 79469                |
| Nitrotoluene (p-)                             | 99990                |
| Nitrotoluene (o-)                             | 88722                |
| Nitrotoluene (m-)                             | 99081                |
| Nitrotoluene (all isomers)                    | 1321126              |
| Nitroxylene                                   | 25168041             |
| Nonene  | 27215958             |
| Nonyl alcohol                                 | 1430808              |
| Nonylbenzene (branched)                       | 1081772              |
| Nonylphenol                                   | 25154523             |
| Nonylphenol, ethoxylated                      | 9016459              |
| Nonylphenol (branched)                        | 25154523             |
| Octane  | 111659               |
| Octene-1                                      | 111660               |
| Octylamine (tert-)                            | 107459               |
| Octylphenol                                   | 27193288             |
| Oil-soluble petroleum sulfonate sodium salt   |                      |
| Oil-soluble petroleum sulfonate calcium salt  |                      |
| Oxalic acid                                   | 144627               |
| Oxamide                                       | 471465               |
| Oxo chemicals                                 |                      |

TABLE 1 TO SUBPART YYY—LIST OF SOCM I CHEMICALS—Continued

| Chemical name <sup>a</sup>            | CAS No. <sup>b</sup> |
|---------------------------------------|----------------------|
| p-tert-Butyl toluene                  | 98511                |
| Paraformaldehyde                      | 30525894             |
| Paraldehyde                           | 123637               |
| Pentachlorophenol                     | 87865                |
| Pentaerythritol tetranitrate          |                      |
| Pentane                               | 109660               |
| Pentanethiol                          | 115775               |
| Pentanol (3-)                         | 584021               |
| Pentanol (2-)                         | 6032297              |
| Pentene (1-)                          | 109671               |
| Pentene (2-)                          | 109682               |
| Peracetic acid                        | 79210                |
| Perchloromethyl mercaptan             | 594423               |
| Phenacetin                            | 62442                |
| Phenanthrene                          | 85018                |
| Phenetidine (p-)                      | 156434               |
| Phenetidine (o-)                      | 94702                |
| Phenol                                | 108952               |
| Phenolphthalein                       | 77098                |
| Phenolsulfonic acids (all isomers)    | 1333397              |
| Phenyl anthranilic acid (all isomers) | 91407                |
| Phenylenediamine (m-)                 | 108452               |
| Phenylenediamine (p-)                 | 106503               |
| Phenylenediamine (o-)                 | 95545                |
| Phenylmethylpyrazolone                |                      |
| Phenylpropane                         | 103651               |
| Phloroglucinol (1,3,5-benzenetriol)   | 108736               |
| Phosgene                              | 75445                |
| Phthalic acid                         | 88993                |
| Phthalic anhydride                    | 85449                |
| Phthalimide                           | 85416                |
| Phthalonitrile                        | 91156                |
| Picoline (b-)                         | 108996               |
| Picoline (a-)                         |                      |
| Picramic acid                         |                      |
| Picric acid                           | 88891                |
| Piperazine                            | 110850               |
| Piperidine                            | 110894               |
| Piperylene                            | 504609               |
| Polybutenes                           | 9003296              |
| Polyethylene glycol                   | 25322683             |
| Polypropylene glycol                  | 25322694             |
| Potassium acetate                     | 127082               |
| Propane                               | 74986                |
| Propiolactone (beta-) (2-Oxetanone)   | 57578                |
| Propionaldehyde                       | 123386               |
| Propionic acid                        | 79094                |
| Propyl acetate (n-)                   | 109604               |
| Propyl chloride                       | 540545               |
| Propyl alcohol (n-)                   | 71238                |
| Propylamine                           | 107108               |
| Propylene chlorohydrin                | 127004               |
| Propylene glycol                      | 57556                |
| Propylene                             | 115071               |
| Propylene oxide                       | 75569                |
| Propylene carbonate                   | 108327               |
| Propylene glycol monomethyl ether     | 107982               |
| 1,2-dichloropropane                   | 78875                |
| Pseudocumene                          | 95636                |
| Pseudocumidine                        |                      |
| Pyrene                                | 129000               |
| Pyridine                              | 110861               |
| Pyrrolidone (2-)                      | 616455               |
| Quinone                               | 106514               |
| Resorcinol (1,3-benzenediol)          | 108463               |
| Salicylic acid                        | 69727                |
| Sebacic acid                          | 111206               |
| Sodium benzoate                       | 532321               |
| Sodium phenate                        | 139026               |
| Sodium acetate                        | 127093               |
| Sodium formate                        | 141537               |
| Sodium methoxide                      | 124414               |



TABLE 1 TO SUBPART YYY—LIST OF SOCM1 CHEMICALS—Continued

| Chemical name <sup>a</sup>              | CAS No. <sup>b</sup> |
|---|----------------------|
| Sodium cyanide                          | 143339               |
| Sodium propionate                       | 137406               |
| Sodium chloroacetate                    | 3926623              |
| Sodium carboxymethyl cellulose          | 9004324              |
| Sodium oxalate                          | 62760                |
| Sodium dodecyl benzene sulfonate        |                      |
| Sorbic acid                             | 110441               |
| Sorbitol (D-Glucitol)                   | 50704                |
| Stilbene                                | 588590               |
| Styrene                                 | 100425               |
| Succinic acid                           | 110156               |
| Succinonitrile                          | 110612               |
| Sulfanilic acid                         | 121573               |
| Sulfolane                               | 126330               |
| Synthesis gas                           |                      |
| Tannic acid                             | 1401554              |
| Tartaric acid                           | 526830               |
| Terephthalic acid                       | 100210               |
| Terephthaloyl chloride                  | 100209               |
| Tetra (methyl-ethyl) lead               |                      |
| Tetrabromophthalic anhydride            | 632791               |
| Tetrachlorobenzene (1,2,3,5-)           |                      |
| Tetrachlorobenzene (1,2,4,5-)           | 95943                |
| Tetrachloroethane (1,1,2,2-)            | 79345                |
| Tetrachloroethylene (Perchloroethylene) | 127184               |
| Tetrachlorophthalic anhydride           | 117088               |
| Tetraethyl lead                         | 78002                |
| Tetraethylene glycol                    | 112607               |
| Tetraethylenepentamine                  | 112572               |
| Tetrafluoroethylene                     |                      |
| Tetrahydrofuran                         | 109999               |
| Tetrahydronaphthalene                   | 119642               |
| Tetrahydrophthalic anhydride            | 85438                |
| Tetramethylenediamine                   | 110601               |
| Tetramethylethylenediamine              | 110189               |
| Tetramethyllead                         | 75741                |
| Thiourea                                | 62566                |
| Tolidines                               |                      |
| Toluene sulfonic acids                  | 104154               |
| Toluene diisocyanate (2,4-)             | 584849               |
| Toluene                                 | 108883               |
| Toluene diamine (2,4-)                  | 95807                |
| Toluene diisocyanates (mixture)         | 26471625             |
| Toluene sulfonamides (o- and p-)        | 1333079              |
| Toluenesulfonyl chloride                | 98599                |
| Toluidine (o-)                          | 95534                |
| Trichloroacetic acid                    | 76039                |
| Trichloroaniline (2,4,6-)               | 634935               |
| Trichlorobenzene (1,2,4-)               | 120821               |
| Trichlorobenzene (1,2,3-)               | 87616                |
| Trichlorobenzene (1,3,5-)               | 108703               |
| Trichloroethane (1,1,2-)                | 79005                |
| Trichloroethane (1,1,1-)                | 71556                |
| Trichloroethylene                       | 79016                |
| Trichlorofluoromethane                  | 75694                |
| Trichlorophenol (2,4,5-)                | 95954                |
| Trichloropropane (1,2,3-)               | 96184                |
| Tricresyl phosphate                     | 1330785              |
| Tridecyl alcohol                        | 112709               |
| Tridecyl mercaptan                      |                      |
| Triethanolamine                         | 102716               |
| Triethylamine                           | 121448               |
| Triethylene glycol monoethyl ether      | 112505               |
| Triethylene glycol                      | 112276               |
| Triethylene glycol dimethyl ether       | 112492               |
| Triethylene glycol monomethyl ether     | 112356               |
| Triisobutylene                          | 7756947              |
| Trimellitic anhydride                   | 552307               |
| Trimethyl-1,3-pentanediol (2,2,4-)      | 144194               |
| Trimethyl-1-pentanol (2,4,4-)           | 16325636             |
| Trimethylamine                          | 75503                |
| Trimethylcyclohexanol                   | 933482               |

TABLE 1 TO SUBPART YYY—LIST OF SOCMI CHEMICALS—Continued

| Chemical name <sup>a</sup>                       | CAS No. <sup>b</sup> |
|--|----------------------|
| Trimethylcyclohexanone .....                     | 2408379              |
| Trimethylcyclohexylamine .....                   | 34216347             |
| Trimethylolpropane .....                         | 77996                |
| Trimethylpentane (2,2,4-) .....                  | 540841               |
| Tripropylene glycol .....                        | 24800440             |
| Urea .....                                       | 57136                |
| Vinyl chloride (Chloroethylene) .....            | 75014                |
| Vinyl acetate .....                              | 108054               |
| Vinyl toluene .....                              | 25013154             |
| Vinyl (N-)-pyrrolidone (2-) .....                | 88120                |
| Vinylcyclohexene (4-) .....                      | 100403               |
| Vinylidene chloride (1,1-dichloroethylene) ..... | 75354                |
| Vinylpyridine (2-) .....                         | 100696               |
| Xanthates .....                                  | 140896               |
| Xylene sulfonic acid .....                       | 25321419             |
| Xylene (m-) .....                                | 108383               |
| Xylene (o-) .....                                | 95476                |
| Xylene (p-) .....                                | 106423               |
| Xylenes (NOS) <sup>c</sup> .....                 | 1330207              |
| Xylenols (Mixed) .....                           | 1300716              |
| Xylidene (dimethylbenzene diamine) .....         | 1300738              |
| Xylidene (2,3-) .....                            | 1300738              |
| Xylidene (2,6-) .....                            | 1300738              |
| Xylidene (2,5-) .....                            | 1300738              |
| Xylidene (3,5-) .....                            | 1300738              |
| Xylidene (2,4-) .....                            | 1300738              |
| Xylidene (3,4-) .....                            | 1300738              |
| Zinc acetate .....                               | 5970456              |

<sup>a</sup> Isomer means all structural arrangements for the same number of atoms of each element and does not mean salts, esters, or derivatives.

<sup>b</sup> CAS Number = Chemical Abstract Service number.

TABLE 2 TO SUBPART YYY—APPLICABILITY OF 40 CFR PART 60 GENERAL PROVISIONS TO SUBPART YYY

| Reference                    | Applies to subpart YYY | Subject/comment  |
|------------------------------|------------------------|--|
| 60.1 .....                   | Yes .....              | Applicability.   |
| 60.2 .....                   | Yes .....              | Definitions. If a term is defined in both the General Provisions and subpart YYY, the definition in YYY shall override the definition in the General provisions.   |
| 60.3 .....                   | Yes .....              | Units and abbreviations.   |
| 60.4 .....                   | Yes .....              | Address.   |
| 60.5 .....                   | Yes .....              | Determination of construction or modification.   |
| 60.6 .....                   | Yes .....              | Review of plans.   |
| 60.7(a)(1) .....             | Yes .....              | Submit a notification of the date construction or reconstruction commences.  |
| 60.7(a)(2) .....             | Yes .....              | Submit a notification of anticipated date of initial startup.  |
| 60.7(a)(3) .....             | Yes .....              | Submit a notification of actual date of initial startup.   |
| 60.7(a)(4) .....             | Yes .....              | Submit a notification of any physical or operational change to an existing facility which increases the emission rate of any air pollutant.  |
| 60.7(a)(5) .....             | No .....               | Continuous monitoring requirements and associated reporting and recordkeeping are specified in §§60.781, 60.784, and 60.785 of subpart YYY.  |
| 60.7(a)(6)–160.7(a)(7) ..... | No .....               | Subpart YYY is not an opacity standard.  |
| 60.7(b) .....                | No .....               | §60.785(f)(7) of subpart YYY specify which records to maintain to document periods of startup, shutdown, or malfunction; and periods when a continuous monitoring system is inoperative.   |
| 60.7(c), (d), and (e) .....  | No .....               | The semiannual report required in §60.784(d) includes reports of all excursions and all periods when monitoring parameters are above the maximum or below the minimum established value. §60.784(e) specifies semiannual reporting for treatment processes. Excess emissions are discussed and defined in §60.787(a)(3) and must be recorded in the startup, shutdown, and malfunction plan as specified in §60.787. |
| 60.7(f) .....                | No .....               | §60.785 specifies data retention and the types of records that must be maintained.   |
| 60.7(g)–(h) .....            | Yes .....              |  |
| 60.8 .....                   | No .....               | The performance testing requirements in 40 CFR 63.7 apply to subpart YYY as specified in Table 2A of this subpart.   |
| 60.9 .....                   | Yes .....              | Availability of information.   |
| 60.10 .....                  | Yes .....              | State authority.   |
| 60.11 .....                  | No .....               | Subpart YYY is not an opacity standard. Operation and maintenance requirements are specified throughout subpart YYY.   |
| 60.12 .....                  | Yes .....              | Circumvention.   |
| 60.13(a) .....               | No .....               | Continuous monitoring requirements are specified in §60.781.   |
| 60.13(b) .....               | Yes .....              | Except use 40 CFR part 63.11 in place of the §60.8 reference to performance tests.   |

TABLE 2 TO SUBPART YYY—APPLICABILITY OF 40 CFR PART 60 GENERAL PROVISIONS TO SUBPART YYY—Continued

| Reference                  | Applies to subpart YYY | Subject/comment  |
|----------------------------|------------------------|--|
| 60.13(c)–(d)               | No                     | Opacity and continuous emission monitoring not required in subpart YYY.  |
| 60.13(e)                   | No                     | Monitoring frequency is specified in § 60.781.   |
| 60.13(f)                   | Yes                    | Except § 60.781(g), which pertains to installation, calibration, and maintenance of monitoring equipment, applies, also. |
| 60.13(g)                   | No                     | The locations to install CMS are specified in § 60.781.  |
| 60.13(h)                   | No                     | Provisions explaining how to calculate continuous parameter monitoring values are specified in § 60.785.                 |
| 60.13(i)(1)–(i)(7), (i)(9) | Yes                    | Alternatives to monitoring methods or procedures must be approved by the Administrator.                                  |
| 60.13(i)(8)                | No                     | Opacity monitoring not required in subpart YYY.  |
| 60.13(j)                   | No                     | Continuous emission monitoring not required in subpart YYY.  |
| 60.14(a)                   | No                     | A modification is determined as specified in § 60.772(b).  |
| 60.14(b)                   | No                     | VOC emissions are determined as specified in § 60.772(f).  |
| 60.14(c)                   | No                     |  |
| 60.14(d)                   | No                     | Reserved.  |
| 60.14(e)                   | No                     | Types of physical and operational changes that are not modifications are specified in § 60.772(c).                       |
| 60.14(f)–(g)               | Yes                    |  |
| 60.14(h)–60.14(l)          | Yes                    | List of changes and projects that are exempt from modification provisions.   |
| 60.15                      | Yes                    | Except 60.772(g) specifies additional requirements for “fixed capital cost of the new components.”                       |
| 60.16                      | Yes                    | Prioritized major source categories.   |
| 60.17                      | Yes                    | Incorporations by reference.   |
| 60.18                      | No                     | Control device requirements are specified in 40 CFR part 63.11, as specified in Table 2A of this subpart.                |
| 60.19                      | Yes                    | General notification and reporting requirements.   |

TABLE 2A TO SUBPART YYY—APPLICABILITY OF 40 CFR PART 63 GENERAL PROVISIONS TO SUBPART YYY

| Reference          | Applies to subpart YYY | Subject/comment  |
|--------------------|------------------------|--|
| 63.6(e)(3)(i)      | Yes                    | The startup, shutdown, malfunction plan may include written procedures that identify conditions that justify a delay of repair.  |
| 63.6(e)(3)(i)(B)   | Yes                    |  |
| 63.6(e)(3)(i)(C)   | Yes                    |  |
| 63.6(e)(3)(ii)     | Yes                    |  |
| 63.6(e)(3)(v)      | Yes                    |  |
| 63.6(e)(3)(vi)     | Yes                    |  |
| 63.6(e)(3)(vii)    | Yes                    |  |
| 63.6(e)(3)(vii)(A) | Yes                    |  |
| 63.6(e)(3)(vii)(B) | Yes                    | Except the plan shall provide for operation in compliance with § 60.787(a)(3).   |
| 63.6(e)(3)(vii)(C) | Yes                    |  |
| 63.6(e)(3)(viii)   | Yes                    |  |
| 63.7(a)(3)         | Yes                    |  |
| 63.7(d)            | Yes                    |  |
| 63.7(e)(1)–(e)(2)  | Yes                    | Except § 60.783(a)(8), representative process unit operating conditions, and (a)(9), representative treatment process or control device operating conditions, also address this issue. |
| 63.7(e)(4)         | Yes                    |  |
| 63.7(h)(1)–(h)(2)  | Yes                    |  |
| 63.7(h)(5)         | Yes                    |  |
| 63.9(a)(4)         | Yes                    |  |
| 63.9(b)(5)         | Yes                    |  |
| 63.11              | Yes                    |  |

TABLE 3 TO SUBPART YYY—CONTROL REQUIREMENT OPTIONS FOR WASTEWATER TANKS, SURFACE IMPOUNDMENT, CONTAINERS, INDIVIDUAL DRAIN SYSTEMS, AND OIL-WATER SEPARATORS

| Unit operation       | Part 63, subpart G (HON) | Part 63 (standard-standards)        | Part 60, subpart QQQ (petroleum refinery) | Part 264, subpart CC (RCRA CC) | Part 265, subpart CC (RCRA CC) | Part 61, subpart FF (benzene waste) |
|----------------------|--------------------------|-------------------------------------|---|--------------------------------|--------------------------------|-------------------------------------|
| Wastewater Tanks     | § 63.133                 | .....                               | .....                                     | § 264.1084                     | § 265.1085                     | § 61.343                            |
| Surface Impoundments | § 63.134                 | §§ 63.942 and 63.943 of Subpart QQ. | .....                                     | § 264.1085                     | § 265.1086                     | § 61.344                            |

TABLE 3 TO SUBPART YYY—CONTROL REQUIREMENT OPTIONS FOR WASTEWATER TANKS, SURFACE IMPOUNDMENT, CONTAINERS, INDIVIDUAL DRAIN SYSTEMS, AND OIL-WATER SEPARATORS—Continued

| Unit operation             | Part 63, subpart G (HON) | Part 63 (standard-standards)                    | Part 60, subpart QQQ (petroleum refinery) | Part 264, subpart CC (RCRA CC) | Part 265, subpart CC (RCRA CC) | Part 61, subpart FF (benzene waste) |
|----------------------------|--------------------------|---|---|--------------------------------|--------------------------------|-------------------------------------|
| Containers .....           | § 63.135                 | §§ 63.922 and 63.923 of Subpart PP.             | .....                                     | § 264.1086                     | § 265.1087                     | § 61.345                            |
| Individual Drain System    | § 63.136                 | § 63.962 of Subpart RR.                         | § 60.693-1                                | .....                          | .....                          | § 61.346                            |
| Oil-water separators ..... | § 63.137                 | §§ 63.1042, 63.1043, and 63.1044 of Subpart VV. | .....                                     | .....                          | .....                          | § 61.347                            |

TABLE 4 TO SUBPART YYY—WASTEWATER TANKS REQUIRING CONTROLS AND CONTROL REQUIREMENTS

| Capacity (m <sup>3</sup> ) | Vapor pressure (kPa) | Control requirements         |
|----------------------------|----------------------|------------------------------|
| <75 .....                  | .....                | FR, IFR, EFR, or CVS and CD. |
| ≥75 and <151 .....         | <13.1                | FR, IFR, EFR, or CVS and CD. |
| .....                      | ≥13.1                | IFR, EFR, or CVS and CD.     |
| ≥151 .....                 | <5.2                 | FR, IFR, EFR, or CVS and CD. |
| .....                      | ≥5.2                 | IFR, EFR, or CVS and CD.     |

FR means fixed roof requirements in the applicable compliance option.  
 IFR means internal floating roof requirements in the applicable compliance option.  
 EFR means external floating roof requirements in the applicable compliance option.  
 CVS and CD means closed vent system routed to a control device requirements in the applicable compliance option.

TABLE 5 TO SUBPART YYY—COMPLIANCE OPTIONS FOR WASTEWATER TANKS, SURFACE IMPOUNDMENTS, CONTAINERS, INDIVIDUAL DRAIN SYSTEMS, AND OIL-WATER SEPARATORS

| Headings within § 60.774 | Part 63, subpart G (HON)  | Part 63 (standard-standards)   | Part 60, subpart QQQ (petroleum refinery)   | Part 264, subpart CC (RCRA CC)   | Part 265, subpart CC (RCRA CC)  | Part 61, subpart FF (benzene waste)  |
|--------------------------|---|--|---|--|---|--|
| Control Requirements.    | If WMU subject to YYY or HON, then comply with HON control requirements.  | Comply with subpart QQ, PP, RR, or VV control requirements, as applicable. | If WMU subject to YYY or Petroleum Refinery for individual drain systems, then comply with Petroleum Refinery control requirements.   | If WMU subject to YYY and RCRA, part 264, then comply with RCRA, part 264 control requirements.  | If WMU subject to YYY and RCRA, part 265, then comply with RCRA, part 265 control requirements.   | If WMU subject to YYY and Benzene Waste, then comply with Benzene Waste control requirements.  |
| Monitoring .....         | If WMU subject to HON, can comply with either HON or YYY provisions; if WMU not subject to HON, comply with YYY provisions. | Comply with standards provisions or YYY provisions.                        | If WMU subject to Petroleum Refinery, then can comply with either Petroleum Refinery or YYY provisions; if WMU not subject to Petroleum Refinery, comply with YYY provisions. | If WMU subject to part RCRA, part 264, then can comply with either RCRA, part 264 or YYY provisions; if WMU not subject to RCRA, part 264, comply with YYY provisions. | If WMU subject to RCRA, part 265, then can comply with either RCRA, part 265 or YYY provisions; if WMU not subject to RCRA, part 265, comply with YYY provisions.   | If WMU subject to Benzene Waste, then can comply with either Benzene Waste or YYY provisions; if WMU not subject to Benzene Waste, comply with YYY provisions. |
| Reporting .....          | If WMU subject to HON, can comply with either HON or YYY provisions; if WMU not subject to HON, comply with YYY provisions. | Comply with standards provisions or YYY provisions*.                       | If WMU subject to Petroleum Refinery, then can comply with either Petroleum Refinery or YYY provisions; if WMU not subject to Petroleum Refinery, comply with YYY provisions. | If WMU subject to RCRA, part 264, then can comply with either RCRA, part 264 or YYY provisions; if WMU not subject to RCRA, part 264, comply with YYY provisions.      | If WMU subject to RCRA, part 265, then can comply with either RCRA, part 265 or YYY provisions; if WMU not subject to RCRA, part 265, comply with YYY provisions**. | If WMU subject to Benzene Waste, then can comply with either Benzene Waste or YYY provisions; if WMU not subject to Benzene Waste, comply with YYY provisions. |

TABLE 5 TO SUBPART YYY—COMPLIANCE OPTIONS FOR WASTEWATER TANKS, SURFACE IMPOUNDMENTS, CONTAINERS, INDIVIDUAL DRAIN SYSTEMS, AND OIL-WATER SEPARATORS—Continued

| Headings within § 60.774 | Part 63, subpart G (HON)  | Part 63 (standards)                                  | Part 60, subpart QQQ (petroleum refinery)   | Part 264, subpart CC (RCRA CC)  | Part 265, subpart CC (RCRA CC)  | Part 61, subpart FF (benzene waste)  |
|--------------------------|---|--|---|---|---|--|
| Recordkeeping .....      | If WMU subject to HON, can comply with either HON or YYY provisions; if WMU not subject to HON, comply with YYY provisions. | Comply with standards provisions or YYY provisions*. | If WMU subject to Petroleum Refinery, then can comply with either Petroleum Refinery or YYY provisions; if WMU not subject to Petroleum Refinery, comply with YYY provisions. | If WMU subject to RCRA, part 264, then can comply with either RCRA, part 264 or YYY provisions; if WMU not subject to RCRA, part 264, comply with YYY provisions. | If WMU subject to RCRA, part 265, then can comply with either RCRA, part 265 or YYY provisions; if WMU not subject to RCRA, part 265, comply with YYY provisions. | If WMU subject to Benzene Waste, then can comply with either Benzene Waste or YYY provisions; if WMU not subject to Benzene Waste, comply with YYY provisions. |
| Leak Detection .....     | If WMU subject to HON, can comply with either HON or YYY provisions; if WMU not subject to HON, comply with YYY provisions. | Comply with standards provisions or YYY provisions.  | If WMU subject to Petroleum Refinery, then can comply with either Petroleum Refinery or YYY provisions; if WMU not subject to Petroleum Refinery, comply with YYY provisions. | If WMU subject to RCRA, part 264, then can comply with either RCRA, part 264 or YYY provisions; if WMU not subject to RCRA, part 264, comply with YYY provisions. | If WMU subject to RCRA, part 265, then can comply with either RCRA, part 265 or YYY provisions; if WMU not subject to RCRA, part 265, comply with YYY provisions. | If WMU subject to Benzene Waste, then can comply with either Benzene Waste or YYY provisions; if WMU not subject to Benzene Waste, comply with YYY provisions. |
| Delay of repair .....    | If WMU subject to HON, can comply with either HON or YYY provisions; if WMU not subject to HON, comply with YYY provisions. | Comply with standards provisions or YYY provisions.  | If WMU subject to Petroleum Refinery, then can comply with either Petroleum Refinery or YYY provisions; if WMU not subject to Petroleum Refinery, comply with YYY provisions. | If WMU subject to RCRA, part 264, then can comply with either RCRA, part 264 or YYY provisions; if WMU not subject to RCRA, part 264, comply with YYY provisions. | If WMU subject to RCRA, part 265, then can comply with either RCRA, part 265 or YYY provisions; if WMU not subject to RCRA, part 265, comply with YYY provisions. | If WMU subject to Benzene Waste, then can comply with either Benzene Waste or YYY provisions; if WMU not subject to Benzene Waste, comply with YYY provisions. |
| Control device .....     | If WMU subject to HON, can comply with either HON or YYY provisions; if WMU not subject to HON, comply with YYY provisions. | Comply with YYY provisions.                          | If WMU subject to Petroleum Refinery, then can comply with either Petroleum Refinery or YYY provisions; if WMU not subject to Petroleum Refinery, comply with YYY provisions. | If WMU subject to part RCRA 264, then can comply with either RCRA part 264 or YYY provisions; if WMU not subject to RCRA part 264, comply with YYY provisions.    | If WMU subject to RCRA, part 265, then can comply with either RCRA part 265 or YYY provisions; if WMU not subject to RCRA part 265, comply with YYY provisions.   | If WMU subject to Benzene Waste, then can comply with either Benzene Waste or YYY provisions; if WMU not subject to Benzene Waste, comply with YYY provisions. |

\* The owner or operator shall comply with the recordkeeping and reporting provisions in §§ 60.784 and 60.785 of this subpart when complying with the provisions for containers in subpart PP.

