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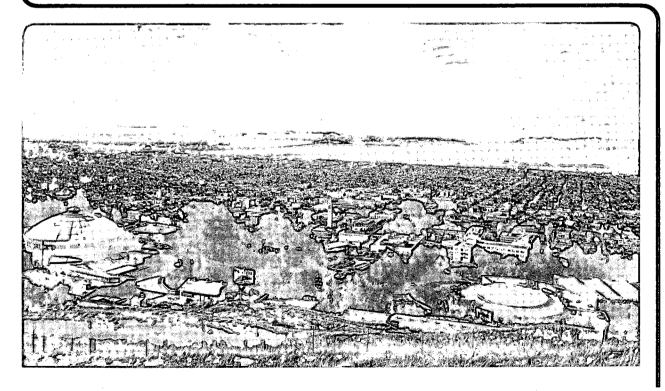
Groupings of Organic Waste Chemicals Based on Sorption, Biotransformation and Hydrolysis at Standard Conditions for Application to the Deep Subsurface Environment

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January 1988

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Groupings of Organic Waste Chemicals Based on Sorption, Biotransformation and Hydrolysis at Standard Conditions for Application to the Deep Subsurface Environment

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ABSTRACT

Chemical and biological reactions of organic hazardous wastes disposed to injection wells at depths down to 7000 feet deep are assessed. At these depths, the major reactions involving organic wastes include hydrolysis, biotransformation and sorption. However, experimental data on reactions of organics in the deep subsurface are sparse or nonexistent.

Tables are given on values for hydrolysis, sorption and biotransformation reactions obtained from research publications, mostly at 25 °C and atmospheric pressure. It is suggested that the more plentiful data on reactions at the land surface be used to approximate the rates of hydrolysis and biotransformation, and sorption equilibrium by taking into account the expected subsurface environment. For example, predictive methods will take into consideration a deep subsurface environment which has: a higher temperature and pressure than the land surface; fewer kinds of microorganisms and a substantially lower concentration of microorganisms; lower organic carbon levels; and, highly saline ground waters.

1.0 INTRODUCTION

Injection wells are used to isolate organic liquid wastes resulting from industrial processes by pumping the liquids into carefully selected subsurface geologic formations (1-3.57). This is an attractive method for disposal of hazardous liquid organic wastes because the injected material is efficiently removed from man's immediate air, water and land environment. Other beneficial aspects sought are the isolation of these organic wastes for geologic times (56), and eventual transformation into stable products which might include inorganic substances, carbon dioxide and water (4). The method is economical when compared with other optional procedures, especially those where pretreatment is required (2). Most injection wells were drilled in the mid-1960's and mid-1970's; there has been no significant increase since 1980 (57, p.6). Injection of hazardous wastes is not a widespread practice (57, p.11). However, the question has been raised, can these wastes enter our source of underground drinking water by migration from the disposal well sites. Implicit in this question is the time required for the waste constituents to travel by groundwater movement or by diffusion, because it is likely that the original organic substances will undergo reactions outside the immediate disposal well system.

Understanding the reactions (fate) of organic chemicals disposed to the deep subsurface environment at depths of 1000 to 7000 feet is important to their management and regulation (13). These reactions are best understood by experimental measurements of chemical reactions most likely to occur in a heterogeneous system consisting of minerals, organic matter and ground water. However, little information is available for all hazardous organic substances and their reactions below about 10 meters, so that predictive methods are mandatory. Predictive methods ideally would use the plentiful data compiled from laboratory measurements or calculations for 25 °C and for pressures near one bar. These values of equilibrium constants and kinetic rate constants can then be extrapolated to conditions which reflect a deep subsurface environment using theoretically based methods and correlations, to account for the effects of differing temperatures and pressures.

The