\*\* The owner or operator shall comply with the reporting provisions in § 60.784 of this subpart when complying with the provisions of RCRA, 40 CFR part 265.

TABLE 6 TO SUBPART YYY—CONTROL REQUIREMENTS FOR ITEMS OF EQUIPMENT THAT MEET THE CRITERIA OF § 60.775

| Item of equipment          | Control requirement <sup>a</sup>   |
|----------------------------|--|
| Drain or drain hub .....   | (a) Tight fitting solid cover (TFSC); or<br>(b) TFSC with a vent to either a fuel gas system or to a control device meeting the requirements of § 60.780; or<br>(c) Water seal with submerged discharge or barrier to protect discharge from wind.   |
| Manhole <sup>b</sup> ..... | (a) TFSC; or<br>(b) TFSC with a vent to a control device meeting the requirements of § 60.780; or<br>(c) If the item is vented to the atmosphere, use a TFSC with a properly operating water seal at the entrance or exit to the item to restrict ventilation in the collection system. The vent pipe shall be at least 90 cm in length and not exceeding 10.2 cm in diameter. |
| Lift station .....         | (a) TFSC; or<br>(b) TFSC with a vent to a control device meeting the requirements of § 60.780; or  |

TABLE 6 TO SUBPART YYY—CONTROL REQUIREMENTS FOR ITEMS OF EQUIPMENT THAT MEET THE CRITERIA OF § 60.775—Continued

| Item of equipment         | Control requirement <sup>a</sup>  |
|---------------------------|---|
| Trench .....              | (c) If the lift station is vented to the atmosphere, use a TFSC with a properly operating water seal at the entrance or exit to the item to restrict ventilation in the collection system. The vent pipe shall be at least 90 cm in length and not exceeding 10.2 cm in nominal inside diameter. The lift station shall be level controlled to minimize changes in the liquid level.<br>(a) TFSC; or<br>(b) TSFC with a vent to a control device meeting the requirements of § 60.780; or<br>(c) If the item is vented to the atmosphere, use a TFSC with a properly operating water seal at the entrance or exit to the item to restrict ventilation in the collection system. The vent pipe shall be at least 90 cm in length and not exceeding 10.2 cm in nominal inside diameter. |
| Pipe .....                | Each pipe shall have no visible gaps in joints, seals, or other emission interfaces.  |
| Oil/Water separator ..... | (a) Equip with a fixed roof and closed vent system that routes vapors to a control device meeting the requirements of § 60.780; or<br>(b) Equip with a floating roof that meets the equipment specifications of § 60.693 (a)(1)(i), (a)(1)(ii), (a)(2), (a)(3), and (a)(4).   |
| Tank <sup>c</sup> .....   | Maintain a fixed roof <sup>d</sup> . If the tank is sparged <sup>e</sup> or used for heating or treating by means of an exothermic reaction, a fixed roof and a closed vent system shall be maintained that routes the VOC vapors to a control device that meets the requirements of § 60.780.  |

<sup>a</sup> Where a tight fitting solid cover (TFSC) is required, it shall be maintained with no visible gaps or openings, except during periods of sampling, inspection, or maintenance.

<sup>b</sup> Manhole includes sumps and other points of access to a conveyance system.

<sup>c</sup> Applies to tanks with capacities of 38 m<sup>3</sup> or greater.

<sup>d</sup> A fixed roof may have openings necessary for proper venting of the tank, such as pressure/vacuum vent, j-pipe vent.

<sup>e</sup> The liquid in the tank is agitated by injecting compressed air or gas.

TABLE 7 TO SUBPART YYY—MONITORING REQUIREMENTS FOR TREATMENT PROCESSES

| To comply with  | Parameters to be monitored   | Frequency  | Methods  |
|---|--|--|--|
| 1. Required mass removal of each organic treated in a properly operated biological treatment unit § 60.779. | Appropriate parameters as specified in § 60.781(c) and approved by the permitting authority.   | Appropriate frequency as specified in § 60.781(c) and as approved by permitting authority. | Appropriate methods as specified in § 60.781(c) and as approved by permitting authority.                     |
| 2. Design steam stripper § 60.779(d).   | Steam flow rate .....  | Continuously .....   | Integrating steam flow monitoring device equipped with a continuous recorder.                                |
|   | Wastewater feed mass flow rate ..  | Continuously .....   | Liquid flow meter installed at stripper influent and equipped with a continuous recorder.                    |
|   | Wastewater feed temperature .....  | Continuously .....   | Liquid temperature monitoring device installed at stripper influent and equipped with a continuous recorder. |
| 3. Alternative monitoring parameters.   | Other parameters may be monitored upon approval from the Administrator in accordance with the requirements specified in § 60.781(d). | .....  |  |

TABLE 8 TO SUBPART YYY—MONITORING REQUIREMENTS FOR CONTROL DEVICES

| Control device         | Monitoring equipment required  | Parameters to be monitored  | Frequency   |
|------------------------|--|---|---|
| All control devices .. | 1. Flow indicator installed at all bypass lines to the atmosphere and equipped with continuous recorder <sup>b</sup> or.<br><br>2. Valves sealed closed with car-seal or lock-and-key configuration. | 1. Diversion to the atmosphere from the control device or.<br><br>2. Monthly inspections of sealed valves | Hourly records of whether the flow indicator was operating and whether a diversion was detected at any time during each hour.<br>Record and report the times of all periods when emissions are diverted through a bypass line or the flow indicator is not operating.<br>Monthly. |
| Thermal Incinerator    | Temperature monitoring device installed in firebox or in ductwork immediately downstream of firebox <sup>a</sup> and equipped with a continuous recorder <sup>b</sup> .                              | Firebox temperature .....   | Continuous.   |

TABLE 8 TO SUBPART YYY—MONITORING REQUIREMENTS FOR CONTROL DEVICES—Continued

| Control device   | Monitoring equipment required  | Parameters to be monitored   | Frequency  |
|--|--|--|--|
| Catalytic Incinerator  | Temperature monitoring device installed in gas stream immediately before and after catalyst bed and equipped with a continuous recorder <sup>b</sup> . | 1. Temperature upstream and downstream of catalyst bed.<br>2. Temperature difference across catalyst bed.  | Continuous.  |
| Flare .....  | Heat sensing device installed at the pilot light and equipped with a continuous recorder <sup>b</sup> .  | Presence of a flame at the pilot light ...   | Hourly records of whether the monitor was continuously operating and whether the pilot flame was continuously present during each hour.  |
| Boiler or process heater <44 megawatts and vent stream is not mixed with the primary fuel. | Temperature monitoring device installed in firebox <sup>a</sup> and equipped with continuous recorder <sup>b</sup> .                                   | Combustion temperature .....   | Continuous.  |
| Condenser .....  | Temperature monitoring device installed at condenser exit and equipped with continuous recorder <sup>b</sup> .   | Condenser exit (product side) temperature.   | Continuous.  |
| Carbon Adsorber (Regenerative).  | Integrating regeneration stream flow monitoring device having an accuracy of ±10 percent, and.<br>Carbon bed temperature monitoring device.            | 1. Total regeneration stream mass or volumetric flow during carbon bed regeneration cycle(s).<br>2. Temperature of carbon bed after regeneration [and within 15 minutes of completing any cooling cycle(s)]. | For each regeneration stream mass or volumetric flow.<br>For each regeneration cycle and within 15 minutes of completing any cooling cycle, record the carbon bed temperature. |
| Carbon Adsorber (Non-regenerative).  | Organic compound concentration monitoring device <sup>c</sup> .  | Organic compound concentration of adsorber exhaust.  | Daily or at intervals no greater than 20 percent of the design carbon replacement interval, whichever is greater.  |
| Alternative monitoring parameters.   | Other parameters may be monitored upon approval from the Administrator in accordance with the requirements in § 60.781(e)(3).                          | .....  |  |

<sup>a</sup> Monitor may be installed in the firebox or in the ductwork immediately downstream of the firebox before any substantial heat exchange is encountered.

<sup>b</sup> "Continuous recorder" is defined in § 60.771 of this subpart.

<sup>c</sup> As an alternative to conducting this monitoring, an owner or operator may replace the carbon in the carbon adsorption system with fresh carbon at a regular predetermined time interval that is less than the carbon replacement interval that is determined by the maximum design flow rate and organic concentration in the gas stream vented to the carbon adsorption system.

TABLE 9 TO SUBPART YYY—INFORMATION ON PROCESS WASTEWATER STREAMS TO BE SUBMITTED WITH NOTIFICATION OF COMPLIANCE STATUS<sup>a, b</sup>

| Process unit identification code <sup>c</sup> | Stream identification code | VOC concentration (ppmw) <sup>d, e</sup> | Flow rate (lpm) <sup>e, f</sup> | Group 1 or Group 2 <sup>g</sup> | Compliance approach <sup>h</sup> | Treatment process(es) identification <sup>i</sup> | Waste management unit(s) identification | Intended control device |
|---|----------------------------|--|---------------------------------|---------------------------------|----------------------------------|---|---|-------------------------|
|   |                            |  |                                 |                                 |                                  |   |   |                         |

<sup>a</sup> The information specified in this table 9 must be submitted; however, it may be submitted in any format. This table 9 presents an example format.

<sup>b</sup> Other requirements for the Notification of Compliance Status are specified in § 60.784(c).

<sup>c</sup> Also include a description of the process unit (e.g., benzene process unit).

<sup>d</sup> Except when § 60.773(c) is used, annual average concentration as specified in § 60.773(b) and § 60.782.

<sup>e</sup> When § 60.773(c) is used, indicate the wastewater stream is a designated Group 1 wastewater stream.

<sup>f</sup> Except when § 60.773(c) is used, annual average flowrate as specified in § 60.773(b) and § 60.782.

<sup>g</sup> Indicate whether stream is Group 1 or Group 2.

<sup>h</sup> Cite § 60.779 compliance option used.

<sup>i</sup> Identification codes should correspond to those listed in Table 10 of this subpart.



TABLE 10 TO SUBPART YYY—INFORMATION FOR TREATMENT PROCESSES TO BE SUBMITTED WITH NOTIFICATION OF COMPLIANCE STATUS <sup>a, b</sup>

| Treatment process identification <sup>c</sup> | Description <sup>d</sup> | Wastewater stream(s) treated <sup>e</sup> | Monitoring parameters <sup>f</sup> |
|---|--------------------------|---|------------------------------------|
|   |                          |   |                                    |

<sup>a</sup>The information specified in this table 10 must be submitted; however, it may be submitted in any format. This table 10 presents an example format.

<sup>b</sup>Other requirements for the Notification of Compliance Status are specified in § 60.784(c) of this subpart.

<sup>c</sup>Identification codes should correspond to those listed in Table 9 of this subpart.

<sup>d</sup>Description of treatment process (e.g., steam stripper).

<sup>e</sup>Stream identification code for each wastewater stream treated by each treatment unit. Identification codes should correspond to entries listed in Table 9 of this subpart.

<sup>f</sup>Parameter(s) to be monitored or measured in accordance with Table 7 and § 60.781 of this subpart.

TABLE 11 TO SUBPART YYY—INFORMATION FOR WASTE MANAGEMENT UNITS TO BE SUBMITTED WITH NOTIFICATION OF COMPLIANCE STATUS <sup>a, b</sup>

| Waste management unit Identification <sup>c</sup> | Description <sup>d</sup> | Wastewater stream(s) received or managed <sup>e</sup> |
|---|--------------------------|---|
|   |                          |   |

<sup>a</sup>The information specified in this table 11 must be submitted; however, it may be submitted in any format. This table 11 presents an example format.

<sup>b</sup>Other requirements for the Notification of Compliance Status are specified in § 60.784(c) of this subpart.

<sup>c</sup>Identification codes should correspond to those listed in Table 9 of this subpart.

<sup>d</sup>Description of waste management unit.

<sup>e</sup>Stream identification code for each wastewater stream received or managed by each waste management unit. Identification codes should correspond to entries listed in Table 9 of this subpart.

TABLE 12 TO SUBPART YYY—INFORMATION ON RESIDUALS TO BE SUBMITTED WITH NOTIFICATION OF COMPLIANCE STATUS <sup>a, b</sup>

| Residual identification <sup>c</sup> | Residual description <sup>d</sup> | Wastewater stream identification <sup>e</sup> | Treatment process <sup>f</sup> | Fate <sup>g</sup> | Control device identification code | Control device description <sup>h</sup> | Control device efficiency <sup>i</sup> |
|--------------------------------------|-----------------------------------|---|--------------------------------|-------------------|------------------------------------|---|--|
|                                      |                                   |   |                                |                   |                                    |   |  |

<sup>a</sup>The information specified in this table 12 must be submitted; however, it may be submitted in any format. This table 12 presents an example format.

<sup>b</sup>Other requirements for the Notification of Compliance Status are specified in § 60.784(c) of this subpart.

<sup>c</sup>Name or identification code of residual removed from Group 1 wastewater stream.

<sup>d</sup>Description of residual (e.g., steam stripper A-13 overhead condensates).

<sup>e</sup>Identification of stream from which residual is removed.

<sup>f</sup>Treatment process from which residual originates.

<sup>g</sup>Indicate whether residual is sold, returned to production process, or returned to waste management unit or treatment process; or whether VOC mass of residual is destroyed by 99 percent.

<sup>h</sup>If the fate of the residual is such that the VOC mass is destroyed by 99 percent, give description of device used for VOC destruction.

<sup>i</sup>The fate of the residual is such that the VOC mass is destroyed by 99 percent, provide an estimate of control device efficiency and attach substantiation in accordance with § 60.784(c)(5) of this subpart.

TABLE 13 TO SUBPART YYY—SEMIANNUAL REPORTING REQUIREMENTS FOR CONTROL DEVICES [§ 60.784(F)]

| Control Device              | Reporting Requirements   |
|-----------------------------|--|
| Thermal Incinerator .....   | 1. Report all daily average <sup>a</sup> temperatures that are above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit and all operating days when insufficient monitoring data are collected. <sup>c</sup> |
| Catalytic Incinerator ..... | 1. Report all daily average <sup>a</sup> upstream temperatures that are above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit.  |

TABLE 13 TO SUBPART YYY—SEMIANNUAL REPORTING REQUIREMENTS FOR CONTROL DEVICES [§ 60.784(F)]—Continued

| Control Device  | Reporting Requirements  |
|---|---|
| Boiler or Process Heater with a design heat input capacity less than 44 megawatts and vent stream is not mixed with the primary fuel. | 2. Report all daily average <sup>a</sup> temperature differences across the catalyst bed that are above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit.<br>3. Report all operating days when insufficient monitoring data are collected. <sup>c</sup>   |
| Flare .....   | 1. Report all daily average <sup>a</sup> firebox temperatures that are above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit and all operating days when insufficient monitoring data are collected. <sup>c</sup>  |
| Condenser .....   | 1. Report the duration of all periods when all pilot flames are absent.   |
| Carbon Adsorber .....   | 1. Report all daily average <sup>a</sup> exit temperatures that are above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit and all operating days when insufficient monitoring data are collected. <sup>c</sup>   |
| All Control Devices .....   | 1. Report all carbon bed regeneration cycles when the total regeneration stream mass or volumetric flow is above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit.<br>2. Report all carbon bed regeneration cycles during which the temperature of the carbon bed after regeneration is above the maximum or below the minimum operating parameter value established in the NCS <sup>b</sup> or operating permit.<br>3. Report all operating days when insufficient monitoring data are collected. <sup>c</sup> |
|   | 1. Report the times and durations of all periods when the vent stream is diverted through a bypass line or the monitor is not operating, or<br>2. Report all monthly inspections that show the valves are moved to the diverting position or the seal has been changed.   |

<sup>a</sup> The daily average is the average of all values recorded during the operating day, as specified in § 60.785(e) of this subpart.

<sup>b</sup> NCS = Notification of Compliance Status described in § 60.784(c) of this subpart.

<sup>c</sup> The semiannual reports shall include the duration of periods when monitoring data are not collected for each excursion as defined in § 60.784(d)(3) of this subpart.

TABLE 14 TO SUBPART YYY—COMPOUND AND DEFAULT BIORATES USED FOR COMPLIANCE DEMONSTRATIONS FOR ENHANCED BIOLOGICAL TREATMENT PROCESSES (SEE § 60.783(H))

| Compound name  | Biorate, K1<br>L/g MLVSS-hr |
|--|-----------------------------|
| Acetonitrile .....                                   | 0.100                       |
| Acetophenone .....                                   | 0.538                       |
| Acrylonitrile .....                                  | 0.750                       |
| Biphenyl .....                                       | 5.643                       |
| Chlorobenzene .....                                  | 10.000                      |
| Dichloroethyl Ether (bis (2-chloroethyl ether) ..... | 0.246                       |
| Diethyl Sulfate .....                                | .0105                       |
| Dimethyl Hydrazine (1,1-) .....                      | 0.227                       |
| Dimethyl Sulfate .....                               | 0.178                       |
| Dinitrophenol (2,4-) .....                           | 0.620                       |
| Dinitrotoluene (2,4-) .....                          | 0.784                       |
| Dioxane (1,4-) (1,4-diethylene oxide) .....          | 0.393                       |
| Ethylene Glycol Dimethyl Ether .....                 | 0.364                       |
| Ethylene Glycol Monobutyl Ether Acetate .....        | 0.496                       |
| Ethylene Glycol Monomethyl Ether Acetate .....       | 0.159                       |
| Hexachlorobenzene .....                              | 16.179                      |
| Isophorone .....                                     | 0.598                       |
| Methanol .....                                       | 0.200                       |
| Methyl Methacrylate .....                            | 4.300                       |
| Nitrobenzene .....                                   | 2.300                       |
| Toluidine (-o) .....                                 | 0.859                       |
| Trichlorobenzene (1,2,4-) .....                      | 4.393                       |
| Trichlorophenol (2,4,5-) .....                       | 4.477                       |
| Triethylamine .....                                  | 1.064                       |

3. Part 60 is amended by adding appendix J to read as follows:

**Appendix J to Part 60—How to Determine Henry's Law Constants, Fm Values, Fr Values, and Fe Values for Organic Compounds**

1. *Use of Appendix and General Information.* This appendix has four sections.

Section 2 contains the procedures for determining Henry's law constants, fraction measured (Fm) values, fraction removed values (Fr), and fraction emitted (Fe) values for an individual chemical. Section 3 describes how to locate certain resources. Section 4 contains five tables and thirteen forms.

1.1 You should use this appendix if you need to:

1. Determine whether a chemical has a Henry's law constant at 25° C that is less than 0.1 y/x atmosphere per mole fraction (see section 2.1).

2. Determine a fraction measured (Fm) value for a chemical (see section 2.2).

3. Subtract the concentration of a chemical from a Method 25D concentration (see section 2.3).

4. Determine the fraction removed (Fr) value for a chemical that has a Henry's law constant at 25° C that is greater than or equal to 0.1 y/x atmosphere per mole fraction (see section 2.4).

5. Determine the fraction emitted (Fe) value for a chemical that has a Henry's law constant at 25° C that is greater than or equal to 0.1 y/x atmosphere per mole fraction (see section 2.5).

6. Calculate a Henry's law constant at a specific temperature using a Henry's law constant at a different temperature for the same chemical (see section 2.6).

1.2 This appendix requires documentation for some procedures. The referencing subpart, i.e., the rule to which you are complying, may require additional recordkeeping and may specify records concerning this appendix that are to be included in reports.

1.3 When the term "WATER8" is used in this appendix, the term "WATER8, or updates to WATER8" must be used for the purposes of this appendix. When the term "CHEM9" is used in this appendix, the term "CHEM9, or updates to CHEM9" must be used for the purposes of this appendix. When the terms "waste" or "wastewater" are used in this appendix, the term "waste or wastewater, as applicable to the referencing subpart" must be used for the purposes of this appendix. When the terms "Henry's law constant" or "Henry's law constants" are used in this appendix, the terms "Henry's law constant(s) with units of atmosphere per mole fraction" must be used for the purposes of this subpart.

## 2. Procedures.

2.1 *How to determine whether a chemical has a Henry's law constant at 25° C that is less than 0.1 y/x.* You must use one of the following to determine whether a chemical has a Henry's law constant that is less than 0.1 y/x atmosphere per mole fraction.

2.1.1 *Use Table 1.* The chemicals listed in Table 1 have a Henry's law constant at 25° C that is less than 0.1 y/x atmosphere per mole fraction.

2.1.2 *Use CHEM9 or WATER8.* Use CHEM9 or WATER8 to determine the Henry's law constant at 25° C. You must know compound properties, such as solubility in water and vapor pressure, and the structure of the compound to estimate a Henry's law constant using CHEM9 or WATER8.

2.1.3 *Determine experimentally.* The Henry's law constant may be measured by several laboratory techniques. These techniques can be categorized as either two phase closed systems techniques or open system techniques.

2.1.3.1 *Two phase closed systems.* For two phase closed system techniques, the volume of each phase and two concentration measurements are needed. The concentration measurements are: (1) concentration in one of the phases, and (2) either the concentration in the other phase or the total concentration in both phases. Use Form 1 to calculate the Henry's law constant for two phase closed systems.

2.1.3.2 *Open systems.* For open systems, gas is passed through a liquid volume containing

the compound. The Henry's law constant is calculated from the rate of stripping of the compound from the water. Use Form 2 to calculate the Henry's law constant for open systems.

2.1.4 *Calculate a Henry's law constant at 25° C from a Henry's law constant at a different temperature for the same chemical.* Use the procedures specified in section 2.6 to calculate a Henry's law constant at 25° C from a Henry's law constant at a different temperature for the same chemical.

2.2. *How to determine a Fm value for a chemical.* Fm means compound-specific fraction measured factor, and it has the units of mass measured by Method 25D divided by the total mass in the wastewater. You must use one of the following to determine the Fm value for a chemical.

2.2.1 *Use Table 1 or Table 2.* To determine the Fm value for a chemical with a Henry's law constant at 25° C that is less than 0.1 y/x atmosphere per mole fraction, use the Fm value listed for the chemical in Table 1. To determine the Fm value for a chemical with a Henry's law constant at 25° C that is greater than or equal to 0.1 y/x, use the Fm value listed for the chemical in Table 2.

Note to section 2.2.1: Table 1 and Table 2 include Fm values for Method 25D and for Method 305. Unless otherwise specified in this appendix or the referencing subpart, use the Fm values for Method 25D.

2.2.2 *Use CHEM9.* Use CHEM9 to determine an Fm value. You must know the structure of the chemical and certain other compound properties, e.g., boiling point, Antoine's coefficients, vapor pressure, and solubility in water, to estimate an Fm value using CHEM9. The accuracy of the computer estimation procedure depends on the nature of the compound and the quality of the available data. The procedure is flexible in that the method can be used with a variety of different types of compound data. You must confirm and document the compound properties used as inputs for CHEM9 and the lack of availability for missing compound properties. In some cases, this method is not accurate, especially with missing compound properties. Before accepting the estimation values of CHEM9 in these cases, you must document the consistency of the predicted values with other related experimental data.

2.2.3 *Measure the Fm value.* Spike a sample of waste with a known amount of the compound of interest. Measure the concentration of the sample using Method 25D. The Fm value for the recovery of a specific chemical is the ratio of the Method 25D concentration to the actual concentration in the waste sample. You must minimize loss of organic compounds during sample collection and analysis, and maintain sample integrity. An example of acceptable sampling and handling procedures are the sampling and handling requirements in Method 25D.

2.2.4 *Extrapolating a Method 25D Fm Value from a Method 305 Fm value.* Method 305 measures the recovered concentration, not the actual concentration in the wastewater. The Method 25D correction value may be obtained from the Method 305 value and the ratio of the Method 25D value to the Method 305 value for that compound.

This ratio for a compound is independent of the wastewater and may be determined once for each compound.

2.3 *How to subtract a chemical from a Method 25D concentration.* You must follow the procedures specified in sections 2.3.1 through 2.3.5 to subtract a chemical's concentration from the total concentration measured by Method 25D. You may only subtract from the total Method 25D concentration compounds for which you have a measured concentration (i.e., you must not subtract compounds for which test results are below the quantification limit.) If an Fm value cannot be determined for a chemical, the concentration of the chemical cannot be subtracted from the Method 25D results. You must follow the procedures in Form 3 to subtract a chemical from a Method 25D concentration. Form 4 provides an example.

2.3.2 *Determine the concentration for each chemical in the wastewater stream that will be subtracted from the Method 25D concentration.* The concentration for each chemical must be determined using a method and sampling procedure specified in the referencing subpart. Methods other than Method 25D and Method 305 are considered alternative methods for the purposes of this appendix.

2.3.3 *Determine the correct Fm value.* If an Fm value is needed, use the procedures in section 2.2 of this appendix to determine the correct Fm value.

2.3.4 *Adjust the concentration of chemicals which may be subtracted from the Method 25D concentration.* You must multiply the concentration of the chemical measured by the alternative method (i.e., a method that is not Method 25D or Method 305 and that is specified in the referencing subpart) by the Method 25D Fm. The product will be the adjusted concentration for that chemical. This adjustment must be done for each chemical you subtract from the concentration measured by Method 25D.

2.3.5 *Subtract.* Subtract the product(s) you calculated from the Method 25D concentration.

2.4 *How to determine an Fr value for a chemical with a Henry's law constant at 25° C that is greater than or equal to 0.1 y/x.* Fr means fraction removal value and is unitless. You must use one of the following to determine a Fr value.

2.4.1 *Use Table 2.* Use the Fr value listed for the chemical in Table 2. The chemicals listed in table 2 have a Henry's law constant at 25° C that is greater than or equal to 0.1 y/x.

2.4.2 *Use 0.99.* Assign an Fr value of 0.99 to any chemical. This is the highest Fr value that is assigned to a chemical.

2.4.3 *Use CHEM9.* Use CHEM9 to determine the Fr value of a chemical. You must know the compound structure and the Henry's law constant at 100° C to estimate an Fr value using CHEM9. The Henry's law constant at 100° C for a chemical must be determined as specified in either section 2.4.3.1, 2.4.3.2, or 2.4.3.3. The method used to determine the Henry's law constant at 100° C for a chemical must be documented.

2.4.3.1 *Determine Henry's law at 100° C experimentally.* The Henry's law constant

may be measured by several laboratory techniques. These techniques can be categorized as either two phase closed systems techniques or open system techniques.

2.4.3.1.1 *Two phase closed systems.* For two phase closed system techniques, the volume of each phase and two concentration measurements are needed. The concentration measurements are: (1) concentration in one of the phases, and (2) either the concentration in the other phase or the total concentration in both phases. Use Form 1 to calculate the Henry's law constant for two phase closed systems.

2.4.3.1.2 *Open systems.* For open systems, gas is passed through a liquid volume containing the compound. The Henry's law constant is calculated from the rate of stripping of the compound from the water. Use Form 2 to calculate the Henry's law constant for open systems.

2.4.3.2 *Calculate a Henry's law constant at 100° C from a Henry's law constant at a different temperature for the same chemical.* Use the procedures in section 2.6 to calculate a Henry's law constant at 100° C from a Henry's law constant at a different temperature for the same chemical.

2.4.3.3 *Literature Value.* Experimental values of Henry's law constants at a 100° C for some chemicals are available in data bases or reported in the literature. You must provide the reference for and description of any database or literature you used.

2.5 *How to determine an Fe value for a chemical that has a Henry's law constant at 25° C that is greater than or equal to 0.1 y/x.* Use the appropriate Fe value as specified in the referencing subpart.

2.5.1 *Default Fe values for emissions from both the individual drain system and the treatment process.* You must measure the temperature of the wastewater stream at the point of determination, unless another location is specified by the referencing subpart. If the temperature of the wastewater stream is less than or equal to 35° C, you may

use the default Fe values listed in either Table 2 or Table 3. If the temperature of the wastewater stream is greater than 35° C, you must use the default Fe values listed in Table 3.

2.5.1.1 *Use Table 2.* To use Table 2, use the default Fe value listed for the chemical in Table 2.

2.5.1.2 *Use Table 3.* You must either use a default Fe listed in Table 3 or use Table 3 to interpolate an Fe value. To use Table 3, you must determine the chemical's Henry's law constant at the temperature you measured for the wastewater stream. You must find this Henry's law constant in the table and select an Fe value greater than or equal to the Fe value that corresponds to the Henry's law constant.

2.5.2 *Site-specific Fe values for emissions from the individual drain system.* Use WATER8 and Forms 6 and 7 for each type of waste management unit modeled and Forms 8 through 13, as appropriate for the different types of waste management units. (Note that this Fe value does not include Fe values for the treatment process.)

2.5.3 *Default Fe values for emissions from the biological treatment process (Fet).* The default Fe values in Table 4 and Table 5 are Fe values for the biological treatment system (i.e., the wastewater treatment plant) and have been assigned the abbreviation "Fet." You must measure the temperature of the wastewater stream(s) treated in the biological treatment system at the inlet to the biological treatment system (e.g., at the bar screen). If the temperature of the wastewater stream(s) is less than or equal to 35° C, you must use either Table 4 or Table 5 to determine the Fet value. If the temperature of the wastewater stream is greater than 35° C, you must use Table 5 to determine the Fet value.

2.5.3.1 *Use Table 4.* To use Table 4, use the default Fet value listed for the chemical in Table 4.

2.5.3.2 *Use Table 5.* To use Table 5, you must either use a default Fet listed in Table 5 or use Table 5 to interpolate an Fet value.

You must determine the chemical's Henry's law constant at the temperature you measured for the wastewater stream. You must find this Henry's law constant in the table and select an Fet value greater than or equal to the Fet value that corresponds to the Henry's law constant.

2.6 *How to calculate a Henry's law constant from a Henry's law constant at a different temperature for the same chemical.* Use WATER8 and Form 5 to estimate a Henry's law constant from a Henry's law constant at a different temperature for the same chemical.

3. *Location of resources.*

3.1 *Where to find information on CHEM9 and WATER8.*

3.1.1 *CHEM9 and WATER8 access via Internet.* You can find CHEM9 and WATER8 on the Internet by accessing EPA's Technology Transfer Network (TTN) via the Internet. The Internet address is: <http://www.epa.gov/ttn/chief/software.html>. If you need more information on the TTN, contact the systems operator at (919) 541-5384.

3.1.2 *Procedures used in CHEM9.* Reports describing the CHEM9 procedures for estimating Fm, Fr, and Fe values are in Docket Number A-94-32, Item IV-A-1. The database for CHEM9 is not available as a hard copy.

Docket No. A-94-32 is available for public inspection and copying between 8:00 a.m. and 5:30 p.m., Monday through Friday, at the EPA's Air and Radiation Docket and Information Center, Waterside Mall, Room M-1500, first floor, 401 M Street SW, Washington, DC 20460, or by calling (202) 260-7548 or 260-7549. A reasonable fee may be charged for copying.

3.2 *Methods.*

*Method 25D* can be found in 40 CFR part 60, Appendix A.

*Method 305* can be found in 40 CFR part 63, Appendix A.

4. *Tables and Forms.* This section contains 5 tables and 13 forms.

TABLE 1 OF APPENDIX J.—FM VALUES FOR HENRY'S LAW CONSTANTS AT 25°C LESS THAN 0.1 (Y/X) ATMOSPHERES PER MOLE FRACTION

[Use with Section 2.1]

| Compound                               | Y/X      | Fm 25D | Fm 305 |
|--|----------|--------|--------|
| 1H IMIDAZOLE .....                     | 0.000004 | 0.001  | 0.001  |
| 2,4 D .....                            | 0.000000 | 0.151  | 0.167  |
| 2,4,5 BENZOIC ACID .....               | 0.000007 | 0.000  | 0.000  |
| 2-HYDROXYETHANAL .....                 | 0.001400 | 0.031  | 0.059  |
| 3,4-DIMETHYLPHENOL xyleneol .....      | 0.004200 | 0.018  | 0.017  |
| 3,5-DIBROMO-4HYDROXYBENZONITRILE ..... | 0.011700 | 0.021  | 0.033  |
| 3-OXOPROPANOIC ACID .....              | 0.007900 | 0.002  | 0.004  |
| 4-OXOBUTANOIC ACID .....               | 0.011100 | 0.004  | 0.006  |
| 5-OXOPENTANOIC ACID .....              | 0.013900 | 0.005  | 0.007  |
| ACETALDOL .....                        | 0.001900 | 0.011  | 0.016  |
| ACETAMIDE .....                        | 0.000100 | 0.305  | 0.463  |
| ACETYL-2-THIOUREA, 1- .....            | 0.001600 | 0.034  | 0.053  |
| ACETYL-5-HYDROXYPIPERIDINE 3 .....     | 0.038900 | 0.001  | 0.001  |
| ACETYLAMINOFLUORENE, 2- .....          | 0.074400 | 0.020  | 0.018  |
| ACETYLPYPERIDINE 3 .....               | 0.006900 | 0.151  | 0.175  |
| ACRIDINE ORANGE* .....                 | 0.013300 | 0.050  | 0.049  |
| ACRIDINE YELLOW* .....                 | 0.000400 | 0.001  | 0.001  |
| ACRYLAMIDE .....                       | 0.000015 | 0.003  | 0.003  |
| ACRYLIC ACID .....                     | 0.011000 | 0.431  | 0.643  |
| ADAMANTANE DICARBOXYLIC ACID .....     | 0.002600 | 0.001  | 0.001  |

TABLE 1 OF APPENDIX J.—FM VALUES FOR HENRY'S LAW CONSTANTS AT 25°C LESS THAN 0.1 (Y/X) ATMOSPHERES PER MOLE FRACTION—Continued

[Use with Section 2.1]

| Compound                             | Y/X      | Fm 25D | Fm 305 |
|--------------------------------------|----------|--------|--------|
| ADENINE                              | 0.000005 | 0.001  | 0.002  |
| ADIPIC ACID                          | 0.000003 | 0.001  | 0.001  |
| ADIPONITRILE                         | 0.000700 | 0.004  | 0.004  |
| ALACHLOR (M)                         | 0.001800 | 0.090  | 0.090  |
| alpha-PICOLINE                       | 0.025900 | 0.870  | 0.842  |
| AMETRYN                              | 0.000001 | 0.001  | 0.001  |
| AMINOBIIPHENYL, 4-                   | 0.017200 | 0.012  | 0.011  |
| AMINOETHYLPIPERAZINE                 | 0.000021 | 0.001  | 0.001  |
| AMINOPHENOL, 3-                      | 0.003400 | 0.035  | 0.040  |
| AMINOPYRIDINE, 4-                    | 0.000005 | 0.000  | 0.001  |
| ANILINE                              | 0.097800 | 0.142  | 0.138  |
| ANISIDINE, o-                        | 0.097200 | 0.011  | 0.013  |
| ANTHRAQUINONE                        | 0.000200 | 0.001  | 0.001  |
| ATRAZINE (M)                         | 0.000200 | 0.117  | 0.117  |
| BENZENE ACETIC ACID                  | 0.025500 | 0.014  | 0.015  |
| BENZENE ARSONIC ACID (M)             | 0.000006 | 0.124  | 0.124  |
| BENZENE DICARBOXYLIC ACID            | 0.000900 | 0.001  | 0.001  |
| BENZENE SULFONIC ACID (M)            | 0.043900 | 0.146  | 0.146  |
| BENZIDINE                            | 0.000001 | 0.000  | 0.000  |
| BENZO (A) ANTHRACENE                 | 0.000077 | 0.121  | 0.095  |
| BENZO(A) PYRENE                      | 0.000077 | 1.267  | 1.000  |
| BENZO (ghi) PERYLENE                 | 0.002800 | 0.006  | 0.005  |
| BENZO (k) FLUORANTHENE               | 0.000059 | 0.001  | 0.001  |
| BENZOIC ACID                         | 0.001000 | 0.003  | 0.003  |
| BENZOTHAZOLONE 2 (2H)-*              | 0.065600 | 0.121  | 0.123  |
| BENZYL ALCOHOL                       | 0.033900 | 0.069  | 0.067  |
| BHC, gamma-                          | 0.027400 | 1.035  | 0.973  |
| BIS (2-ETHYLHEXYL) PHTHALATE         | 0.016700 | 0.317  | 0.327  |
| BROMOCHLOROMETHYL ACETATE            | 0.010400 | 0.342  | 0.541  |
| BUTYL CELLOSOLVE                     | 0.014600 | 0.095  | 0.120  |
| BUTYL-m-CRESOL MONO T                | 0.052100 | 0.042  | 0.039  |
| BUTYL-p-CRESOL MONO T                | 0.052100 | 0.042  | 0.039  |
| BUTYRIC ACID                         | 0.096100 | 0.089  | 0.124  |
| CAPROLACTAM                          | 0.000200 | 0.002  | 0.003  |
| CAPROLACTONE                         | 0.071100 | 0.205  | 0.248  |
| CATECHOL                             | 0.000002 | 0.000  | 0.000  |
| CHLORACETOPHENONE, 2-                | 0.048400 | 0.161  | 0.152  |
| CHLORO (-p) CRESOL (-m)              | 0.009100 | 0.029  | 0.028  |
| CHLORO-1, 2-ETHANE DIOL (M)          | 0.005400 | 0.999  | 0.999  |
| CHLORO-2, 5-DIKETOPYRROLIDINE 3 (M)  | 0.003700 | 0.430  | 0.430  |
| CHLOROACETIC ACID                    | 0.003600 | 0.020  | 0.028  |
| CHLOROANILINE, p-                    | 0.014700 | 0.069  | 0.067  |
| CHLOROBENZOPHENONE (PARA)            | 0.000200 | 0.313  | 0.283  |
| CHLOROBENZYLATE                      | 0.000028 | 0.000  | 0.000  |
| CHLOROHYDRIN, a 3 CHLORO 1, 2 PROPAN | 0.000300 | 0.003  | 0.004  |
| CHLOROPHENOL POLYMERS (M)            | 0.005600 | 0.000  | 0.000  |
| CHLOROPHENOL-4                       | 0.062200 | 0.032  | 0.031  |
| CHOLINE CHLORIDE                     | 0.000600 | 0.012  | 0.015  |
| CHRYSENE                             | 0.000066 | 0.006  | 0.004  |
| CITRIC ACID                          | 0.000000 | 0.000  | 0.000  |
| CREOSOTE (M)                         | 0.004400 | 0.025  | 0.025  |
| CRESOL                               | 0.090000 | 0.049  | 0.047  |
| CRESOL (-m)                          | 0.039400 | 0.035  | 0.033  |
| CRESOL (-o)                          | 0.091200 | 0.057  | 0.055  |
| CRESOL (-p)                          | 0.039700 | 0.028  | 0.027  |
| CUMYLPHENOL-4                        | 0.093300 | 0.002  | 0.002  |
| CYANIDE methyl                       | 0.001500 | 0.328  | 0.417  |
| CYANOMETHYL BENZOATE 4 (M)           | 0.000700 | 0.128  | 0.128  |
| DIAZINON                             | 0.001200 | 0.001  | 0.001  |
| DIBENZO (a,h ) ANTHRACENE            | 0.002100 | 0.001  | 0.001  |
| DIBUTYLPHTHALATE                     | 0.015600 | 0.002  | 0.002  |
| DICHLORO—(2,6)-NITROANILINE (4) (M)  | 0.000400 | 0.122  | 0.122  |
| DICHLOROANILINE 2, 3                 | 0.029900 | 0.049  | 0.047  |
| DICHLOROBENZONITRILE,2 ,6-           | 0.064400 | 0.338  | 0.322  |
| DICHLOROPHENOL 2, 5                  | 0.086100 | 0.151  | 0.148  |
| DICHLOROTETRAHYDROFURAN 3, 4 (M)     | 0.007800 | 0.303  | 0.303  |
| DICHLORVOS                           | 0.019000 | 0.008  | 0.011  |
| DIETHANOLAMINE                       | 0.000000 | 0.000  | 0.000  |
| DIETHYL (N, N) ANILINE               | 0.003200 | 0.964  | 0.907  |

TABLE 1 OF APPENDIX J.—FM VALUES FOR HENRY'S LAW CONSTANTS AT 25°C LESS THAN 0.1 (Y/X) ATMOSPHERES PER MOLE FRACTION—Continued

[Use with Section 2.1]

| Compound                            | Y/X      | Fm 25D | Fm 305 |
|-------------------------------------|----------|--------|--------|
| DIETHYL PROPIONAMIDE, 2aN (M)       | 0.001100 | 0.089  | 0.089  |
| DIETHYLENE GLYCOL                   | 0.077800 | 0.000  | 0.000  |
| DIETHYLENE GLYCOL DIMETHYL ETHER    | 0.083800 | 0.105  | 0.150  |
| DIETHYLENE GLYCOL MONOBUTYL ETHER   | 0.001200 | 0.003  | 0.003  |
| DIETHYLENE GLYCOL MONOETHYL ETHER   | 0.002700 | 0.005  | 0.007  |
| DIETHYLENE GLYCOL MONOETHYL ETHER A | 0.035800 | 0.007  | 0.010  |
| DIETHYLENE GLYCOL MONOMETHYL ETHER  | 0.003200 | 0.004  | 0.007  |
| DIETHYLENETRIAMINE                  | 0.000001 | 0.000  | 0.000  |
| DIETHYLHYDRAZINE N, N               | 0.019000 | 0.184  | 0.253  |
| DIETHYLTHIOPHOSPHATEBENZO M ETHYL P | 0.001200 | 0.000  | 0.000  |
| DIMETHOATE (M)                      | 0.050900 | 0.110  | 0.110  |
| DIMETHYL CARBAMOYL CHLORIDE         | 0.024700 | 0.116  | 0.151  |
| DIMETHYL DISULFIDE                  | 0.083300 | 0.455  | 1.000  |
| DIMETHYL FORMAMIDE                  | 0.010600 | 0.009  | 0.013  |
| DIMETHYL HYDRAZINE (1, 1)           | 0.091100 | 0.277  | 0.382  |
| DIMETHYL PHTHALATE                  | 0.054800 | 0.006  | 0.007  |
| DIMETHYLAMINOAZOBENZENE, 4-         | 0.004100 | 0.022  | 0.023  |
| DIMETHYLBENZ (A) ANTHRACENE (7, 12) | 0.000015 | 0.008  | 0.006  |
| DIMETHYLBENZIDINE 3,3               | 0.000075 | 0.000  | 0.000  |
| DIMETHYLSULFONE                     | 0.001300 | 0.002  | 0.003  |
| DIMETHYLSULFOXIDE                   | 0.026900 | 0.037  | 0.057  |
| DINITRO- <i>o</i> -CRESOL (4, 6)    | 0.078000 | 0.009  | 0.016  |
| DIPHENYLHYDRAZINE (1, 2)            | 0.013600 | 0.462  | 0.448  |
| DIPROPYLENE GLYCOL                  | 0.000900 | 0.002  | 0.003  |
| ENDRIN                              | 0.084400 | 0.005  | 0.004  |
| EPINEPHRINE (M)                     | 0.020300 | 0.133  | 0.133  |
| ETHANOLAMINE (mono-)                | 0.017800 | 0.004  | 0.007  |
| ETHYL CARBAMATE                     | 0.000600 | 0.004  | 0.008  |
| ETHYL MORPHOLINE, ethyl diethylene  | 0.011300 | 0.048  | 0.059  |
| ETHYLENE GLYCOL                     | 0.000100 | 0.002  | 0.005  |
| ETHYLENE GLYCOL MONOBUTYL ETHER     | 0.029200 | 0.056  | 0.071  |
| ETHYLENE GLYCOL MONOETHYL ETHER     | 0.061700 | 0.111  | 0.144  |
| ETHYLENE GLYCOL MONOETHYL ETHER AC  | 0.098600 | 0.057  | 0.089  |
| ETHYLENE GLYCOL MONOMETHYL ETHER    | 0.045800 | 0.101  | 0.163  |
| ETHYLENE GLYCOL MONOPHENYL ETHER    | 0.003800 | 0.005  | 0.005  |
| ETHYLENE GLYCOL MONOPROPYL ETHER    | 0.047400 | 0.182  | 0.242  |
| ETHYLENE THIOUREA                   | 0.000008 | 0.001  | 0.002  |
| ETHYLPHENOL, 3-                     | 0.005600 | 0.021  | 0.020  |
| FLUOROACETIC ACID, SODIUM SALT*     | 0.000300 | 0.750  | 1.000  |
| FORMALDEHYDE                        | 0.018700 | 0.533  | 1.000  |
| FORMAMIDE                           | 0.065600 | 0.092  | 0.170  |
| FORMIC ACID                         | 0.038900 | 0.078  | 0.225  |
| FUMARIC ACID                        | 0.092200 | 0.000  | 0.000  |
| GLUTARIC ACID                       | 0.001100 | 0.000  | 0.000  |
| GLYCERIN (GLYCEROL)                 | 0.000700 | 0.000  | 0.000  |
| GLYCINAMIDE                         | 0.008200 | 0.019  | 0.089  |
| GLYOXYLIC ACID                      | 0.006200 | 0.001  | 0.002  |
| GLYPHOSATE                          | 0.000400 | 0.005  | 0.009  |
| GUANIDINE, NITROSO*                 | 0.048900 | 0.000  | 0.001  |
| GUTHION                             | 0.000093 | 0.001  | 0.001  |
| GYLCIDOL                            | 0.050100 | 0.024  | 0.032  |
| HEXAMETHYLENE 1, 6 DIISOCYANATE     | 0.014800 | 0.005  | 0.007  |
| HEXAMETHYLPHOSPHORAMIDE             | 0.000000 | 0.000  | 0.000  |
| HEXANOIC ACID                       | 0.058900 | 0.061  | 0.075  |
| HYDRAZINE                           | 0.037000 | 0.190  | 0.332  |
| HYDROCYANIC ACID (M)                | 0.025800 | 0.999  | 0.999  |
| HYDROQUINONE                        | 0.000080 | 0.000  | 0.000  |
| HYDROXY-(2)-PROPIONITRILE           | 0.004200 | 0.003  | 0.004  |
| HYDROXYPROPIONALDEHYDE              | 0.013200 | 0.066  | 0.102  |
| INDENO(1,2,3- <i>cd</i> )-PYRENE    | 0.000000 | 0.000  | 0.000  |
| LEAD ACETATE (M)                    | 0.000041 | 0.062  | 0.062  |
| LEAD SUBACTEATE (M)                 | 0.000800 | 0.000  | 0.000  |
| LEUCINE (M)                         | 0.030000 | 0.469  | 0.469  |
| MALATHION (M)                       | 0.006700 | 0.060  | 0.060  |
| MALEIC ACID                         | 0.000800 | 0.000  | 0.000  |
| MALEIC ANHYDRIDE                    | 0.012200 | 0.027  | 0.043  |
| MALIC ACID (hydroxybutaneic)        | 0.000000 | 0.000  | 0.000  |
| MESITYL OXIDE (M)                   | 0.019500 | 0.999  | 0.999  |
| METHANE SULFONIC ACID*              | 0.026700 | 0.000  | 0.001  |

TABLE 1 OF APPENDIX J.—FM VALUES FOR HENRY'S LAW CONSTANTS AT 25°C LESS THAN 0.1 (Y/X) ATMOSPHERES PER MOLE FRACTION—Continued

[Use with Section 2.1]

| Compound                                    | Y/X      | Fm 25D | Fm 305 |
|---|----------|--------|--------|
| METHOMYL                                    | 0.045100 | 0.008  | 0.013  |
| METHOXYPHENOL P                             | 0.017200 | 0.003  | 0.003  |
| METHYL HYDRAZINE                            | 0.024800 | 0.082  | 0.155  |
| METHYL METHANESULFONATE                     | 0.000039 | 0.001  | 0.001  |
| METHYL PARATHION                            | 0.000007 | 0.012  | 0.020  |
| METHYL SULFURIC ACID (M)                    | 0.031200 | 0.794  | 0.794  |
| METHYL THIOPHENOL 4                         | 0.024400 | 0.885  | 1.000  |
| METHYL-2-METHOXYAZIRIDINE 1                 | 0.024200 | 0.727  | 0.998  |
| METHYLENE DIPHENYL DIISOCYANATE             | 0.002700 | 0.010  | 0.011  |
| METHYLENE DIPHENYLAMINE (MDA)               | 0.001600 | 0.002  | 0.002  |
| METHYLENE-BIS (2-CHLOROANILINE), 4, 4       | 0.018700 | 0.008  | 0.008  |
| METHYLENEDIANILINE 4, 4                     | 0.028500 | 0.001  | 0.001  |
| METHYLETHYLIDENE BISPHENOL, 4, 4'           | 0.000001 | 0.000  | 0.000  |
| METHYLFURFURAL 5                            | 0.012200 | 0.859  | 1.000  |
| METHYLIMINOACETIC ACID                      | 0.055600 | 0.002  | 0.004  |
| MONOMETHYL FORMANIDE                        | 0.000054 | 0.003  | 0.005  |
| NABAM                                       | 0.000000 | 0.000  | 0.000  |
| NAPHTHOL, alpha-                            | 0.001400 | 0.004  | 0.004  |
| NAPHTHOL, beta-                             | 0.000800 | 0.003  | 0.003  |
| NAPHTHYLAMINE, alpha-                       | 0.002800 | 0.005  | 0.005  |
| NAPHTHYLAMINE, beta-                        | 0.002000 | 0.004  | 0.004  |
| NEOPENTYL GLYCOL                            | 0.000900 | 0.004  | 0.005  |
| NIACIN (M)                                  | 0.034200 | 0.606  | 0.606  |
| NIACINAMIDE(M)                              | 0.067800 | 0.623  | 0.623  |
| NITROANILINE (-o) (M)                       | 0.027800 | 0.351  | 0.351  |
| NITROGLYCERIN                               | 0.000000 | 0.013  | 0.047  |
| NITROPHENOL, 2-                             | 0.006500 | 0.011  | 0.016  |
| NITROPHENOL, 4-                             | 0.000073 | 0.001  | 0.001  |
| NITROSODIMETHYLAMINE N                      | 0.048800 | 0.103  | 0.285  |
| NITROSODI-n-PROPYLAMINE N                   | 0.025200 | 0.088  | 0.105  |
| NITROSODIPHENYLAMINE N*                     | 0.046000 | 0.026  | 0.025  |
| NITROSOMORPHOLINE                           | 0.004700 | 0.011  | 0.019  |
| NITROSO-N-METHYLUREA N                      | 0.001400 | 0.015  | 0.037  |
| ODDIETH . O2ETH . THIOETH . PHOSPHORATE (M) | 0.000000 | 0.096  | 0.096  |
| OXALIC ACID                                 | 0.000200 | 0.010  | 0.028  |
| PARATHION                                   | 0.034000 | 0.001  | 0.001  |
| PENTAERYTHRITOL                             | 0.000021 | 0.000  | 0.000  |
| PHENACETIN (M)                              | 0.012400 | 0.135  | 0.135  |
| PHENOL                                      | 0.072200 | 0.036  | 0.035  |
| PHENYL MERCURIC ACETATE (M)                 | 0.000700 | 0.057  | 0.057  |
| PHENYLACETIC ACID (M)                       | 0.045600 | 0.385  | 0.385  |
| PHENYLENE DIAMINE (-m)                      | 0.000600 | 0.000  | 0.000  |
| PHENYLENE DIAMINE (-o)                      | 0.000600 | 0.001  | 0.002  |
| PHENYLENE DIAMINE (-p)                      | 0.000070 | 0.001  | 0.001  |
| PHORATE (M)                                 | 0.024300 | 0.095  | 0.095  |
| PHTHALIC ANHYDRIDE                          | 0.044100 | 0.016  | 0.019  |
| PROPANE SULTONE, 1, 3-                      | 0.000500 | 0.001  | 0.002  |
| PROPANONAL (methylglyoxal)                  | 0.001700 | 0.161  | 0.242  |
| PROPIOLACTONE b                             | 0.006400 | 0.199  | 0.304  |
| PROPORUR (Baygon)                           | 0.003200 | 0.004  | 0.004  |
| PROPYLENE GLYCOL                            | 0.083300 | 0.005  | 0.008  |
| PYRIDINIUM BROMIDE (M)                      | 0.091700 | 0.060  | 0.060  |
| PYRUVIC ACID                                | 0.000200 | 0.003  | 0.005  |
| QUINOLINE                                   | 0.015000 | 0.002  | 0.002  |
| QUINONE                                     | 0.057700 | 0.868  | 1.000  |
| RESORCINOL                                  | 0.001000 | 0.000  | 0.000  |
| SIMAZINE (M)                                | 0.000045 | 0.124  | 0.124  |
| SODIUM ACETATE                              | 0.000200 | 0.042  | 0.079  |
| SODIUM ACRYLATE                             | 0.076100 | 0.073  | 0.108  |
| SODIUM FORMATE                              | 0.000094 | 0.356  | 0.988  |
| STRYCHNIDIN-10-ONE, 2, 3-DIMETHOXY(M)       | 0.000800 | 0.028  | 0.028  |
| STRYCHNINE (M)                              | 0.000002 | 0.058  | 0.058  |
| SUCCINIC ACID                               | 0.000097 | 0.000  | 0.001  |
| SUCCINIMIDE *                               | 0.001800 | 0.000  | 0.001  |
| SULFANILIC ACID (M)                         | 0.088900 | 0.138  | 0.138  |
| TEREPHTHALIC ACID                           | 0.000600 | 0.001  | 0.001  |
| TETRAETHYLDITHIOPYROPHOSPHATE               | 0.00040  | .....  | .....  |
| TETRAETHYLENE GLYCOL MONOMETHYL ETH         | 0.000200 | 0.000  | 0.001  |
| TETRAETHYLENE PENTAMINE                     | 0.000000 | 0.000  | 0.000  |



TABLE 1 OF APPENDIX J.—FM VALUES FOR HENRY'S LAW CONSTANTS AT 25°C LESS THAN 0.1 (Y/X) ATMOSPHERES PER MOLE FRACTION—Continued

[Use with Section 2.1]

| Compound                                  | Y/X      | Fm 25D | Fm 305 |
|---|----------|--------|--------|
| TETRAETHYLENE PENTAMINE .....             | 0.000000 | 0.000  | 0.000  |
| TETRAHYDRO 3-FURANOL .....                | 0.034400 | 0.095  | 0.134  |
| THIOFANOX (M) .....                       | 0.000500 | 0.116  | 0.116  |
| THIOSEMICARBAZIDE* .....                  | 0.003300 | 0.000  | 0.000  |
| THIOUREA, 1- (o-CHLOROPHENYL)- .....      | 0.000001 | 0.000  | 0.001  |
| TOLUENE DIAMINE (2, 4) .....              | 0.000070 | 0.001  | 0.001  |
| TOLUENE DIISOCYANATE (2, 4) .....         | 0.009200 | 0.000  | 0.000  |
| TOLUENEDIAMINE (2, 6) .....               | 0.000001 | 0.000  | 0.000  |
| TOLUENEDIAMINE (3, 4) .....               | 0.000200 | 0.002  | 0.002  |
| TOLUIC ACID (para-) .....                 | 0.000300 | 0.011  | 0.012  |
| TOLUIDINE m .....                         | 0.089400 | 0.123  | 0.118  |
| TRICHLORO (1, 1, 2) TRIFLUOROETHANE ..... | 0.000008 | 1.042  | 1.000  |
| TRIETHANOLAMINE .....                     | 0.000008 | 0.000  | 0.000  |
| TRIETHYLENE GLYCOL DIMETHYL ETHER .....   | 0.002600 | 0.017  | 0.025  |
| TRIETHYLENE GLYCOL MONOMETHYL ETHER ..... | 0.001900 | 0.004  | 0.005  |
| TRIETHYLENE TETRAMINE .....               | 0.000000 | 0.000  | 0.000  |
| TRIPROPYLENE GLYCOL .....                 | 0.005300 | 0.004  | 0.005  |
| WARFARIN .....                            | 0.000000 | 0.000  | 0.000  |

\* Molecular structure only approximate.

(M) fraction measured (fm) estimated from Mwt correlation.

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION

| Compound                          | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS      |
|-----------------------------------|-------|-------|-------|-----------------|----------|
| 1 BROMO 2 CHLORO 2 BUTENE .....   | 0.990 | 0.786 | 1.000 | 0.761           |          |
| 1 BUTYENE .....                   | 0.990 | 1.172 | 1.000 | 0.872           |          |
| 1 ETHYL 4 METHYLBENZENE .....     | 0.990 | 1.219 | 1.000 | 0.748           |          |
| 1 HEPTANOL .....                  | 0.946 | 0.525 | 0.564 | 0.186           |          |
| 1 HEPTYNE .....                   | 0.990 | 1.138 | 1.000 | 0.980           |          |
| 1 HEXYNE .....                    | 0.990 | 1.145 | 1.000 | 0.924           |          |
| 1 ISOCYANO 3-METHYLBENZENE .....  | 0.990 | 0.870 | 0.913 | 0.210           |          |
| 1 ISOPROPYL 4 METHYLBENZENE ..... | 0.990 | 1.193 | 1.000 | 0.804           |          |
| 1 METHYLCYCLOHEXENE .....         | 0.990 | 1.138 | 1.000 | 0.980           |          |
| 1 METHYLNAPHTHALENE .....         | 0.990 | 1.237 | 1.000 | 0.384           |          |
| 1 NONYNE .....                    | 0.990 | 1.128 | 1.000 | 0.980           |          |
| 1 OCTENE .....                    | 0.990 | 1.112 | 1.000 | 0.980           |          |
| 1 OCTYNE .....                    | 0.990 | 1.132 | 1.000 | 0.980           |          |
| 1 PENTYNE .....                   | 0.990 | 1.156 | 1.000 | 0.885           |          |
| 1,1 DIETHOXYETHANE .....          | 0.985 | 0.810 | 0.996 | 0.320           |          |
| 1,1,3 TRIMETHYLCYCLOPENTANE ..... | 0.990 | 1.124 | 1.000 | 0.980           |          |
| 1,1-DIFLUOROETHANE .....          | 0.990 | 1.077 | 1.000 | 0.876           |          |
| 1,2 DIETHOXYETHANE .....          | 0.932 | 0.762 | 0.999 | 0.309           |          |
| 1,2,4,5 TETRAMETHYLBENZENE .....  | 0.990 | 1.194 | 1.000 | 0.887           |          |
| 1,3-DIOXOLANE .....               | 0.642 | 0.764 | 1.000 | 0.232           | 646-06-0 |
| 1,4 PENTADIENE .....              | 0.990 | 1.176 | 1.000 | 0.980           |          |
| 1,5 HEXADIENE .....               | 0.990 | 1.155 | 1.000 | 0.980           |          |
| 1-NITROPROPANE .....              | 0.966 | 0.522 | 0.982 | 0.374           |          |
| 1-PENTANOL .....                  | 0.990 | 0.708 | 0.807 | 0.579           |          |
| 1-PENTENE .....                   | 0.990 | 1.124 | 1.000 | 0.980           |          |
| 1-PROPOXY 2-PROPANOL .....        | 0.430 | 0.134 | 0.167 | 0.070           |          |
| 2 BUTEN 1 OL .....                | 0.207 | 0.703 | 0.801 | 0.095           |          |
| 2 HEPTANONE .....                 | 0.990 | 0.955 | 0.991 | 0.356           |          |
| 2 METHYL 1 BUTANOL .....          | 0.797 | 0.721 | 0.807 | 0.201           |          |
| 2 METHYL 2 BUTENE .....           | 0.990 | 1.143 | 1.000 | 0.980           |          |
| 2 METHYL 2 PENTANOL .....         | 0.959 | 0.806 | 0.869 | 0.257           |          |
| 2 METHYL 3 PENTANOL .....         | 0.989 | 0.539 | 0.565 | 0.241           |          |
| 2 METHYLHEXANE C7H16 .....        | 0.990 | 1.099 | 1.000 | 0.980           |          |
| 2 METHYLNAPHTHALENE .....         | 0.990 | 1.237 | 1.000 | 0.449           |          |
| 2 NONANONE .....                  | 0.990 | 0.959 | 0.970 | 0.441           |          |
| 2 OCTANONE .....                  | 0.990 | 0.961 | 0.983 | 0.350           |          |
| 2 PENTANONE .....                 | 0.942 | 0.919 | 0.998 | 0.350           |          |
| 2 PENTENE .....                   | 0.990 | 1.131 | 1.000 | 0.980           |          |
| 2 PROPYLBENZENE .....             | 0.990 | 1.198 | 1.000 | 0.582           |          |
| 2 UNDECANONE .....                | 0.990 | 0.927 | 0.922 | 0.495           |          |
| 2-(1-METHOXY)-1-PROPANOL .....    | 0.648 | 0.202 | 0.251 | 0.093           |          |
| 2,2 DIMETHYL PROPANOIC ACID ..... | 0.131 | 0.296 | 0.376 | 0.074           |          |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                            | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|-------------------------------------|-------|-------|-------|-----------------|-----------|
| 2,2 DIMETHYLBUTANE C6H14            | 0.990 | 1.108 | 1.000 | 0.901           |           |
| 2,2 DIMETHYLPENTANE                 | 0.990 | 1.106 | 1.000 | 0.980           |           |
| 2,2,5 TRIMETHYLHEXANE C9H20         | 0.990 | 1.114 | 1.000 | 0.980           |           |
| 2,3 DIMETHYL 1,3 BUTADIENE          | 0.990 | 1.168 | 1.000 | 0.942           |           |
| 2,3 DIMETHYLBUTANE C6H14            | 0.990 | 1.115 | 1.000 | 0.980           |           |
| 2,3 DIMETHYLBUTANOL                 | 0.978 | 0.648 | 0.694 | 0.259           |           |
| 2,3 DIMETHYLPENTANE C7H16           | 0.990 | 1.112 | 1.000 | 0.980           |           |
| 2,3,4 TRIMETHYLPENTANE C8H18        | 0.990 | 1.121 | 1.000 | 0.980           |           |
| 2,3-DIMETHYLPYRIDINE                | 0.048 | 1.048 | 1.000 | 0.110           |           |
| 2,4 DIMETHYLPENTANE C7H16           | 0.990 | 1.112 | 1.000 | 0.980           |           |
| 2,4,5 T                             | ..... | 0.024 | 0.028 | 0.000           | 93-76-5   |
| 2,4-DIMETHYLPYRIDINE                | 0.044 | 1.048 | 1.000 | 0.105           |           |
| 2,5-DIMETHYLPYRIDINE                | 0.055 | 1.048 | 1.000 | 0.122           |           |
| 2,6-DIMETHYL2,5-HEPTADIEN4-ONE      | 0.990 | 0.906 | 0.882 | 0.354           |           |
| 2,6-DIMETHYL2,5-HEPTADIEN 4-ONE     | 0.990 | 0.682 | 0.649 | 0.278           |           |
| 2,6-DIMETHYLPYRIDINE                | 0.067 | 1.048 | 1.000 | 0.137           |           |
| 2-CHLORO 2-METHYLBUTANE             | 0.990 | 1.078 | 1.000 | 0.726           |           |
| 2-ETHYL 3-METHOXYPYRAZINE           | 0.990 | 0.039 | 0.050 | 0.151           |           |
| 2-ETHYLPYRAZINE                     | 0.746 | 0.452 | 0.527 | 0.070           |           |
| 2-ETHYLPYRIDINE                     | 0.080 | 1.041 | 1.000 | 0.141           |           |
| 2-FLUOROPROPANE                     | 0.990 | 1.099 | 1.000 | 0.980           |           |
| 2-ISOBUTYL 3-METHOXYPYRAZINE        | 0.990 | 0.044 | 0.057 | 0.256           |           |
| 2-ISOBUTYLPYRAZINE                  | 0.969 | 0.362 | 0.395 | 0.096           |           |
| 2-METHYL PENTANE C6H14              | 0.990 | 1.100 | 1.000 | 0.899           |           |
| 2-METHYLPYRAZINE                    | 0.626 | 0.505 | 0.613 | 0.068           |           |
| 2-PENTANOL                          | 0.810 | 0.721 | 0.807 | 0.205           |           |
| 3 METHYL 1 BUTENE                   | 0.990 | 1.143 | 1.000 | 0.980           |           |
| 3 METHYL PYRIDINE                   | 0.630 | 0.685 | 0.663 | 0.131           |           |
| 3 METHYLHEPTANE C8H18               | 0.990 | 1.098 | 1.000 | 0.980           |           |
| 3 METHYLHEXANE C7H16                | 0.990 | 1.099 | 1.000 | 0.980           |           |
| 3,3 DIMETHYLPENTANE C7H16           | 0.990 | 1.106 | 1.000 | 0.980           |           |
| 3,4-DIMETHYLPYRIDINE                | 0.025 | 1.048 | 1.000 | 0.083           |           |
| 3,5-DIMETHYLPYRIDINE                | 0.044 | 1.048 | 1.000 | 0.105           |           |
| 3-ETHYLPYRIDINE                     | 0.080 | 1.041 | 1.000 | 0.141           |           |
| 3-HEXANOL                           | 0.990 | 0.638 | 0.694 | 0.294           |           |
| 3-PENTEN-2-OL                       | 0.860 | 0.610 | 0.656 | 0.230           |           |
| 4 METHYL 1 PENTENE                  | 0.990 | 1.134 | 1.000 | 0.980           |           |
| 4 METHYL 2 PENTANOL                 | 0.990 | 0.539 | 0.565 | 0.264           |           |
| 4 METHYL 2 PENTANONE                | 0.385 | 0.923 | 0.968 | 0.145           |           |
| 4 METHYLOCTANE C9H20                | 0.990 | 1.098 | 1.000 | 0.980           |           |
| 4-ETHYLPYRIDINE                     | 0.064 | 1.041 | 1.000 | 0.123           |           |
| 4-METHYLPYRIDINE                    | 0.990 | 1.033 | 1.000 | 0.109           |           |
| 5 METHOXY 2 PENTANONE               | 0.798 | 0.327 | 0.382 | 0.142           |           |
| ACENAPHTHENE                        | 0.990 | 1.111 | 0.899 | 0.804           | 83-32-9   |
| ACENAPHTHYLENE                      | 0.990 | 1.094 | 0.868 | 0.312           | 208-96-8  |
| ACETAL                              | 0.990 | 0.813 | 1.000 | 0.432           |           |
| ACETALDEHYDE                        | 0.953 | 0.724 | 1.000 | 0.485           | 75-07-0   |
| ACETATE (M)                         | 0.990 | 0.558 | 0.558 | 0.794           |           |
| ACETIC ACID                         | 0.066 | 0.101 | 0.189 | 0.120           | 64-19-7   |
| ACETIC ANHYDRIDE                    | 0.524 | 0.165 | 0.262 | 0.214           | 108-24-7  |
| ACETONE                             | 0.843 | 0.827 | 0.997 | 0.261           | 67-64-1   |
| ACETONITRILE                        | 0.641 | 0.778 | 0.989 | 0.359           | 75-05-8   |
| ACETOPHENONE                        | 0.735 | 0.334 | 0.314 | 0.137           | 96-86-2   |
| ACETYL CHLORIDE                     | 0.990 | 0.923 | 1.000 | 0.531           | 79-36-5   |
| ACETYL DIETHYLMALONATE              | 0.978 | 0.018 | 0.025 | 0.156           |           |
| ACETYLENE                           | 0.990 | 1.280 | 1.000 | 0.711           | 74-86-2   |
| ACETYLFURAN 2 *                     | 0.990 | 0.365 | 0.423 | 0.382           | 1192-62-7 |
| ACETYLMETHYLPHTHALATE 4             | 0.990 | 0.036 | 0.048 | 0.127           |           |
| ACETYLPYRIDINE 3                    | 0.990 | 0.927 | 1.000 | 0.980           | 1122-54-9 |
| ACIFLUORFEN                         | 0.990 | 0.198 | 0.223 | 0.601           |           |
| ACROLEIN                            | 0.968 | 0.855 | 1.000 | 0.427           | 107-02-8  |
| ACRYLONITRILE                       | 0.969 | 0.876 | 0.999 | 0.429           | 107-13-1  |
| ADAMANTANE DICHLORIDE               | 0.990 | 1.097 | 0.986 | 0.562           |           |
| AFLATOXINS (M)                      | 0.990 | 0.063 | 0.063 | 0.406           | 1402-68-2 |
| ALDICARB                            | 0.027 | 0.002 | 0.002 | 0.007           | 116-06-3  |
| ALDRIN                              | 0.990 | 0.056 | 0.051 | 0.469           | 509-00-2  |
| ALKYLIMINE CARBOXYLIC ACID N,SUB(M) | 0.848 | 0.125 | 0.125 | 0.111           |           |
| ALLYL ALCOHOL                       | 0.783 | 0.538 | 0.659 | 0.276           | 107-18-6  |
| ALLYL CHLORIDE                      | 0.990 | 1.092 | 1.000 | 0.887           | 107-05-1  |
| ALLYL ETHER, diallyl ether          | 0.990 | 0.974 | 1.000 | 0.663           |           |
| ALPHA METHYL STYRENE                | 0.990 | 1.217 | 1.000 | 0.767           | 98-83-9   |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                                  | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|---|-------|-------|-------|-----------------|-----------|
| ALPHA METHYL STYRENE DIMERS .....         | 0.990 | 1.186 | 0.975 | 0.855           |           |
| alpha-CHLORO-beta-METHYLNAPHTHALENE ..... | 0.990 | 1.197 | 1.000 | 0.828           | 86-52-2   |
| ALPHA-HYDROXYACETALDEHYDE .....           | 0.990 | 0.031 | 0.059 | 0.515           |           |
| ALPHA-HYDROXYADIPIMIDE (M) .....          | 0.925 | 0.144 | 0.144 | 0.135           |           |
| AMINO-2-CHLOROTOLUENE 4 .....             | 0.990 | 0.020 | 0.020 | 0.790           |           |
| AMINO-3-CHLORO-5-PHENYLCYCLOHEXA(M) ..... | 0.622 | 0.143 | 0.143 | 0.086           |           |
| AMINO-4-CHLORO-6-CYANOPYRIDINE 2(M) ..... | 0.990 | 0.148 | 0.148 | 0.411           |           |
| AMINO-4'-CHLOROBIPHENYL 4(M) .....        | 0.990 | 0.123 | 0.123 | 0.980           |           |
| AMINO-4-CHLOROPYRIDINE 2 (M) .....        | 0.990 | 0.514 | 0.514 | 0.710           | 1072-98-6 |
| AMINO-4-NITROBENZYL ALCOHOL 2 (M) .....   | 0.742 | 0.149 | 0.149 | 0.102           |           |
| AMINO-4-NITROTOLUENE 2 .....              | 0.990 | 0.000 | 0.001 | 0.802           | 99-55-8   |
| AMINO-5-CHLOROPYRIDINE 2 (M) .....        | 0.990 | 0.514 | 0.514 | 0.384           | 1072-98-6 |
| AMINOBENZOIC ACID (-p) (M) .....          | 0.624 | 0.368 | 0.368 | 0.086           | 150-13-0  |
| AMINOCYCLOHEXANE .....                    | 0.934 | 0.929 | 0.996 | 0.416           | 108-91-8  |
| AMINOMETHYL-3-ISOXAZOLOL 5 (M) .....      | 0.990 | 0.760 | 0.760 | 0.287           | 2763-96-4 |
| AMINOPHENOL(-o) .....                     | 0.641 | 0.034 | 0.039 | 0.083           | 95-55-6   |
| AMINOPHENOL(-p) .....                     | 0.265 | 0.001 | 0.001 | 0.180           | 101-80-4  |
| AMINO-p'-METHYLAZOBENZENE P (M) .....     | 0.990 | 0.119 | 0.119 | 0.852           |           |
| AMINOPROPIONITRILE 3 (M) .....            | 0.834 | 0.999 | 0.999 | 0.163           | 151-18-8  |
| AMITROLE (M) .....                        | 0.618 | 0.999 | 0.999 | 0.085           | 61-82-5   |
| AMMONIA .....                             | 0.990 | 0.520 | 1.000 | 0.732           | 7664-41-7 |
| AMPHETAMINE(M) .....                      | 0.990 | 0.401 | 0.401 | 0.323           | 60-15-1   |
| AMYL ACETATE(-n) .....                    | 0.990 | 0.426 | 0.504 | 0.462           | 628-63-7  |
| ANETHOLE (M) .....                        | 0.990 | 0.180 | 0.180 | 0.406           | 104-46-1  |
| ANISOLE .....                             | 0.990 | 1.036 | 1.000 | 0.731           | 100-66-3  |
| ANTHRACENE .....                          | 0.990 | 0.109 | 0.087 | 0.513           | 120-12-7  |
| ARAMITE (M) .....                         | 0.990 | 0.058 | 0.058 | 0.406           | 140-57-8  |
| AURAMINE (M) .....                        | 0.990 | 0.091 | 0.091 | 0.980           | 492-80-8  |
| AZASERINE (M) .....                       | 0.986 | 0.138 | 0.138 | 0.206           | 115-02-6  |
| AZEPINE (M) .....                         | 0.990 | 0.058 | 0.058 | 0.817           | 111-49-9  |
| AZIRIDINE ethyleneimine .....             | 0.990 | 0.628 | 0.867 | 0.685           | 151-56-4  |
| BENXENEDICARBOXYLIC ACID DIHEPTYL .....   | 0.990 | 0.113 | 0.119 | 0.667           |           |
| BENZ(c)ACRIDINE (M) .....                 | 0.990 | 0.110 | 0.110 | 0.853           | 225-51-4  |
| BENZAL CHLORIDE .....                     | 0.990 | 1.159 | 0.996 | 0.798           | 98-87-3   |
| BENZALDEHYDE .....                        | 0.980 | 0.516 | 0.490 | 0.283           | 100-52-7  |
| BENZALKONIUM CHLORIDE (M) .....           | 0.408 | 0.129 | 0.129 | 0.065           |           |
| BENZEN SULFONATE (M) .....                | 0.990 | 0.642 | 0.642 | 0.894           |           |
| BENZENE .....                             | 0.990 | 1.227 | 1.000 | 0.797           | 71-43-2   |
| BENZETHONIUM CHLORIDE (M) .....           | 0.956 | 0.001 | 0.001 | 0.140           | 121-54-0  |
| BENZIDINE DIHYDROCHLORIDE(M) .....        | 0.990 | 0.096 | 0.096 | 0.980           | 531-85-1  |
| BENZO(B)FLUORANTHENE .....                | 0.990 | 1.219 | 0.962 | 0.135           | 205-99-2  |
| BENZO(j)FLUORANTHENE (M) .....            | 0.990 | 0.099 | 0.099 | 0.853           | 205-82-3  |
| BENZODIOXANE-1,3 (M) .....                | 0.668 | 0.108 | 0.108 | 0.093           |           |
| BENZOFUORANTHENE,3,4-(M) .....            | 0.990 | 0.099 | 0.099 | 0.853           | 205-99-2  |
| BENZOFURAN 2,3 .....                      | 0.990 | 1.061 | 0.988 | 0.374           |           |
| BENZOIC ACID, 4 METHYL .....              | 0.642 | 0.102 | 0.108 | 0.103           |           |
| BENZONITRILE .....                        | 0.990 | 0.397 | 0.373 | 0.170           | 100-47-0  |
| BENZOPHENONE .....                        | 0.990 | 0.052 | 0.046 | 0.834           | 119-61-9  |
| BENZOPYRENE 3,4 (M) .....                 | 0.990 | 0.099 | 0.099 | 0.318           | 50-32-8   |
| BENZOQUINONE,p-(M) .....                  | 0.990 | 0.862 | 0.862 | 0.794           | 106-51-4  |
| BENZOTHIAZOLE * .....                     | 0.990 | 0.059 | 0.060 | 0.341           | 95-16-9   |
| BENZOTRICHLORIDE .....                    | 0.990 | 1.069 | 0.958 | 0.558           | 98-07-7   |
| BENZOYL CHLORIDE .....                    | 0.990 | 1.132 | 0.979 | 0.468           | 98-88-4   |
| BENZYL CHLORIDE .....                     | 0.990 | 1.164 | 1.000 | 0.415           | 100-44-7  |
| BENZYL METHYL ETHER .....                 | 0.990 | 1.047 | 1.000 | 0.587           | 538-86-3  |
| BHC,alpha- .....                          | 0.990 | 1.063 | 1.000 | 0.729           | 319-84-6  |
| BHC,beta- .....                           | 0.990 | 1.063 | 1.000 | 0.854           | 319-85-7  |
| BHC,delta- .....                          | 0.990 | 1.063 | 1.000 | 0.588           | 319-86-8  |
| BICYCLO(4,2,0) OCTA 1.3.5 TRIENE .....    | 0.990 | 1.222 | 1.000 | 0.759           |           |
| BICYCLO[2.2.1]-2,5-HEPTADIENE DI(M) ..... | 0.990 | 0.146 | 0.146 | 0.980           |           |
| BIPHENYL .....                            | 0.990 | 1.074 | 0.864 | 0.445           | 92-52-4   |
| BIS (2-CHLOROETHOXY) METHANE .....        | 0.282 | 0.170 | 0.196 | 0.067           | 111-91-1  |
| BIS(1,1,2,2-TETRACHLOROPROPYL) ETHE ..... | 0.990 | 0.960 | 1.000 | 0.980           |           |
| BIS(2-CHLOROETHYL)ETHER .....             | 0.656 | 0.806 | 0.858 | 0.162           | 111-44-4  |
| BIS(2-CHLOROISOPROPYL)ETHER .....         | 0.990 | 0.948 | 0.972 | 0.310           | 108-60-1  |
| BIS(CHLOROMETHYL)ETHER .....              | 0.975 | 0.888 | 0.999 | 0.459           | 542-88-1  |
| BISPENOL(A) .....                         | 0.990 | 0.011 | 0.011 | 0.665           | 80-05-7   |
| BROMACIL .....                            | 0.990 | 0.582 | 1.000 | 0.980           |           |
| BROMO-(1)-CHLOROETHANE-2 .....            | 0.990 | 0.711 | 1.000 | 0.995           | 107-04-0  |
| BROMO-3-CHLOROBUTADIENE 2 .....           | 0.990 | 0.803 | 1.000 | 0.820           |           |
| BROMO-4-CHLORO-6-CYANOBENZYL ALC(M) ..... | 0.941 | 0.131 | 0.131 | 0.136           |           |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                             | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|--------------------------------------|-------|-------|-------|-----------------|-----------|
| BROMO-4-CHLOROCYCLOHEXANE 1          | 0.990 | 0.819 | 0.986 | 0.980           |           |
| BROMO-4-CYANOMETHYL BENZOATE 2 (M)   | 0.990 | 0.105 | 0.105 | 0.980           |           |
| BROMO-4-CYANOMETHYL BENZOATE 3 (M)   | 0.990 | 0.105 | 0.105 | 0.885           |           |
| BROMOACETONE                         | 0.520 | 0.356 | 0.590 | 0.145           | 598-31-2  |
| BROMOBENZENE                         | 0.990 | 1.182 | 1.000 | 0.745           | 108-86-1  |
| BROMOBENZYL ALCOHOL-(m)              | 0.371 | 0.012 | 0.015 | 0.083           | 15852-73- |
| BROMOBENZYL ALCOHOL-(o)              | 0.371 | 0.012 | 0.015 | 0.083           | 18982-34- |
| BROMOBENZYL ALCOHOL-(p)              | 0.371 | 0.012 | 0.015 | 0.083           | 873-75-6  |
| BROMOCHLOROBENZENE P                 | 0.990 | 0.870 | 1.000 | 0.980           | 106-39-8  |
| BROMOCHLOROBENZYL ALCOHOL            | 0.420 | 0.007 | 0.009 | 0.107           |           |
| BROMOCHLOROMETHANE                   | 0.990 | 1.017 | 1.000 | 0.992           | 74-97-5   |
| BROMODICHLOROMETHANE                 | 0.990 | 0.735 | 1.000 | 0.980           | 75-27-4   |
| BROMOETHYL ACETATE                   | 0.911 | 0.470 | 0.801 | 0.458           | 927-68-4  |
| BROMOETHYLENE                        | 0.990 | 0.629 | 1.000 | 0.990           | 543-60-2  |
| BROMOFORM                            | 0.990 | 0.480 | 0.998 | 0.494           | 75-25-2   |
| BROMOMETHANE                         | 0.990 | 0.539 | 1.000 | 0.852           | 74-83-9   |
| BROMOPHENYL PHENYL ETHER,4-          | 0.990 | 0.240 | 0.265 | 0.269           | 101-55-3  |
| BROMOPROPIONITRILE 3 (M)             | 0.990 | 0.422 | 0.422 | 0.856           | 2417-90-5 |
| BROMOTOLUENE 4                       | 0.990 | 1.164 | 1.000 | 0.676           | 106-38-7  |
| BROMOURACIL,5-(M)                    | 0.990 | 0.130 | 0.130 | 0.980           | 51-20-7   |
| BUTADIENE-(1,3)                      | 0.990 | 1.187 | 1.000 | 0.979           | 106-99-0  |
| BUTANE                               | 0.990 | 1.080 | 1.000 | 0.980           | 106-97-8  |
| BUTANEDINITRILE                      | 0.990 | 0.007 | 0.009 | 0.182           | 110-61-2  |
| BUTANENITRILE (M)                    | 0.521 | 0.999 | 0.999 | 0.266           | 109-74-0  |
| BUTANOL ISO                          | 0.821 | 0.647 | 0.756 | 0.068           | 78-83-1   |
| BUTANOL(S)                           | 0.846 | 0.502 | 0.600 | 0.253           | 78-92-2   |
| BUTANOL-1                            | 0.818 | 0.502 | 0.600 | 0.177           | 71-36-3   |
| BUTENE                               | 0.990 | 1.131 | 1.000 | 0.980           |           |
| BUTYL ACETATE(-n)                    | 0.990 | 0.808 | 0.995 | 0.368           | 123-86-4  |
| BUTYL ACRYLATE                       | 0.990 | 0.781 | 0.910 | 0.492           | 141-32-2  |
| BUTYL BENZENE                        | 0.990 | 1.181 | 1.000 | 0.980           | 104-51-8  |
| BUTYL BENZYL PHTHALATE               | 0.990 | 0.052 | 0.053 | 0.852           | 85-68-7   |
| BUTYL CARBITOL                       | 0.990 | 0.006 | 0.008 | 0.980           | 112-34-5  |
| BUTYL MERCAPTAN                      | 0.990 | 0.692 | 1.000 | 0.980           |           |
| BUTYL-3-METHOXY PYRAZINE, 2-ISO (M)  | 0.990 | 0.142 | 0.142 | 0.980           | 24683-00- |
| BUTYLAMINE                           | 0.904 | 0.813 | 0.948 | 0.241           | 109-73-9  |
| BUTYLBUTOXY PROPIONATE               | 0.990 | 0.263 | 0.276 | 0.266           |           |
| BUTYLENE GLYCOL-(1,3)                | 0.780 | 0.003 | 0.004 | 0.096           | 107-88-0  |
| BUTYLISOBUTYRATE                     | 0.990 | 0.873 | 1.000 | 0.794           |           |
| BUTYRALDEHYDE                        | 0.989 | 0.861 | 0.992 | 0.490           | 123-72-8  |
| BUTYRALDEHYDE ISO                    | 0.989 | 0.886 | 1.000 | 0.438           | 78-84-2   |
| c10 linear                           | 0.990 | 1.088 | 1.000 | 0.980           |           |
| c11 linear                           | 0.990 | 1.088 | 1.000 | 0.980           |           |
| CACODYLIC ACID (M)                   | 0.983 | 0.354 | 0.354 | 0.219           | 75-60-5   |
| CAMPHENE (M)                         | 0.990 | 0.383 | 0.383 | 0.588           | 79-92-5   |
| CAPTAN                               | 0.990 | 0.007 | 0.008 | 0.196           |           |
| CARBARYL sevin                       | 0.990 | 0.015 | 0.016 | 0.202           | 63-25-2   |
| CARBAZOLE (M)                        | 0.990 | 0.141 | 0.141 | 0.980           | 86-74-8   |
| CARBENDAZIM                          | 0.957 | 0.023 | 0.038 | 0.070           |           |
| CARBON DIOXIDE (M)                   | 0.990 | 0.999 | 0.999 | 0.896           |           |
| CARBON DISULFIDE                     | 0.990 | 0.213 | 1.000 | 0.918           | 75-15-0   |
| CARBON OXYFLUORIDE*                  | 0.990 | 0.884 | 1.000 | 0.993           | 353-50-4  |
| CARBON TETRACHLORIDE                 | 0.990 | 1.027 | 1.000 | 0.900           | 56-23-5   |
| CARBONYL FLUORIDE *                  | 0.658 | 0.884 | 1.000 | 0.358           |           |
| CARBONYL SULFIDE                     | 0.886 | 0.547 | 1.000 | 0.500           |           |
| CHLORAL                              | 0.990 | 0.938 | 1.000 | 0.556           | 302-17-0  |
| CHLORAMBEN                           | 0.962 | 0.545 | 0.633 | 0.229           |           |
| CHLORAMBUCIL                         | 0.957 | 0.031 | 0.031 | 0.101           | 305-03-3  |
| CHLORDANE                            | 0.990 | 0.438 | 0.407 | 0.151           | 57-74-9   |
| CHLORENDIC ANHYDRIDE (M)             | 0.990 | 0.558 | 0.558 | 0.794           | 115-27-5  |
| CHLORINATED TARS (M)                 | 0.990 | 0.050 | 0.050 | 0.343           |           |
| CHLORNAPHAZINE                       | 0.990 | 0.422 | 0.385 | 0.158           |           |
| CHLORO 2 BUTENE,1 trans              | 0.990 | 1.098 | 1.000 | 0.632           |           |
| CHLORO(-p)PHENYLHYDRAZINE(M)         | 0.990 | 0.286 | 0.286 | 0.398           |           |
| CHLORO-1,3-CYCLOPENTADIENE 5         | 0.990 | 1.148 | 1.000 | 0.948           |           |
| CHLORO-2,2-DIBROMOETHANE 1           | 0.990 | 0.569 | 0.919 | 0.526           |           |
| CHLORO-2,3-EPOXYPROPANE,1-(M)        | 0.977 | 0.999 | 0.999 | 0.321           | 106-89-8  |
| CHLORO-2-METHOXYBENZOIC ACID 4 (M)   | 0.990 | 0.132 | 0.132 | 0.722           | 57479-70- |
| CHLORO-2-NITROBENZYL ALCOHOL 4 (M)   | 0.601 | 0.132 | 0.132 | 0.083           | 22996-18- |
| CHLORO-3-NITRO-5-PHENYLCYCLOHEXA (M) | 0.631 | 0.131 | 0.131 | 0.087           |           |
| CHLORO-3-NITROANILINE 4 (M)          | 0.990 | 0.139 | 0.139 | 0.342           | 635-22-3  |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                             | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|--------------------------------------|-------|-------|-------|-----------------|-----------|
| CHLORO-4AMINOCOUMARAN-6 CARBOXYLI(M) | 0.990 | 0.118 | 0.118 | 0.980           |           |
| CHLORO-4-CYANOBENZYL ALCOHOL 2 (M)   | 0.743 | 0.149 | 0.149 | 0.102           |           |
| CHLORO-4-HYDROXYBIPHENYL 3 (M)       | 0.990 | 0.123 | 0.123 | 0.980           | 92-04-6   |
| CHLORO-4-METHOXY-6-AMINOBENZOIC(M)   | 0.990 | 0.125 | 0.125 | 0.449           |           |
| CHLORO-4-METHYL-N-METHYLBENZAMID(M)  | 0.832 | 0.134 | 0.134 | 0.109           |           |
| CHLORO-4-NITROANISOLE 2 (M)          | 0.990 | 0.131 | 0.131 | 0.980           |           |
| CHLORO-4-PHENYLPYRIDINE 2(M)         | 0.839 | 0.130 | 0.130 | 0.110           |           |
| CHLORO-5AMINO3PYRIDINE CARB.ACID (M) | 0.990 | 0.134 | 0.134 | 0.439           |           |
| CHLORO-5-CYANOPHTHALIC ACID 4 (M)    | 0.990 | 0.112 | 0.112 | 0.980           |           |
| CHLORO-5-CYANOTOLUENE 3 (M)          | 0.990 | 0.150 | 0.150 | 0.601           |           |
| CHLORO-5-FLUOROTOLUENE 3             | 0.990 | 1.150 | 1.000 | 0.400           | 443-83-4  |
| CHLORO-5-PHENOXYDIMETHYL PHTHALA(M)  | 0.990 | 0.065 | 0.065 | 0.980           |           |
| CHLOROACETALDEHYDE                   | 0.762 | 0.855 | 0.997 | 0.324           | 107-20-0  |
| CHLOROALLYL ALCOHOL 2                | 0.926 | 0.270 | 0.291 | 0.244           | 5976-47-6 |
| CHLOROANILINE(2)                     | 0.990 | 0.245 | 0.238 | 0.867           | 95-51-2   |
| CHLOROANILINE(3)                     | 0.990 | 0.108 | 0.105 | 0.867           | 108-42-9  |
| CHLOROAZOBENZENE                     | 0.990 | 1.204 | 1.000 | 0.852           |           |
| CHLOROBENZENE                        | 0.990 | 1.157 | 1.000 | 0.728           | 108-90-7  |
| CHLOROBENZENESULFONIC ACID (-p)(M)   | 0.826 | 0.137 | 0.137 | 0.108           | 100-03-8  |
| CHLOROBENZILATE                      | 0.876 | 0.000 | 0.000 | 0.030           | 510-15-6  |
| CHLOROBENZOIC ACID,2                 | 0.629 | 0.083 | 0.089 | 0.105           | 118-91-2  |
| CHLOROBENZOIC ACID,3-                | 0.535 | 0.083 | 0.089 | 0.092           | 535-80-8  |
| CHLOROBENZOIC ACID,4-                | 0.535 | 0.083 | 0.089 | 0.092           | 74-11-3   |
| CHLOROBENZOTRICHLORIDE P             | 0.990 | 1.103 | 1.000 | 0.980           | 5216-25-1 |
| CHLOROBENZOTRIFLUORIDE, P            | 0.990 | 1.131 | 1.000 | 0.980           |           |
| CHLOROBENZYL ALCOHOL-(m)             | 0.852 | 0.035 | 0.033 | 0.074           | 873-63-2  |
| CHLOROBENZYL ALCOHOL-(o)             | 0.275 | 0.058 | 0.056 | 0.074           | 17849-38- |
| CHLOROBENZYL ALCOHOL-(p)             | 0.251 | 0.040 | 0.039 | 0.074           | 873-76-7  |
| CHLOROBIPHENYL (-p)                  | 0.990 | 1.204 | 1.000 | 0.840           | 2051-62-9 |
| CHLOROBUTADIENE,1                    | 0.990 | 1.124 | 1.000 | 0.850           |           |
| CHLOROCOUMARAN 2 (M)                 | 0.990 | 0.135 | 0.135 | 0.832           | 2051-59-4 |
| CHLOROCYANOBENZENE (1,4) (M)         | 0.990 | 0.362 | 0.362 | 0.980           | 873-32-5  |
| CHLOROCYCLOHEXANE                    | 0.990 | 1.081 | 1.000 | 0.980           | 542-18-7  |
| CHLOROCYCLOHEXANOL 2                 | 0.990 | 0.102 | 0.107 | 0.428           | 1561-86-0 |
| CHLOROCYCLOHEXANOL 4                 | 0.990 | 0.102 | 0.107 | 0.587           |           |
| CHLORODIACETYL (M)                   | 0.990 | 0.651 | 0.651 | 0.980           |           |
| CHLORODIMETHYL PHTHALATE 3 (M)       | 0.990 | 0.111 | 0.111 | 0.980           |           |
| CHLORODIPHENYL THIOETHER P (M)       | 0.990 | 0.123 | 0.123 | 0.851           | 7005-72-3 |
| CHLOROETHANE (ethyl chloride)        | 0.990 | 1.046 | 1.000 | 0.901           | 75-00-3   |
| CHLOROETHANOL (ETHYLENE CHLOROXYDR)  | 0.480 | 0.256 | 0.309 | 0.221           | 107-07-3  |
| CHLOROETHYL(2-) VINYL ETHER          | 0.990 | 0.934 | 1.000 | 0.910           | 110-75-8  |
| CHLOROETHYLENE                       | 0.990 | 1.064 | 1.000 | 0.757           |           |
| CHLOROFLUOROBENZENE P                | 0.990 | 1.152 | 1.000 | 0.980           | 352-33-0  |
| CHLOROFLUOROMETHANE *                | 0.355 | 1.075 | 1.000 | 0.980           | 593-70-4  |
| CHLOROFORM                           | 0.990 | 1.023 | 1.000 | 0.775           | 67-66-3   |
| CHLOROHYDROXYPHENYL4 METHYLBENZ(M)   | 0.990 | 0.094 | 0.094 | 0.980           |           |
| CHLOROMETHYL ACETYLENE *             | 0.990 | 1.121 | 1.000 | 0.980           |           |
| CHLOROMETHYL BENZOATE P (M)          | 0.990 | 0.140 | 0.140 | 0.980           | 1126-46-1 |
| CHLOROMETHYL ETHYL KETONE            | 0.990 | 0.873 | 0.935 | 0.697           |           |
| CHLOROMETHYL METHYL ETHER            | 0.937 | 0.840 | 1.000 | 0.494           | 107-30-2  |
| CHLOROMETHYL PHENYL KETONE           | 0.290 | 0.715 | 0.673 | 0.077           | 532-27-4  |
| CHLOROMETHYL PHENYLHYDRAZINE P (M)   | 0.990 | 0.147 | 0.147 | 0.413           |           |
| CHLOROMETHYLAMINOIMINE (M)           | 0.990 | 0.999 | 0.999 | 0.913           |           |
| CHLORONAPHTHALENE,2-                 | 0.990 | 1.177 | 0.980 | 0.870           | 91-58-7   |
| CHLORONITROALKOXYIMINE (M)           | 0.958 | 0.110 | 0.110 | 0.142           |           |
| CHLORONITROBENZENE(-o)               | 0.990 | 0.519 | 0.625 | 0.808           | 88-73-3   |
| CHLORONITROBENZENE, p                | 0.990 | 0.591 | 0.713 | 0.301           |           |
| CHLORO-N-METHYLBENZAMIDE P (M)       | 0.818 | 0.140 | 0.140 | 0.107           |           |
| CHLOROPHENOL-2                       | 0.323 | 0.245 | 0.240 | 0.107           | 95-97-8   |
| CHLOROPHENOL-3                       | 0.635 | 0.057 | 0.057 | 0.078           | 108-43-0  |
| CHLOROPHENYL PHENYL ETHER,4*         | 0.990 | 0.861 | 0.775 | 0.389           | 7005-72-3 |
| CHLOROPHENELETHANOL 1,1              | 0.990 | 0.057 | 0.054 | 0.807           |           |
| CHLOROPHTHALIC ANHYDRIDE 4 (M)       | 0.595 | 0.133 | 0.133 | 0.083           |           |
| CHLORO-p'-METHYLBIPHENYL P (M)       | 0.990 | 0.124 | 0.124 | 0.850           | 1667-11-4 |
| CHLOROPRENE                          | 0.990 | 1.124 | 1.000 | 0.677           | 126-99-8  |
| CHLOROPROPANE-1                      | 0.990 | 1.055 | 1.000 | 0.858           | 540-54-5  |
| CHLOROPROPANE-2                      | 0.990 | 1.050 | 1.000 | 0.867           | 75-29-6   |
| CHLOROPROPENE 3                      | 0.990 | 1.092 | 1.000 | 0.980           | 557-98-2  |
| CHLOROPROPIONITRILE,3-               | 0.359 | 0.580 | 0.622 | 0.111           | 542-76-7  |
| CHLOROPROPYLENE-2                    | 0.990 | 1.090 | 1.000 | 0.980           | 557-98-2  |
| CHLORO-p-XYLENE                      | 0.987 | 1.163 | 1.000 | 0.592           | 104-82-5  |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                            | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|-------------------------------------|-------|-------|-------|-----------------|-----------|
| CHLOROPYRIDINE 2 (M)                | 0.990 | 0.769 | 0.769 | 0.599           | 109-09-1  |
| CHLOROSTYRENE (-4)                  | 0.990 | 1.179 | 1.000 | 0.788           | 1331-28-8 |
| CHLOROTETRAHYDROFURAN 3 (M)         | 0.990 | 0.642 | 0.642 | 0.407           |           |
| CHLOROTHIOPHENOL P *                | 0.990 | 0.893 | 1.000 | 0.980           | 106-54-7  |
| CHLOROTOLUENE-4                     | 0.990 | 1.164 | 1.000 | 0.741           | 106-43-4  |
| CHLOROURACIL,5-(M)                  | 0.990 | 0.138 | 0.138 | 0.980           | 1820-81-1 |
| cis 1,2 DIMETHYLCYCLOHEXANE         | 0.990 | 1.117 | 1.000 | 0.980           |           |
| CITRUS RED #2 (M)                   | 0.990 | 0.071 | 0.071 | 0.853           | 6358-53-8 |
| COPPER PHTHALOCYANINE (M)           | 0.990 | 0.000 | 0.000 | 0.764           | 147-14-8  |
| COUMARAN (M)                        | 0.990 | 0.215 | 0.215 | 0.980           | 91-64-5   |
| CROTONALDEHYDE                      | 0.578 | 0.887 | 0.974 | 0.212           | 470-30-3  |
| CROTONYLENE (2-BUTYNE)              | 0.990 | 1.185 | 1.000 | 0.980           | 503-17-3  |
| CUMENE (isopropylbenzene)           | 0.990 | 1.197 | 1.000 | 0.876           | 98-82-8   |
| CUMENE HYDROPEROXIDE                | 0.987 | 0.478 | 0.464 | 0.204           |           |
| CYANOBENZYL ALCOHOL P *             | 0.147 | 0.002 | 0.002 | 0.070           |           |
| CYANOGEN                            | 0.990 | 0.800 | 1.000 | 0.747           | 460-19-5  |
| CYANOGEN BROMIDE *                  | 0.990 | 0.558 | 1.000 | 0.462           | 506-68-3  |
| CYANOGEN CHLORIDE(M)                | 0.990 | 0.999 | 0.999 | 0.704           | 506-77-4  |
| CYANO GUANIDINE (M)                 | 0.990 | 0.999 | 0.999 | 0.648           | 461-58-5  |
| CYANOMETHYLPHTHALATE 4 (M)          | 0.990 | 0.071 | 0.071 | 0.980           |           |
| CYANOPYRIDINE (-4) *                | 0.990 | 0.118 | 0.124 | 0.980           | 100-48-1  |
| CYANOPYRIDINE 3 *                   | 0.990 | 0.113 | 0.119 | 0.980           | 100-54-9  |
| CYANOTOLUENE 4                      | 0.990 | 0.450 | 0.419 | 0.980           |           |
| CYANURIC ACID (M)                   | 0.491 | 0.505 | 0.505 | 0.072           | 108-80-5  |
| CYCASIN (M)                         | 0.990 | 0.099 | 0.099 | 0.794           | 14901-08- |
| CYCLOHEXADIENE1,4DIONE2,6BIS11DIMET | 0.753 | 0.027 | 0.026 | 0.072           |           |
| CYCLOHEXANE                         | 0.990 | 1.093 | 1.000 | 0.859           | 110-82-7  |
| CYCLOHEXANOL                        | 0.851 | 0.456 | 0.493 | 0.159           |           |
| CYCLOHEXANOL                        | 0.925 | 0.243 | 0.262 | 0.136           | 108-93-0  |
| CYCLOHEXANONE                       | 0.198 | 0.703 | 0.740 | 0.088           | 108-94-1  |
| CYCLOHEXENE                         | 0.990 | 1.136 | 1.000 | 0.980           | 110-83-8  |
| CYCLOHEXENE 1 ONE, 2                | 0.759 | 0.498 | 0.507 | 0.183           |           |
| CYCLOHEXYL ACETATE                  | 0.990 | 0.846 | 0.963 | 0.273           | 622-45-7  |
| CYCLOHEXYL-2,2-DIPHENYLETHYLAMIN(M) | 0.990 | 0.097 | 0.097 | 0.384           |           |
| CYCLOHEXYL-4,6-DINITROPHENOL,2-(M)  | 0.990 | 0.092 | 0.092 | 0.980           | 131-89-5  |
| CYCLOHEXYLAMINE                     | 0.978 | 0.878 | 0.940 | 0.280           | 108-91-8  |
| CYCLOHEXYLCYCLOHEXANONE 4           | 0.990 | 0.732 | 0.707 | 0.727           | 56025-96- |
| CYCLOPENTADIENE                     | 0.990 | 1.198 | 1.000 | 0.980           |           |
| CYCLOPENTADIENE 1,3                 | 0.990 | 1.198 | 1.000 | 0.713           |           |
| CYCLOPENTANE                        | 0.990 | 1.093 | 1.000 | 0.980           |           |
| CYCLOPENTENE                        | 0.990 | 1.144 | 1.000 | 0.979           |           |
| CYCLOPHOSPHAMIDE (M)                | 0.990 | 0.094 | 0.094 | 0.610           | 50-18-0   |
| CYCLOPROPANE C3H6                   | 0.990 | 1.093 | 1.000 | 0.980           |           |
| CYCLOHEXYL o,o-DIMETHYL PHOS.DIT(M) | 0.99  | 0.105 | 0.980 | 0.980           |           |
| CYMENE,para                         | 0.990 | 1.193 | 1.000 | 0.871           |           |
| CYTOSINE (M)                        | 0.990 | 0.811 | 0.811 | 0.831           | 71-30-7   |
| DAUNOMYCIN(M)                       | 0.990 | 0.000 | 0.000 | 0.853           | 20830-81- |
| DAZOMET                             | 0.900 | 0.085 | 0.153 | 0.066           |           |
| DDD,p,p'-                           | 0.950 | 1.150 | 1.000 | 0.394           | 72-54-8   |
| DDE,p,p'-                           | 0.990 | 1.138 | 0.990 | 0.621           | 72-55-9   |
| DDT                                 | 0.990 | 1.131 | 1.000 | 0.980           | 50-29-3   |
| DECANAL                             | 0.990 | 0.918 | 0.928 | 0.612           |           |
| DECENE, 8 METHYL 1-                 | 0.990 | 1.116 | 1.000 | 0.980           |           |
| DIACETYL (M)                        | 0.990 | 0.999 | 0.999 | 0.318           | 431-03-8  |
| DIAMINO-5-SULFONYL BENZYL 2,4 (M)   | 0.990 | 0.133 | 0.133 | 0.628           |           |
| DIAMINODIPHENYLMETHANE P,P' (M)     | 0.990 | 0.126 | 0.126 | 0.980           | 101-77-9  |
| DIAZOMETHANE                        | 0.575 | 0.573 | 1.000 | 0.356           |           |
| DIBENZOFURANS                       | 0.990 | 1.112 | 0.967 | 0.740           |           |
| DIBENZOPYRENE 1,2,7,8               | 0.990 | 0.803 | 0.633 | 0.720           |           |
| DIBROMO-3-CHLOROPROPANE,1,2         | 0.709 | 1.048 | 1.000 | 0.185           | 96-12-8   |
| DIBROMOCHLOROMETHANE                | 0.990 | 0.585 | 1.000 | 0.643           | 124-48-1  |
| DIBROMOETHANE-1,2                   | 0.990 | 1.114 | 1.000 | 0.852           | 106-93-4  |
| DIBROMOMETHANE                      | 0.990 | 0.493 | 1.000 | 0.558           | 74-95-3   |
| DIBUTYL ETHER                       | 0.990 | 0.958 | 1.000 | 0.727           | 142-96-1  |
| DIBUTYLAMINE                        | 0.990 | 0.949 | 0.984 | 0.300           |           |
| DICHLORO 2-PROPANOL 1,3             | 0.990 | 0.237 | 0.257 | 0.570           | 96-23-1   |
| DICHLORO PROPANOL 2,3               | 0.507 | 0.119 | 0.130 | 0.255           | 616-23-9  |
| DICHLORO-1,3-CYCLOPENTADIENE 5,5(M) | 0.990 | 0.413 | 0.413 | 0.980           |           |
| DICHLORO-2-BUTENE 1,2               | 0.990 | 1.079 | 1.000 | 0.562           |           |
| DICHLORO-2-BUTENE(1,4)              | 0.990 | 1.079 | 1.000 | 0.453           | 764-41-0  |
| DICHLORO-2-BUTENE, 1,4              | 0.990 | 1.079 | 1.000 | 0.612           |           |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                             | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|--------------------------------------|-------|-------|-------|-----------------|-----------|
| DICHLOROANILINE(2,3)                 | 0.527 | 0.121 | 0.117 | 0.064           |           |
| DICHLOROBENZENE(1,2) (-o)            | 0.990 | 1.134 | 1.000 | 0.637           | 95-50-1   |
| DICHLOROBENZENE(1,3) (-m)            | 0.990 | 1.134 | 1.000 | 0.719           | 541-73-1  |
| DICHLOROBENZENE(1,4) (-p)            | 0.990 | 1.134 | 1.000 | 0.724           | 106-46-7  |
| DICHLOROBENZIDINE,3,3'               | 0.001 | 0.055 | 0.053 | 0.026           | 91-94-1   |
| DICHLOROBENZOPHENONE P,P             | 0.978 | 0.366 | 0.332 | 0.093           | 90-98-2   |
| DICHLOROBIPHENYL (PARA)              | 0.990 | 1.177 | 1.000 | 0.914           | 213029-08 |
| DICHLOROBUTANE (1,4)                 | 0.990 | 1.052 | 1.000 | 0.980           | 110-56-5  |
| DICHLORODIPHENYLMETHANE (M)          | 0.990 | 0.107 | 0.107 | 0.855           | 2051-90-3 |
| DICHLOROETHANE(1,1)                  | 0.990 | 1.024 | 1.000 | 0.792           | 75-34-3   |
| DICHLOROETHANE(1,2)                  | 0.990 | 1.040 | 1.000 | 0.640           | 107-06-2  |
| DICHLOROETHENE 1,2 trans             | 0.990 | 1.061 | 1.000 | 0.981           | 156-60-5  |
| DICHLOROETHENE(1,1)                  | 0.990 | 1.061 | 1.000 | 0.937           | 75-35-4   |
| DICHLOROETHYL ETHER                  | 0.872 | 0.711 | 0.757 | 0.212           |           |
| DICHLOROETHYLENE(1,2) cis            | 0.990 | 1.061 | 1.000 | 0.904           | 156-54-2  |
| DICHLOROIODOMETHANE                  | 0.990 | 0.553 | 0.975 | 0.362           |           |
| DICHLOROMONOFUOROMETHANE             | 0.990 | 1.023 | 1.000 | 0.989           | 75-43-4   |
| DICHLOROPHENOL                       | 0.990 | 0.940 | 0.920 | 0.227           |           |
| DICHLOROPHENOL(2,4)                  | 0.945 | 0.158 | 0.154 | 0.094           | 120-83-2  |
| DICHLOROPHENOL(2,6)                  | 0.846 | 0.213 | 0.209 | 0.094           | 87-65-0   |
| DICHLOROPHENOXYACETIC ACID(2,4)      | 0.990 | 0.922 | 1.000 | 0.978           | 94-75-7   |
| DICHLOROPROPANE 1,2                  | 0.990 | 1.054 | 1.000 | 0.720           | 78-87-5   |
| DICHLOROPROPENE(1,3)                 | 0.990 | 1.071 | 1.000 | 0.759           | 542-75-6  |
| DICHLOROPROPYLENE,1,2-(cis)          | 0.990 | 1.062 | 1.000 | 0.831           |           |
| DICHLOROPROPYLENE,1,2-(trans)        | 0.990 | 1.072 | 1.000 | 0.853           | 563-54-2  |
| DICHLOROPROPYLENE-2,3                | 0.990 | 1.071 | 1.000 | 0.857           | 78-88-6   |
| DICHLOROSTYRENE 2,6                  | 0.990 | 1.149 | 1.000 | 0.823           |           |
| DICHLORO-TRANS-ETHYLENE(1,2)         | 0.990 | 1.061 | 1.000 | 0.980           | 540-59-0  |
| DIELDRIN                             | 0.990 | 0.259 | 0.235 | 0.225           | 60-57-1   |
| DIETHYL AMINE                        | 0.828 | 0.865 | 1.000 | 0.286           | 109-89-7  |
| DIETHYL ETHER                        | 0.990 | 0.856 | 1.000 | 0.423           | 602-97-6  |
| DIETHYL ETHER ACID CHLORIDE (M)      | 0.990 | 0.379 | 0.379 | 0.980           |           |
| DIETHYL PHTHALATE                    | 0.990 | 0.054 | 0.063 | 0.853           | 84-66-2   |
| DIETHYL SULFATE                      | 0.909 | 0.001 | 0.002 | 0.107           |           |
| DIETHYL THIOETHER (M)                | 0.990 | 0.999 | 0.999 | 0.980           | 352-93-2  |
| DIETHYLBENZENE P                     | 0.990 | 1.191 | 1.000 | 0.784           | 105-05-5  |
| DIETHYLDIPHENYL UREA SYM(M)          | 0.990 | 0.091 | 0.091 | 0.859           | 85-98-3   |
| DIETHYLENE GLYCOL DIETHYL ETHER      | 0.316 | 0.168 | 0.217 | 0.033           |           |
| DIETHYLUREA 1,1 (M)                  | 0.729 | 0.726 | 0.726 | 0.101           | 634-95-7  |
| DIHYDRO-5-OXAZALONE (DIHYDROAZLA (M) | 0.990 | 0.982 | 0.982 | 0.722           |           |
| DIISOBUTYLENE                        | 0.990 | 1.127 | 1.000 | 0.980           |           |
| DIISODECYL PHTHALATE                 | 0.990 | 0.007 | 0.007 | 0.451           |           |
| DIISOPROPYL BENZENE (PARA)           | 0.990 | 1.184 | 1.000 | 0.980           | 100-18-5  |
| DIISOPROPYL KETONE                   | 0.990 | 0.973 | 1.000 | 0.483           |           |
| DIISOPROPYLAMINE                     | 0.990 | 0.939 | 1.000 | 0.409           |           |
| DIMETHOXY METHANE                    | 0.878 | 0.594 | 0.950 | 0.442           | 109-87-5  |
| DIMETHOXY-(3,3')-BENZIDINE           | 0.990 | 0.000 | 0.000 | 0.660           | 119-90-4  |
| DIMETHYL AMINE                       | 0.321 | 0.709 | 0.996 | 0.198           | 124-40-3  |
| DIMETHYL BENZ(A)ANT 7,12             | 0.990 | 1.214 | 0.973 | 0.857           |           |
| DIMETHYL BENZOIC ACID, 2,4           | 0.854 | 0.101 | 0.105 | 0.115           |           |
| DIMETHYL BENZOIC ACID, 3,5           | 0.854 | 0.101 | 0.105 | 0.115           |           |
| DIMETHYL BENZYLAMINE N,N             | 0.990 | 0.003 | 0.003 | 0.587           | 103-83-3  |
| DIMETHYL METHYLTHIOCARBAMATE N,N(M)  | 0.990 | 0.676 | 0.676 | 0.863           |           |
| DIMETHYL NITROISOPROPYLAMINE N,N(M)  | 0.990 | 0.439 | 0.439 | 0.389           |           |
| DIMETHYL NITROSAMINE (M)             | 0.990 | 0.999 | 0.999 | 0.980           |           |
| DIMETHYL SULFATE                     | 0.549 | 0.034 | 0.086 | 0.079           | 77-78-1   |
| DIMETHYL SULFIDE                     | 0.990 | 0.508 | 1.000 | 0.829           | 75-18-3   |
| DIMETHYL TRISULFIDE                  | 0.990 | 0.354 | 1.000 | 0.980           |           |
| DIMETHYL-1-NITROBENZENE 2,4          | 0.990 | 0.564 | 0.669 | 0.801           | 25168-04- |
| DIMETHYLACETAMIDE                    | 0.547 | 0.707 | 0.994 | 0.284           |           |
| dimethylaniline N,N                  | 0.990 | 0.000 | 0.001 | 0.342           | 57-14-7   |
| DIMETHYLBENZYL HYDROPEROXIDE (M)     | 0.990 | 0.149 | 0.149 | 0.466           | 80-15-9   |
| DIMETHYLETHYLAMINE                   | 0.990 | 0.865 | 1.000 | 0.523           | 75-64-9   |
| DIMETHYLGLYCOL                       | 0.990 | 0.102 | 0.136 | 0.483           |           |
| DIMETHYLHYDANTOIN,5,5-(M)            | 0.990 | 0.521 | 0.521 | 0.980           | 77-71-4   |
| DIMETHYLPHENOL (2,4)                 | 0.990 | 0.050 | 0.047 | 0.552           | 105-67-9  |
| DIMETHYLPHENYLCARBINOL (M)           | 0.990 | 0.385 | 0.385 | 0.794           | 617-94-7  |
| DIMETHYLSULFOXIDE                    | 0.854 | 0.821 | 0.990 | 0.419           |           |
| DINITROBENZENE M                     | 0.023 | 0.564 | 1.000 | 0.285           | 99-65-0   |
| DINITROPHENOL 2,4                    | 0.990 | 0.004 | 0.008 | 0.059           | 51-28-5   |
| DINITROTOLUENE 2,6                   | 0.990 |       |       | 0.109           | 606-20-2  |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                                 | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|--|-------|-------|-------|-----------------|-----------|
| DINITROTOLUENE (2,4)                     | 0.390 | 0.052 | 0.085 | 0.178           | 121-14-2  |
| DINOCAP (M)                              | 0.990 | 0.043 | 0.043 | 0.980           | 39300-45- |
| DI-n-OCTYL PHTHALATE                     | 0.990 | 0.000 | 0.000 | 0.980           | 117-84-0  |
| DINOSEB (M)                              | 0.990 | 0.105 | 0.105 | 0.575           | 88-85-7   |
| DIOXANE (1,4)                            | 0.387 | 0.618 | 0.869 | 0.181           | 123-91-1  |
| DIOXIN (M)                               | 0.990 | 0.064 | 0.064 | 0.279           | 828-00-2  |
| DIPHENYL ETHER (M)                       | 0.990 | 0.140 | 0.140 | 0.662           | 101-84-8  |
| DIPHENYL THIOETHER (M)                   | 0.990 | 0.132 | 0.132 | 0.838           | 139-66-2  |
| DIPHENYLAMINE (M)                        | 0.513 | 0.140 | 0.140 | 0.074           | 122-39-4  |
| DIPHENYLBUTADIENE 1,3 (M)                | 0.990 | 0.122 | 0.122 | 0.647           | 886-65-7  |
| DIPHENYLCHLOROMETHANE (M)                | 0.990 | 0.124 | 0.124 | 0.850           | 90-99-3   |
| DIPHENYLDIKETONE (M)                     | 0.990 | 0.120 | 0.120 | 0.851           | 134-81-6  |
| DIPHENYLETHANE 1,1 (M)                   | 0.990 | 0.134 | 0.134 | 0.551           |           |
| DIPHENYLETHANOL 1,1 (M)                  | 0.416 | 0.126 | 0.126 | 0.066           | 599-67-7  |
| DIPHENYLHYDRAZINE,1,1-(M)                | 0.990 | 0.133 | 0.133 | 0.796           | 530-50-7  |
| DIPHENYLMETHANE                          | 0.990 | 0.628 | 0.509 | 0.195           | 101-81-5  |
| DIPROPYLAMINE                            | 0.979 | 0.927 | 0.998 | 0.411           | 142-84-7  |
| DIPROPYLBUTRAL                           | 0.990 | 0.622 | 0.618 | 0.292           |           |
| DIPROPYLFORMAMIDE (M)                    | 0.990 | 0.503 | 0.503 | 0.980           | 6282-00-4 |
| DI-tert-BUTYL-p-CRESOL                   | 0.990 | 0.031 | 0.028 | 0.072           | 128-37-0  |
| DIVINYL KETONE (M)                       | 0.990 | 0.999 | 0.999 | 0.457           |           |
| dodecane                                 | 0.990 | 1.089 | 1.000 | 0.980           |           |
| EDTA (M)                                 | 0.990 | 0.999 | 0.999 | 0.412           | 60-00-4   |
| ENDOSULFAN                               | 0.900 | 0.020 | 0.018 | 0.102           | 115-29-7  |
| ENDOSULFAN SULFATE (M)                   | 0.990 | 0.014 | 0.014 | 0.980           | 1031-07-8 |
| ENDRIN ALDEHYDE (M)                      | 0.990 | 0.999 | 0.999 | 0.412           |           |
| EPICHLOROHYDRIN                          | 0.915 | 0.847 | 0.939 | 0.350           | 106-89-8  |
| EPOXYBUTANE 1,2                          | 0.990 | 0.879 | 1.000 | 0.582           |           |
| ETHANE                                   | 0.990 | 1.067 | 1.000 | 0.946           |           |
| ETHANOL                                  | 0.322 | 0.586 | 0.860 | 0.126           | 64-17-5   |
| ETHENE                                   | 0.990 | 1.187 | 1.000 | 0.980           |           |
| ETHENYL 2 METHYL BENZENE, 1-             | 0.990 | 1.240 | 1.000 | 0.710           |           |
| ETHOXYETHANOL-2                          | 0.545 | 0.144 | 0.207 | 0.134           | 110-80-5  |
| ETHYL 2 METHYL BENZENE, 1-               | 0.990 | 1.198 | 1.000 | 0.731           |           |
| ETHYL ACETATE PEROXIDE (M)               | 0.990 | 0.659 | 0.659 | 0.706           |           |
| ETHYL ACRYLATE                           | 0.990 | 0.788 | 1.000 | 0.483           | 140-88-5  |
| ETHYL BUTANOATE                          | 0.990 | 0.775 | 1.000 | 0.457           |           |
| ETHYL CYANIDE (PROPIONITRILE) (M)        | 0.990 | 0.999 | 0.999 | 0.580           | 107-12-0  |
| ETHYL ETHER                              | 0.990 | 0.856 | 1.000 | 0.506           | 60-29-7   |
| ETHYL HEPTANOATE                         | 0.990 | 0.868 | 1.000 | 0.470           |           |
| ETHYL ISOPROPYL PEROXIDE (M)             | 0.990 | 0.931 | 0.931 | 0.386           |           |
| ETHYL METHANOATE                         | 0.990 | 0.537 | 1.000 | 0.566           |           |
| ETHYL PENTANOATE                         | 0.990 | 0.813 | 1.000 | 0.428           |           |
| ETHYL PEROXIDE                           | 0.341 | 0.146 | 0.283 | 0.112           |           |
| ETHYL PROPYL ETHER                       | 0.990 | 0.894 | 1.000 | 0.571           |           |
| ETHYL S,S-DIPHENYL PHOSPHORODITH (M)     | 0.990 | 0.070 | 0.070 | 0.333           | 1709-49-8 |
| ETHYL TOLUENE, 4                         | 0.990 | 1.198 | 1.000 | 0.857           |           |
| ETHYL VINYL ETHER                        | 0.990 | 0.890 | 1.000 | 0.652           |           |
| ETHYL(2) HEXANOL                         | 0.990 | 0.256 | 0.268 | 0.266           | 104-76-7  |
| ETHYL-(2)-PROPYL-(3) ACROLEIN (M)        | 0.977 | 0.999 | 0.999 | 0.257           | 645-62-5  |
| ETHYLACETATE                             | 0.987 | 0.722 | 1.000 | 0.404           | 141-78-6  |
| ETHYLAMINE                               | 0.358 | 0.711 | 0.999 | 0.280           | 75-04-7   |
| ETHYLBENZENE                             | 0.990 | 1.204 | 1.000 | 0.828           | 100-41-4  |
| ETHYLENE                                 | 0.990 | 1.187 | 1.000 | 0.980           | 74-85-1   |
| ETHYLENE DIAMINE                         | 0.963 | 0.012 | 0.022 | 0.241           | 107-15-3  |
| ETHYLENE DIBROMIDE                       | 0.990 | 0.537 | 0.999 | 0.565           | 106-93-4  |
| ETHYLENE GLYCOL DIMETHYL ETHER           | 0.905 | 0.601 | 0.860 | 0.316           | 110-71-4  |
| ETHYLENE GLYCOL MONOBUTYL ETHER ACETATE  | 0.772 | 0.031 | 0.043 | 0.067           |           |
| ETHYLENE GLYCOL MONOMETHYL ETHER ACETATE | 0.285 | 0.055 | 0.093 | 0.048           | 110-49-6  |
| ETHYLENE OXIDE                           | 0.986 | 0.712 | 1.000 | 0.503           | 75-21-8   |
| ETHYLEHOXY PROPIONATE                    | 0.940 | 0.491 | 0.577 | 0.213           |           |
| ETHYLHEXYL HEXANOL 2                     | 0.990 | 0.065 | 0.064 | 0.125           |           |
| ETHYLHEXYLACRYLATE,2-                    | 0.990 | 0.925 | 0.992 | 0.705           | 103-11-7  |
| FENCHONE,d- (M)                          | 0.990 | 0.149 | 0.149 | 0.406           | 4695-62-9 |
| FLUORANTHENE                             | 0.990 | 0.049 | 0.039 | 0.656           | 206-44-0  |
| FLUORENE                                 | 0.990 | 0.965 | 0.774 | 0.314           | 86-73-7   |
| FLUOROMETHANE                            | 0.990 | 1.130 | 1.000 | 0.873           |           |
| FLUOROURACIL,5- (M)                      | 0.990 | 0.999 | 0.999 | 0.412           | 51-21-8   |
| FORMYL FLUORIDE                          | 0.990 | 0.848 | 1.000 | 0.577           |           |
| FREON 11, fluorotrichloromethane         | 0.990 | 1.053 | 1.000 | 0.954           |           |
| FREON 12 DICHLORODIFLUOROMETHANE         | 0.990 | 1.059 | 1.000 | 0.980           | 75-71-8   |



TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                              | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|---------------------------------------|-------|-------|-------|-----------------|-----------|
| FREON 12, dichlorodifluoromethane     | 0.990 | 1.059 | 1.000 | 0.980           |           |
| FREONS (M)                            | 0.990 | 0.644 | 0.644 | 0.980           |           |
| FURAN                                 | 0.990 | 0.983 | 1.000 | 0.755           | 110-00-9  |
| FURFURAL                              | 0.990 | 0.288 | 0.334 | 0.354           | 98-01-1   |
| FUROIC ACID (M)                       | 0.990 | 0.794 | 0.794 | 0.480           | 88-14-2   |
| GEOSMIN (M)                           | 0.990 | 0.134 | 0.134 | 0.406           | 19700-21- |
| GLYOXAL                               | 0.502 | 0.490 | 0.888 | 0.297           |           |
| GUANINE (M)                           | 0.990 | 0.149 | 0.149 | 0.980           | 73-40-5   |
| HEPTACHLOR                            | 0.990 | 0.619 | 0.566 | 0.647           | 76-44-8   |
| HEPTACHLOR EPOXIDE (M)                | 0.976 | 0.030 | 0.030 | 0.162           | 1024-57-3 |
| HEPTANAL                              | 0.990 | 0.942 | 0.991 | 0.407           |           |
| HEPTANE ISO                           | 0.990 | 1.099 | 1.000 | 0.980           | 31394-54- |
| HEPTANE(-n)                           | 0.990 | 1.085 | 1.000 | 0.980           | 142-82-5  |
| HEXACHLORO BENZENE                    | 0.990 | 1.047 | 0.966 | 0.643           | 118-74-1  |
| HEXACHLOROBUTADIENE                   | 0.990 | 0.937 | 0.883 | 0.855           | 87-68-3   |
| HEXACHLOROCYCLOHEXANE (GAMMA ISOMER)  | 0.990 | 0.141 | 0.132 | 0.106           | 58-89-9   |
| HEXACHLOROCYCLOPENTADIENE             | 0.990 | 0.886 | 0.826 | 0.803           | 77-47-4   |
| HEXACHLOROETHANE                      | 0.990 | 0.515 | 0.499 | 0.852           | 67-72-1   |
| HEXACHLOROPENTADIENE (M)              | 0.990 | 0.088 | 0.088 | 0.860           |           |
| HEXADECANE N (M)                      | 0.990 | 0.112 | 0.112 | 0.980           | 544-76-3  |
| HEXAFLUOROACETONE                     | 0.990 | 0.968 | 1.000 | 0.980           |           |
| HEXAFLUOROPROPENE                     | 0.990 | 1.080 | 1.000 | 0.980           | 116-15-4  |
| HEXAMETHYLENEDIAMINE (M)              | 0.971 | 0.724 | 0.724 | 0.213           | 124-09-4  |
| HEXAMETHYLENIMINE                     | 0.520 | 0.923 | 0.989 | 0.109           |           |
| HEXANAL                               | 0.990 | 0.928 | 0.997 | 0.400           |           |
| HEXANE(-n)                            | 0.990 | 1.084 | 1.000 | 1.000           | 110-54-3  |
| HEXANOL 2 ETHYL                       | 0.942 | 0.256 | 0.268 | 0.134           | 104-76-7  |
| HEXANOL-1                             | 0.963 | 0.322 | 0.355 | 0.180           | 111-27-3  |
| HEXEN-2-ONE 5                         | 0.979 | 0.885 | 0.915 | 0.347           |           |
| HEXENE                                | 0.990 | 1.119 | 1.000 | 0.980           |           |
| HEXYL ETHANOATE                       | 0.990 | 0.865 | 0.998 | 0.475           |           |
| HEXYLAMINE                            | 0.948 | 0.803 | 0.870 | 0.239           |           |
| HYDROFLUORIC ACID (M)                 | 0.990 | 0.558 | 0.558 | 0.537           | 7664-39-3 |
| HYDROGEN SULFIDE                      | 0.990 | 0.333 | 1.000 | 0.882           |           |
| HYDROXY DIMETHYL ETHER (M)            | 0.990 | 0.999 | 0.999 | 0.874           |           |
| HYDROXY-1,3-CYCLOPENTADIENE 5 (M)     | 0.990 | 0.999 | 0.999 | 0.728           |           |
| HYDROXY-4-METHYLTETRAHYDROFURAN(M)    | 0.990 | 0.948 | 0.948 | 0.385           |           |
| HYDROXY-5-METHYLDIMETHYL PHTHALA (M)  | 0.990 | 0.113 | 0.113 | 0.980           |           |
| HYDROXY6METHYLPYRIDINE3 CARBOXYLI (M) | 0.990 | 0.148 | 0.148 | 0.409           | 38116-61- |
| HYDROXYACETIC ACID                    | 0.760 | 0.000 | 0.001 | 0.570           | 79-14-1   |
| HYDROXYCYCLOHEXANONE 4 (M)            | 0.631 | 0.761 | 0.761 | 0.087           |           |
| HYDROXYDIMETHYL PHTHALATE 4 (M)       | 0.990 | 0.120 | 0.120 | 0.980           |           |
| HYDROXYMETHYL ACETYLENE (M)           | 0.990 | 0.999 | 0.999 | 0.980           |           |
| HYDROXYMETHYL ISOPROPYL KETONE (M)    | 0.990 | 0.999 | 0.999 | 0.662           |           |
| HYDROXYMETHYL, N-METHYLETHYL AMI (M)  | 0.990 | 0.999 | 0.999 | 0.980           |           |
| HYDROXYMETHYL-N-CHLOROMETHYLETHY (M)  | 0.990 | 0.838 | 0.838 | 0.980           |           |
| HYDROXYMETHYLPHENYL CARBAMATE N (M)   | 0.920 | 0.147 | 0.147 | 0.137           |           |
| HYDROXYMETHYLTHIOBENZENE (M)          | 0.990 | 0.320 | 0.320 | 0.790           |           |
| HYDROXYMETHYLVINYL ETHER (M)          | 0.990 | 0.490 | 0.490 | 0.905           |           |
| HYDROXPENTANE 3 (M)                   | 0.990 | 0.999 | 0.999 | 0.450           |           |
| INDANOL,5-(M)                         | 0.990 | 0.128 | 0.128 | 0.980           | 1470-94-6 |
| INDOLE (M)                            | 0.990 | 0.708 | 0.708 | 0.980           | 120-72-9  |
| IODOCUMARAN 2 (M)                     | 0.990 | 0.102 | 0.102 | 0.980           |           |
| ISOBUTANE                             | 0.990 | 1.103 | 1.000 | 0.963           |           |
| ISOBUTYL ETHANOATE                    | 0.990 | 0.786 | 1.000 | 0.486           |           |
| ISOBUTYLBENZENE                       | 0.990 | 1.191 | 1.000 | 0.905           |           |
| ISOBUTYLENE                           | 0.990 | 1.141 | 1.000 | 0.916           |           |
| ISOCYANO 4 METHYL BENZENE *           | 0.980 | 0.422 | 0.384 | 0.198           |           |
| ISODECANOL                            | 0.932 | 0.165 | 0.158 | 0.099           |           |
| ISODECYL OCTYL ESTER                  | 0.990 | 1.033 | 1.000 | 0.906           |           |
| ISOPENTANE                            | 0.990 | 1.101 | 1.000 | 0.954           |           |
| ISOPENTYL ETHANOATE                   | 0.990 | 0.852 | 0.999 | 0.487           |           |
| ISOPENTYL METHANOATE                  | 0.990 | 0.941 | 0.997 | 0.503           |           |
| ISOPHORONE                            | 0.616 | 0.525 | 0.506 | 0.108           | 78-59-1   |
| ISOPROPYL AMINE                       | 0.990 | 0.811 | 1.000 | 0.538           | 75-31-0   |
| ISOPROPYL ETHER                       | 0.019 | 0.939 | 1.000 | 0.730           | 108-20-3  |
| ISOPROPYL METHANOATE                  | 0.990 | 0.886 | 1.000 | 0.578           |           |
| ISOPROPYL METHANOATE                  | 0.990 | 0.865 | 1.000 | 0.547           |           |
| ISOPROPYL PROPANOATE                  | 0.990 | 0.825 | 1.000 | 0.487           |           |
| ISOXAZOLOL,5-(AMINOMETHYL)-3-(M)      | 0.990 | 0.760 | 0.760 | 0.980           | 2763-96-4 |
| LINDANE hexachlorocyclohexane         | 0.990 | 1.063 | 1.000 | 0.703           |           |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                             | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|--------------------------------------|-------|-------|-------|-----------------|-----------|
| MELAMINE (M)                         | 0.990 | 0.554 | 0.554 | 0.980           | 108-78-1  |
| MERCAPTOBENZOTHAZOLE,2               | 0.990 | 0.844 | 1.000 | 0.641           |           |
| MERCURY (M)                          | 0.990 | 0.125 | 0.125 | 0.854           | 7439-97-6 |
| METHACRYLIC ACID                     | 0.990 | 0.068 | 0.091 | 0.194           | 79-41-4   |
| METHANE                              | 0.990 | 1.067 | 1.000 | 0.980           | 74-82-8   |
| METHANETHIOL (M)                     | 0.990 | 0.999 | 0.999 | 0.731           | 74-93-1   |
| METHANOL                             | 0.317 | 0.433 | 0.855 | 0.168           | 67-56-1   |
| METHAPYRILENE (M)                    | 0.990 | 0.094 | 0.094 | 0.980           | 91-80-5   |
| METHOXYACETIC ACID                   | 0.593 | 0.005 | 0.010 | 0.064           | 625-45-6  |
| METHOXYACETONITRILE (M)              | 0.990 | 0.999 | 0.999 | 0.382           | 1738-36-9 |
| METHOXYCHLOR                         | 0.990 | 0.085 | 0.081 | 0.333           | 72-43-5   |
| METHYL 1-PENTENE 2                   | 0.990 | 1.125 | 1.000 | 0.980           | 763-29-1  |
| METHYL 2-PROPYL ETHER                | 0.990 | 0.976 | 1.000 | 0.537           |           |
| METHYL ACETATE                       | 0.989 | 0.590 | 0.906 | 0.454           | 79-20-9   |
| METHYL ACRYLATE                      | 0.990 | 0.748 | 1.000 | 0.478           | 96-33-3   |
| METHYL ACRYLONITRILE (M)             | 0.990 | 0.999 | 0.999 | 0.980           | 126-98-7  |
| METHYL AMINE                         | 0.990 | 0.516 | 0.992 | 0.877           | 74-89-5   |
| METHYL AMINOACETYLENE (M)            | 0.990 | 0.999 | 0.999 | 0.980           |           |
| METHYL AZIRIDINE 2                   | 0.900 | 0.838 | 1.000 | 0.360           |           |
| METHYL BENZOATE                      | 0.692 | 0.924 | 0.981 | 0.168           |           |
| METHYL BENZYL ALCOHOL 4              | 0.917 | 0.058 | 0.056 | 0.154           |           |
| METHYL BIPHENYL (-p) (M)             | 0.990 | 0.141 | 0.141 | 0.819           | 644-08-6  |
| METHYL BUTANOATE                     | 0.990 | 0.775 | 1.000 | 0.413           |           |
| METHYL CHLORIDE                      | 0.990 | 1.040 | 1.000 | 0.840           | 74-87-3   |
| METHYL CHLOROACETAMIDE N (M)         | 0.863 | 0.872 | 0.872 | 0.137           |           |
| METHYL CHLOROCARBONATE (M)           | 0.990 | 0.999 | 0.999 | 0.980           | 79-22-1   |
| METHYL CHOLANTHRENE 3                | 0.990 | 1.234 | 0.990 | 0.322           | 56-49-5   |
| METHYL COUMARAN 2 (M)                | 0.990 | 0.145 | 0.145 | 0.811           | 607-71-6  |
| METHYL CYCLOHEXANE                   | 0.990 | 1.107 | 1.000 | 0.980           | 108-87-2  |
| METHYL ETHER dimethyl ether          | 0.990 | 0.698 | 1.000 | 0.730           | 115-10-6  |
| METHYL ETHYL ETHER                   | 0.990 | 0.791 | 1.000 | 0.617           |           |
| METHYL ETHYL KETONE, 2 butanone      | 0.958 | 0.872 | 0.990 | 0.477           | 78-93-3   |
| METHYL FORMATE                       | 0.590 | 0.535 | 0.997 | 0.548           | 107-31-3  |
| METHYL HEXANOATE                     | 0.990 | 0.843 | 1.000 | 0.441           |           |
| METHYL IODIDE                        | 0.990 | 0.354 | 1.000 | 0.711           | 74-88-4   |
| METHYL ISOAMYL KETONE (M)            | 0.990 | 0.761 | 0.761 | 0.318           | 110-12-3  |
| METHYL ISOBUTYL KETONE               | 0.990 | 0.933 | 0.979 | 0.529           | 108-10-1  |
| METHYL ISOCYANATE                    | 0.990 | 0.272 | 1.000 | 0.870           | 624-83-9  |
| METHYL ISOPROPYL KETONE              | 0.986 | 0.922 | 0.991 | 0.523           | 563-80-4  |
| METHYL MERCAPTAN                     | 0.990 | 0.333 | 1.000 | 0.719           |           |
| METHYL METHACRYLATE                  | 0.986 | 0.801 | 0.999 | 0.366           | 80-62-6   |
| METHYL MORPHOLINE                    | 0.435 | 0.365 | 0.475 | 0.078           |           |
| METHYL NAPHTHALENE (1-)              | 0.990 | 1.204 | 0.973 | 0.512           | 90-12-0   |
| METHYL NAPHTHALENE (-12)             | 0.990 | 1.219 | 0.986 | 0.246           | 91-57-6   |
| METHYL OCTANOATE                     | 0.990 | 0.888 | 1.000 | 0.524           |           |
| METHYL PENTANOATE                    | 0.990 | 0.813 | 1.000 | 0.417           |           |
| METHYL PEROXIDE                      | 0.587 | 0.024 | 0.070 | 0.159           |           |
| METHYL PROPANOATE                    | 0.985 | 0.724 | 1.000 | 0.431           |           |
| METHYL PROPENE 2 (M)                 | 0.990 | 0.999 | 0.999 | 0.980           | 115-11-7  |
| METHYL PROPYL ETHER                  | 0.990 | 0.848 | 1.000 | 0.598           |           |
| METHYL TERTIARY-BUTYL ETHER          | 0.990 | 0.911 | 1.000 | 0.573           | 1634-04-4 |
| METHYL TETRAHYDROFURAN 2             | 0.990 | 0.914 | 1.000 | 0.357           |           |
| METHYL THIOURACIL (M)                | 0.990 | 0.283 | 0.283 | 0.753           | 56-04-2   |
| METHYL-1,3-CYCLOPENTADIENE 5 (M)     | 0.990 | 0.999 | 0.999 | 0.924           | 26519-91- |
| METHYL-2,3,4-TRIHYDROQUINOLINE N (M) | 0.912 | 0.218 | 0.218 | 0.137           |           |
| METHYL-2-AMINOETHYLAMINE (M)         | 0.990 | 0.999 | 0.999 | 0.871           | 109-81-9  |
| METHYL-2-HYDROXYETHYLAMINE (M)       | 0.578 | 0.999 | 0.999 | 0.081           | 109-83-1  |
| METHYL-3-ACETYLCYCLOPENTADIENE 1 (M) | 0.990 | 0.897 | 0.897 | 0.754           |           |
| METHYL-3-NITROBENZYL ALCOHOL 4 (M)   | 0.767 | 0.141 | 0.141 | 0.103           | 40870-59- |
| METHYL-4-NITROBENZYL ALCOHOL 2 (M)   | 0.568 | 0.141 | 0.141 | 0.079           | 23876-13- |
| METHYL-5-THIOACETYLDIHYDRO1,3THI (M) | 0.994 | 0.146 | 0.146 | 0.980           |           |
| METHYLACETONITRILE (M)               | 0.990 | 0.999 | 0.999 | 0.980           | 75-86-5   |
| METHYLBUTADIENE (isoprene)           | 0.990 | 1.176 | 1.000 | 0.980           |           |
| METHYLBUTYLAMINE                     | 0.809 | 0.791 | 0.883 | 0.178           |           |
| METHYLCYCLOPENTANE                   | 0.990 | 1.109 | 1.000 | 0.980           |           |
| METHYLENE CHLORIDE, dichloromethane  | 0.990 | 1.017 | 1.000 | 0.770           | 75-09-2   |
| METHYLFURAN 2 (M)                    | 0.509 | 0.999 | 0.999 | 0.073           | 534-22-5  |
| METHYLISOBORNEOL,2-(M)               | 0.990 | 0.141 | 0.141 | 0.794           | NA        |
| METHYLPHENYL CARBAMATE N (M)         | 0.906 | 0.320 | 0.320 | 0.137           |           |
| METHYL-PHENYLETHYLAMINE N (M)        | 0.990 | 0.401 | 0.401 | 0.587           | 589-08-2  |
| METHYL-p'-METHYLTRIPHENYL PHOSPH (M) | 0.990 | 0.079 | 0.079 | 0.862           |           |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                            | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|-------------------------------------|-------|-------|-------|-----------------|-----------|
| METHYLSTYRENE (-4)                  | 0.990 | 1.217 | 1.000 | 0.767           | 98-93-9   |
| METHYLTRIN TRICHLORIDE (M)          | 0.470 | 0.105 | 0.105 | 0.070           | 993-16-8  |
| METHYL-TRIHYDRO-1,3-THIAZOLE 4 (M)  | 0.990 | 0.914 | 0.914 | 0.316           |           |
| MITOMYCIN C (M)                     | 0.990 | 0.058 | 0.058 | 0.980           | 50-07-7   |
| MNNG (M)                            | 0.990 | 0.199 | 0.199 | 0.980           | 70-25-7   |
| MONOCHLORODIFLUOROMETHANE           | 0.990 | 1.023 | 1.000 | 0.990           | 75-45-6   |
| MORPHOLINE                          | 0.990 | 0.148 | 0.207 | 0.437           | 110-91-8  |
| MUSTARD GAS (M)                     | 0.990 | 0.146 | 0.146 | 0.406           | 505-60-2  |
| NAPHTHALENE                         | 0.990 | 1.239 | 0.994 | 0.506           |           |
| NAPHTHALENE ACETIC ACID 2 METHYL,   | 10.99 | 0.863 | 0.830 | 0.567           |           |
| NAPHTHOQUINONE-1,4 (M)              | 0.958 | 0.146 | 0.146 | 0.164           | 130-15-4  |
| NICKEL CYANIDE (M)                  | 0.990 | 0.817 | 0.817 | 0.284           | 557-19-7  |
| NITRO m XYLENE, 2                   | 0.990 | 0.779 | 0.923 | 0.455           |           |
| NITRO-4-METHYLBENZOATE 3 (M)        | 0.990 | 0.128 | 0.128 | 0.980           |           |
| NITROANILINE P                      | 0.990 | 0.000 | 0.000 | 0.411           | 100-01-6  |
| NITROBENZENE                        | 0.808 | 0.305 | 0.394 | 0.228           | 98-95-3   |
| NITROBENZENESULFONYL CHLORIDE P (M) | 0.990 | 0.114 | 0.114 | 0.458           | 98-74-8   |
| NITROBENZYL ALCOHOL P (M)           | 0.990 | 0.149 | 0.149 | 0.356           | 619-73-8  |
| NITROBIPHENYL,4-                    | 0.976 | 0.044 | 0.046 | 0.075           | 92-93-3   |
| NITROCELLULOSE (M)                  | 0.990 | 0.000 | 0.000 | 0.558           | 9004-70-0 |
| NITROETHANE                         | 0.225 | 0.412 | 0.964 | 0.161           |           |
| NITROGEN MUSTARD N-OXIDE (M)        | 0.990 | 0.139 | 0.139 | 0.794           | 126-85-2  |
| NITROMETHANE                        | 0.990 | 0.255 | 0.954 | 0.883           | 75-52-5   |
| NITROMETHYLBENZENE                  | 0.990 | 0.463 | 0.570 | 0.270           |           |
| NITROPROPANE 2                      | 0.985 | 0.531 | 0.989 | 0.437           | 79-46-9   |
| NITROBENZYL ALCOHOL 4 (M)           | 0.901 | 0.405 | 0.405 | 0.136           |           |
| NITROSOPYRROLIDINE N (M)            | 0.990 | 0.997 | 0.997 | 0.980           | 930-55-2  |
| NITROTOLUENE (-p)                   | 0.990 | 0.339 | 0.417 | 0.45199-        | 99-0      |
| NITROTOLUENE, m                     | 0.990 | 0.475 | 0.585 | 0.279           |           |
| NITROTOLUENE, o                     | 0.990 | 0.534 | 0.657 | 0.296           |           |
| NITROTOLUENE, o                     | 0.988 | 0.534 | 0.657 | 0.266           |           |
| NONANAL                             | 0.990 | 0.938 | 0.959 | 0.558           |           |
| NONANOL, n                          | 0.856 | 0.099 | 0.103 | 0.091           |           |
| NONYLPHENOL(M)                      | 0.990 | 0.115 | 0.115 | 0.794           | 25154-52  |
| OCTAMETHYLPYRROPHOSPHORAMIDE (M)    | 0.990 | 0.082 | 0.082 | 0.980           | 152-16-9  |
| OCTANAL                             | 0.990 | 0.946 | 0.979 | 0.465           |           |
| OCTANE                              | 0.990 | 1.086 | 1.000 | 0.980           | 111-65-9  |
| OCTANOL 1                           | 0.990 | 0.184 | 0.195 | 0.240           | 111-87-5  |
| OCTANOL 2                           | 0.983 | 0.381 | 0.398 | 0.136           |           |
| OCTANOL 3                           | 0.990 | 0.514 | 0.536 | 0.104           |           |
| OCTANOL 4                           | 0.990 | 0.446 | 0.466 | 0.118           |           |
| OIL (decane)                        | 0.990 | 1.088 | 1.000 | 0.951           |           |
| OXAMIC ACID (M)                     | 0.990 | 0.999 | 0.999 | 0.317           | 471-47-6  |
| PARABROMOPHENOL (M)                 | 0.925 | 0.139 | 0.139 | 0.135           | 106-41-2  |
| PARAFORMALDEHYDE (M)                | 0.990 | 0.000 | 0.000 | 0.558           | 30525-89- |
| PARALDEHYDE                         | 0.795 | 0.717 | 0.991 | 0.232           | 123-63-7  |
| PCB 1016 (monochlorobiphenyl)       | 0.990 | 1.204 | 1.000 | 0.345           | 12674-11- |
| PCB 1221 (monochlorobiphenyl)       | 0.990 | 1.204 | 1.000 | 0.418           | 11104-28- |
| PCB 1232 (dichlorobiphenyl)         | 0.990 | 1.177 | 1.000 | 0.543           | 11141-16- |
| PCB 1242 (trichlorobiphenyl)        | 0.990 | 1.075 | 0.929 | 0.488           | 53469-21- |
| PCB 1248 (quatrochlorobiphenyl)     | 0.990 | 1.142 | 1.000 | 0.640           | 12672-29- |
| PCB 1254(pentachlorobiphenyl)       | 0.990 | 0.698 | 0.618 | 0.813           | 11097-69- |
| PCB 1260 (hexachlorobiphenyl)       | 0.990 | 0.504 | 0.450 | 0.791           | 11096-82- |
| PCB'S (Aroclors)                    | 0.990 | 1.142 | 1.000 | 0.507           |           |
| PENTACHLOROBENZENE                  | 0.990 | 1.091 | 1.000 | 0.796           | 608-93-5  |
| PENTACHLOROETHANE                   | 0.990 | 0.991 | 0.966 | 0.877           | 76-01-7   |
| PENTACHLORONITROBENZENE             | 0.990 | 0.774 | 0.839 | 0.405           |           |
| PENTACHLOROPHENOL                   | 0.990 | 0.092 | 0.090 | 0.298           | 87-86-5   |
| PENTADIENE 1,2                      | 0.990 | 1.191 | 1.000 | 0.855           |           |
| PENTAERYTHRITOL TETRANITRATE (M)    | 0.976 | 0.067 | 0.067 | 0.162           | 78-11-5   |
| PENTANAL                            | 0.990 | 0.904 | 0.999 | 0.406           |           |
| PENTANE                             | 0.990 | 1.082 | 1.000 | 0.925           |           |
| PENTYL PROPANOATE                   | 0.990 | 0.868 | 1.000 | 0.537           |           |
| PENTYLAMINE                         | 0.903 | 0.822 | 0.917 | 0.254           |           |
| PENTYLBENZENE                       | 0.990 | 1.173 | 1.000 | 0.766           |           |
| PENTYLCYCLOPENTANE                  | 0.990 | 1.103 | 1.000 | 0.980           |           |
| PERCHLOROMETHYL MERCAPTAN (M)       | 0.990 | 0.132 | 0.132 | 0.980           | 594-42-3  |
| PERYLENE (M)                        | 0.990 | 0.099 | 0.099 | 0.853           | 198-55-0  |
| PHENANTHRENE                        | 0.990 | 0.279 | 0.222 | 0.193           | 85-01-8   |
| PHENOL,3-(1,1-DIMETHYLETHYL)-(M)    | 0.990 | 0.558 | 0.558 | 0.794           | 585-34-2  |
| PHENOTHIAZINE (M)                   | 0.990 | 0.125 | 0.125 | 0.874           | 92-84-2   |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                             | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|--------------------------------------|-------|-------|-------|-----------------|-----------|
| PHENYL ISOCYANATE (M)                | 0.990 | 0.674 | 0.674 | 0.855           | 103-71-9  |
| PHENYLACETIC PEROXIDE (M)            | 0.917 | 0.149 | 0.149 | 0.137           |           |
| PHENYLCYCLOHEXANONE 4                | 0.990 | 1.029 | 0.914 | 0.826           | 4894-75-1 |
| PHENYLHYDRAZINE (M)                  | 0.990 | 0.860 | 0.860 | 0.314           | 100-63-0  |
| PHENYLPHENOL P                       | 0.990 | 0.001 | 0.001 | 0.710           | 92-69-3   |
| PHENYLTHIOUREA (M)                   | 0.990 | 0.149 | 0.149 | 0.863           | 103-85-5  |
| PHOSGENE (decomposes)                | 0.990 | 0.868 | 1.000 | 0.872           | 75-44-5   |
| PHOSPHINE                            | 0.990 | 0.213 | 1.000 | 0.996           | 7803-51-2 |
| PHTHALATE, DI N BUTYL-               | 0.971 | 0.006 | 0.006 | 0.095           |           |
| PHTHALATE, DI N OCTYL                | 0.990 | 0.042 | 0.044 | 0.574           |           |
| PHTHALIC ACID                        | 0.990 | 0.714 | 0.924 | 0.858           | 88-99-3   |
| PHTHALIMIDE                          | 0.990 | 0.850 | 0.957 | 0.854           | 85-41-6   |
| PICOLINE(2-) (M)                     | 0.990 | 0.999 | 0.999 | 0.398           | 109-06-8  |
| PINENE(alpha-)                       | 0.990 | 1.165 | 1.000 | 0.890           | 80-56-8   |
| PIPERAZINE                           | 0.990 | 0.031 | 0.042 | 0.339           | 110-85-0  |
| POLYCYCLIC KETONE O (M)              | 0.990 | 0.000 | 0.000 | 0.948           |           |
| PROPANAL                             | 0.902 | 0.813 | 1.000 | 0.436           |           |
| PROPANE                              | 0.990 | 1.075 | 1.000 | 0.880           | 74-98-6   |
| PROPANE), 2,2'-OXYBIS(2-CHLORO-(M)   | 0.990 | 0.138 | 0.138 | 0.980           | 39638-32  |
| PROPANOIC ACID                       | 0.104 | 0.105 | 0.163 | 0.064           | 79-09-4   |
| PROPANOL                             | 0.595 | 0.305 | 0.421 | 0.185           |           |
| PROPANOL ISO                         | 0.451 | 0.740 | 0.926 | 0.190           | 67-63-0   |
| PROPENAL                             | 0.943 | 0.855 | 1.000 | 0.487           |           |
| PROPENE                              | 0.990 | 1.144 | 1.000 | 0.980           |           |
| PROPENYL BENZENE                     | 0.990 | 1.217 | 1.000 | 0.860           |           |
| PROPIONALDEHYDE                      | 0.990 | 0.813 | 0.999 | 0.406           | 123-38-6  |
| PROPIONIC ACID                       | 0.990 | 0.066 | 0.102 | 0.381           | 79-09-4   |
| PROPIONITRILE (M)                    | 0.990 | 0.999 | 0.999 | 0.580           | 107-12-0  |
| PROPYL ACETATE ISO                   | 0.990 | 0.786 | 1.000 | 0.453           | 108-21-4  |
| PROPYL BUTANOATE                     | 0.990 | 0.843 | 1.000 | 0.475           |           |
| PROPYL ETHER                         | 0.990 | 0.921 | 1.000 | 0.716           | 111-43-3  |
| PROPYL METHANOATE                    | 0.990 | 0.714 | 1.000 | 0.506           |           |
| PROPYL PROPANOATE                    | 0.990 | 0.813 | 1.000 | 0.446           |           |
| PROPYL THIOURACIL (M)                | 0.990 | 0.140 | 0.140 | 0.921           | 51-52-5   |
| PROPYL(-n) ACETATE                   | 0.990 | 0.773 | 0.999 | 0.448           | 109-60-4  |
| PROPYL(-n) BENZENE                   | 0.990 | 1.191 | 1.000 | 0.781           | 103-65-1  |
| PROPYL-3-METHOXY PYRAZINE,2-ISO(M)   | 0.990 | 0.149 | 0.149 | 0.980           | 25773-40- |
| PROPYLAMINE                          | 0.563 | 0.778 | 0.971 | 0.249           | 107-10-8  |
| PROPYLCYCLOPENTANE                   | 0.990 | 1.105 | 1.000 | 0.980           |           |
| PROPYLENE                            | 0.990 | 1.144 | 1.000 | 0.980           | 115-07-1  |
| PROPYLENE CHLOROHYDRIN               | 0.274 | 0.338 | 0.383 | 0.069           |           |
| PROPYLENE OXIDE                      | 0.990 | 0.841 | 1.000 | 0.600           | 75-56-9   |
| PROPYLENIMINE 1,2 2 methyl aziri     | 0.609 | 0.792 | 0.944 | 0.239           | 75-55-8   |
| PROPYN-1-OL 2(PROPARLGLYL)           | 0.550 | 0.271 | 0.321 | 0.225           | 107-19-7  |
| PROPYNE                              | 0.990 | 1.200 | 1.000 | 0.853           |           |
| PYRENE                               | 0.990 | 0.046 | 0.036 | 0.113           | 129-00-0  |
| PYRIDINE                             | 0.956 | 0.608 | 0.600 | 0.255           | 110-86-1  |
| PYRROLIDINE                          | 0.198 | 0.814 | 0.936 | 0.072           |           |
| QUINALDINE (M)                       | 0.990 | 0.999 | 0.999 | 0.853           | 91-63-4   |
| RESERPINE (M)                        | 0.990 | 0.000 | 0.000 | 0.648           | 50-55-5   |
| s ACETYLMERCAPTOSUCCINIC ACID        | 0.318 | 0.030 | 0.050 | 0.069           |           |
| S4CHL.CYCLOHEX.00DIMETH .PHOS.DIT(M) | 0.990 | 0.052 | 0.052 | 0.342           |           |
| SACCHARIN (M)                        | 0.990 | 0.133 | 0.133 | 0.850           | 81-07-2   |
| SAFROLE (M)                          | 0.990 | 0.144 | 0.144 | 0.406           | 94-59-7   |
| sec BUTYLBENZENE                     | 0.990 | 1.187 | 1.000 | 0.860           |           |
| SILVEX                               | 0.990 | 1.106 | 1.000 | 0.774           | 93-72-1   |
| SODIUM DODECYL SULFATE (M)           | 0.988 | 0.081 | 0.081 | 0.195           | 151-21-3  |
| SODIUM DODECYLBENZENE SULFONATE (M)  | 0.908 | 0.083 | 0.083 | 0.121           | 25155-30- |
| STREPTOZOTOCIN (M)                   | 0.990 | 0.092 | 0.092 | 0.980           | 18883-66- |
| STYRENE                              | 0.990 | 1.229 | 1.000 | 0.800           | 100-42-5  |
| STYRENE OXIDE                        | 0.990 | 0.883 | 0.830 | 0.341           |           |
| SULFIDE (M)                          | 0.990 | 0.999 | 0.999 | 0.649           |           |
| TAMARON (METHAMIDIPHOS)              | 0.306 | 0.430 | 0.672 | 0.091           |           |
| TARS(M)                              | 0.990 | 0.025 | 0.025 | 0.642           |           |
| t-BUTYL HYDROPEROXIDE                | 0.497 | 0.289 | 0.404 | 0.199           | 75-91-2   |
| TERPINEOL, ALPHA                     | 0.990 | 1.008 | 0.984 | 0.473           |           |
| tert BUTANOL                         | 0.630 | 0.856 | 0.989 | 0.231           |           |
| tert-AMYL BENZENE                    | 0.990 | 1.173 | 1.000 | 0.870           |           |
| tert-BUTYLBENZENE                    | 0.990 | 1.192 | 1.000 | 0.855           |           |
| TETRACHLOROQUINONE (M)               | 0.990 | 0.102 | 0.102 | 0.980           |           |
| TETRACHLOROBENZENE(1,2,3,4)          | 0.990 | 1.101 | 1.000 | 0.700           | 634-66-2  |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                              | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|---------------------------------------|-------|-------|-------|-----------------|-----------|
| TETRACHLOROBENZENE(1,2,3,5)           | 0.990 | 1.101 | 1.000 | 0.732           | 634-90-2  |
| TETRACHLOROBENZENE(1,2,4,5)           | 0.990 | 1.101 | 1.000 | 0.732           | 95-94-3   |
| TETRACHLORODIBENZOFURAN (2,3,7,8) (M) | 0.990 | 0.072 | 0.072 | 0.332           | 51207-31- |
| TETRACHLORODIBENZO-p-DIOXIN(2,3,7,8)  | 0.990 | 0.109 | 0.101 | 0.173           | 1746-01-6 |
| TETRACHLOROETHANE(1,1,1,2) (M)        | 0.990 | 0.141 | 0.141 | 0.459           | 630-20-6  |
| TETRACHLOROETHANE(1,1,2,2)            | 0.990 | 1.015 | 0.999 | 0.397           | 79-34-5   |
| TETRACHLOROETHENE                     | 0.990 | 1.048 | 1.000 | 0.917           | 127-18-4  |
| TETRACHLOROPHENOL(2,3,4,6)            | 0.447 | 1.024 | 1.000 | 0.091           | 58-90-2   |
| TETRACHLOROPHENOL(2,3,5,6)            | 0.990 | 0.010 | 0.010 | 0.980           | 935-95-5  |
| TETRACHLOROPROPENE(1,1,2,3) (M)       | 0.990 | 0.135 | 0.135 | 0.831           | 10436-39- |
| TETRADECANE                           | 0.990 | 1.089 | 1.000 | 0.896           | 629-59-4  |
| TETRAETHYL LEAD                       | 0.990 | 0.958 | 0.889 | 0.980           | 78-00-2   |
| TETRAETHYLENE GLYCOL (M)              | 0.892 | 0.128 | 0.128 | 0.117           | 112-60-7  |
| TETRAETHYLENE PENTANE                 | 0.990 | 1.183 | 1.000 | 0.881           |           |
| TETRAETHYLPYROPHOSPHATE (M)           | 0.990 | 0.080 | 0.080 | 0.980           | 107-49-3  |
| TETRAFLUROETHENE                      | 0.990 | 1.080 | 1.000 | 0.980           |           |
| TETRAFLUOROMETHANE                    | 0.990 | 1.037 | 1.000 | 0.980           |           |
| TETRAHYDROBENZALDEHYDE                | 0.912 | 0.635 | 0.641 | 0.213           |           |
| TETRAHYDROFURAN                       | 0.830 | 0.860 | 1.000 | 0.322           | 109-99-9  |
| TETRAHYDRONAPHTHALENE,1,2,3,4-(M)     | 0.887 | 0.452 | 0.452 | 0.794           | 119-64-2  |
| TETRAHYDROPYRAN                       | 0.980 | 0.898 | 1.000 | 0.381           | 142-68-7  |
| TETRAHYDROTHIOPHENE                   | 0.990 | 0.692 | 1.000 | 0.566           |           |
| TETRALIN                              | 0.990 | 1.189 | 1.000 | 0.632           |           |
| TETRANITROMETHANE                     | 0.990 | 0.267 | 1.000 | 0.852           | 509-14-8  |
| THIOACETAMIDE (M)                     | 0.990 | 0.999 | 0.999 | 0.375           | 62-55-5   |
| THIOBENZYL ALCOHOL P (M)              | 0.887 | 0.588 | 0.588 | 0.136           | 100-53-8  |
| THIOBISETHANE, 1,1'                   | 0.990 | 0.692 | 1.000 | 0.763           |           |
| THIOCYANATE (TOTAL AS SCN-) (M)       | 0.990 | 0.642 | 0.642 | 0.894           | NA        |
| THIOMETHANOL (M)                      | 0.990 | 0.999 | 0.999 | 0.499           | 74-93-1   |
| THIOPHENOL (M)                        | 0.659 | 0.826 | 0.826 | 0.933           | 108-98-5  |
| THIOPROPIONAMIDE 2 (M)                | 0.696 | 0.948 | 0.948 | 0.097           |           |
| THIOUREA                              | 0.892 | 0.011 | 0.024 | 0.472           | 62-56-6   |
| THIRAM (M)                            | 0.990 | 0.105 | 0.105 | 0.980           | 137-26-8  |
| THYMINE (M)                           | 0.990 | 0.556 | 0.556 | 0.806           | 65-71-4   |
| TOLUENE                               | 0.990 | 1.215 | 1.000 | 0.804           | 108-88-3  |
| TOLUENE2,4DIAZOBIS-METATOLUENEDIA(M)  | 0.986 | 0.011 | 0.011 | 0.188           |           |
| TOLUENESULFONYL CHLORIDE              | 0.604 | 0.046 | 0.047 | 0.068           |           |
| TOLUIC ALDEHYDE                       | 0.990 | 0.513 | 0.478 | 0.382           | 122-78-1  |
| TOLUIDINE (-0)                        | 0.459 | 0.159 | 0.152 | 0.052           | 95-53-4   |
| TOLUIDINE HYDROCHLORIDE,o-(M)         | 0.990 | 0.258 | 0.258 | 0.980           | 636-21-5  |
| TOLUIDINE P                           | 0.850 | 0.274 | 0.262 | 0.208           | 106-49-0  |
| TOXAPHENE                             | 0.990 | 0.054 | 0.050 | 0.735           | 8001-35-2 |
| trans 1,4 DIMETHYLCYCLOHEXANE         | 0.990 | 1.117 | 1.000 | 0.980           |           |
| trans 2 BUTENAL                       | 0.387 | 0.911 | 1.000 | 0.267           |           |
| trans 2 HEPTENE                       | 0.990 | 1.121 | 1.000 | 0.980           |           |
| trans 2 HEXENAL                       | 0.856 | 0.963 | 1.000 | 0.295           |           |
| trans 2 OCTENAL                       | 0.990 | 0.985 | 0.993 | 0.381           |           |
| trans, trans 2,4 HEXADIENAL           | 0.233 | 0.996 | 1.000 | 0.151           |           |
| TRIBROMOMETHYLPHOSPHATE (M)           | 0.980 | 0.052 | 0.052 | 0.169           |           |
| TRIBUTYL PHOSPHOROTRITHIOATE SSS      | 0.990 | .     | .     | 0.334           | 78-48-8   |
| TRIBUTYL TIN ACETATE                  | 0.990 | 0.929 | 0.980 | 0.789           |           |
| TRIBUTYLPHOSPHATE                     | 0.990 | 1.073 | 0.988 | 0.980           | 126-73-8  |
| TRICHLORO(1,1,2)TRIFLUOROETHANE(M)    | 0.990 | 0.131 | 0.131 | 0.980           | 76-13-1   |
| TRICHLORO-1,2,2-TRIFLUOROETHANE,1,1   | 0.990 | 1.033 | 1.000 | 0.980           | 76-13-1   |
| TRICHLORO-1,3,5-TRIAZINE 2,4,6 (M)    | 0.990 | 0.133 | 0.133 | 0.552           | 108-77-0  |
| TRICHLOROANISOLE 2,3,6 (M)            | 0.990 | 0.119 | 0.119 | 0.980           | 50375-10- |
| TRICHLOROBENZENE 1,2,3                | 0.990 | 1.114 | 1.000 | 0.808           | 87-61-6   |
| TRICHLOROBENZENE 1,2,4                | 0.990 | 1.114 | 1.000 | 0.637           | 120-82-1  |
| TRICHLOROBENZENE 1,3,5                | 0.990 | 1.114 | 1.000 | 0.877           | 108-70-3  |
| TRICHLOROBUTANE 1,2,3 (M)             | 0.990 | 0.144 | 0.144 | 0.980           | 18338-40- |
| TRICHLOROETHANE 1,1,1                 | 0.990 | 1.037 | 1.000 | 0.913           | 71-55-6   |
| TRICHLOROETHANE 1,1,2                 | 0.990 | 1.025 | 1.000 | 0.597           | 79-00-5   |
| TRICHLOROETHYLENE                     | 0.990 | 1.053 | 1.000 | 0.866           | 79-01-6   |
| TRICHLOROFLUOROMETHANE                | 0.990 | 1.027 | 1.000 | 0.968           | 75-69-4   |
| TRICHLOROPHENOL 2,4,5                 | 0.964 | 0.111 | 0.108 | 0.086           | 95-95-4   |
| TRICHLOROPHENOL 2,4,6                 | 0.990 | 0.135 | 0.132 | 0.167           | 88-06-2   |
| TRICHLOROPROPANE 1,1,1                | 0.990 | 1.048 | 1.000 | 0.897           | 7789-89-1 |
| TRICHLOROPROPANE(1,1,2)               | 0.990 | 1.037 | 1.000 | 0.897           | 598-77-6  |
| TRICHLOROPROPANE(1,2,2)               | 0.990 | 1.047 | 1.000 | 0.897           | 3175-23-3 |
| TRICHLOROPROPANE(1,2,3)               | 0.990 | 1.048 | 1.000 | 0.894           | 96-18-4   |
| TRICHLOROPROPENE (1,1,2)(M)           | 0.990 | 0.228 | 0.228 | 0.795           |           |

TABLE 2 OF APPENDIX J.—FR, FM, AND FE<sup>1</sup> VALUES FOR COMPOUNDS WITH HENRY'S LAW CONSTANTS AT 25° C GREATER THAN OR EQUAL TO 0.1 Y/X ATMOSPHERE PER MOLE FRACTION—Continued

| Compound                            | FR    | Fm25D | Fm305 | Fe <sup>1</sup> | CAS       |
|-------------------------------------|-------|-------|-------|-----------------|-----------|
| TRICOSANE N(M)                      | 0.990 | 0.133 | 0.133 | 0.301           | 629-50-5  |
| TRIETHYLAMINE                       | 0.990 | 0.937 | 1.000 | 0.379           | 121-44-8  |
| TRIETHYLENE GLYCOL(M)               | 0.846 | 0.150 | 0.150 | 0.111           | 112-27-6  |
| TRIETHYLPHOSPHOROTHIOAT E,o,o,o-(M) | 0.989 | 0.126 | 0.126 | 0.794           | 126-68-1  |
| TRIFLUOROETHANE(1,1,1)              | 0.990 | 1.059 | 1.000 | 0.980           |           |
| TRIFLUOROMETHANE                    | 0.990 | 1.057 | 1.000 | 0.980           |           |
| TRIFLURALIN                         | 0.990 | 0.086 | 0.116 | 0.291           |           |
| TRIIISOBUTYLENE                     | 0.990 | 1.117 | 1.000 | 0.980           |           |
| TRIIISOPROPYLAMINE                  | 0.990 | 1.026 | 1.000 | 0.715           |           |
| TRIMELLITIC ANHYDRIDE (M)           | 0.629 | 0.129 | 0.129 | 0.087           | 552-30-7  |
| TRIMETHYL BENZENE, 123              | 0.990 | 1.200 | 1.000 | 0.713           |           |
| TRIMETHYL-4-NITROANILINE 2,3,5 (M)  | 0.990 | 0.135 | 0.135 | 0.831           |           |
| TRIMETHYLAMINE                      | 0.990 | 0.811 | 1.000 | 0.464           | 75-50-3   |
| TRIMETHYLBENZENE (1,3,5)            | 0.990 | 1.200 | 1.000 | 0.766           | 108-67-3  |
| TRIMETHYLPENTANE 2,2,4              | 0.990 | 1.116 | 1.000 | 1.000           | 540-84-1  |
| TRIMETHYLSILANOL                    | 0.990 | 0.533 | 1.000 | 0.980           |           |
| TRINITROBENZENE,sym-(M)             | 0.990 | 0.118 | 0.118 | 0.712           | 99-35-4   |
| TRINITROTOLUENE(2,4,6)              | 0.223 | 0.004 | 0.009 | 0.120           | 118-96-7  |
| TRIPHENYL PHOSPHINE (M)             | 0.990 | 0.094 | 0.094 | 0.321           | 603-35-0  |
| TRIPHENYLMETHANE (M)                | 0.990 | 0.103 | 0.103 | 0.980           | 516-73-3  |
| TRIPHENYLPHOSPHINE NICKEL CARBONM)  | 0.990 | 0.037 | 0.037 | 0.722           |           |
| TRIS (1-AZIRIDINYL) PHOSPHINESU(M)  | 0.990 | 0.130 | 0.130 | 0.379           | 52-24-4   |
| TRIS (2,3-DIBROMOPROPYL)PHOSPHA(M)  | 0.990 | 0.000 | 0.000 | 0.980           | 126-72-7  |
| TRISODIUM NITRILOTRIACETATE (M)     | 0.990 | 0.128 | 0.128 | 0.980           | 5064-31-3 |
| TRYPAN BLUE(M)                      | 0.990 | 0.000 | 0.000 | 0.853           | 72-57-1   |
| URACIL (M)                          | 0.990 | 0.794 | 0.794 | 0.857           | 66-22-8   |
| URACIL MUSTARD (M)                  | 0.990 | 0.099 | 0.099 | 0.853           | 66-75-1   |
| UREA                                | 0.990 | 0.016 | 0.030 | 0.582           | 57-13-6   |
| URETHANE                            | 0.990 | 0.024 | 0.039 | 0.370           | 51-79-6   |
| VALERIC ACID (M)                    | 0.990 | 0.963 | 0.963 | 0.287           | 109-52-4  |
| VINYL ACETATE                       | 0.990 | 0.748 | 1.000 | 0.592           | 108-05-4  |
| VINYL ACETYLENE                     | 0.990 | 1.232 | 1.000 | 0.890           |           |
| VINYL BROMIDE                       | 0.990 | 0.629 | 1.000 | 0.849           |           |
| VINYL CHLORIDE                      | 0.990 | 1.081 | 1.000 | 0.971           | 75-01-4   |
| VINYL DIHYDROPYRAN                  | 0.990 | 0.935 | 1.000 | 0.554           |           |
| VINYL METHYL ETHER                  | 0.990 | 0.831 | 1.000 | 0.590           |           |
| VINYLCYCLOHEXENE 4(M)               | 0.990 | 0.860 | 0.860 | 0.980           | 100-40-3  |
| VINYLDENE CHLORIDE                  | 0.990 | 1.061 | 1.000 | 0.889           | 75-35-4   |
| XYLENE                              | 0.990 | 1.206 | 1.000 | 0.788           | 1330-20-7 |
| XYLENE(-m)                          | 0.990 | 1.206 | 1.000 | 0.821           | 108-38-3  |
| XYLENE(-o)                          | 0.990 | 1.206 | 1.000 | 0.787           | 95-47-6   |
| XYLENE(-p)                          | 0.990 | 1.206 | 1.000 | 0.824           | 106-67-9  |
| XYLIDINE dimethylaniline            | 0.606 | 0.131 | 0.124 | 0.074           |           |
| XYLYL CHLORIDE M (M)                | 0.990 | 0.310 | 0.310 | 0.592           | 620-19-9  |
| XYLYL CHLORIDE O (M)                | 0.990 | 0.310 | 0.310 | 0.592           | 552-45-4  |

\* Molecular structure only approximate.

(M) fraction measured (fm) estimated from Mwt correlation.

<sup>1</sup> The Fe values listed in Table 2 are Fe values for emissions from both the individual drain system and the treatment process. Use these Fe values with Section 2.5.1).

TABLE 3 OF APPENDIX J—FE VALUES FOR EMISSIONS FROM BOTH THE INDIVIDUAL DRAIN SYSTEM AND THE TREATMENT PROCESS

[Use with section 2.5.1]

| Henry's Law Constant | Fe Value |
|----------------------|----------|
| 0.00025              | 0.001    |
| 0.00051              | 0.002    |
| 0.00076              | 0.003    |
| 0.00127              | 0.005    |
| 0.00178              | 0.007    |
| 0.00254              | 0.010    |
| 0.00381              | 0.015    |
| 0.00508              | 0.020    |
| 0.00635              | 0.25     |
| 0.00762              | 0.030    |
| 0.00890              | 0.035    |
| 0.01017              | 0.040    |
| 0.01144              | 0.045    |
| 0.02327              | 0.050    |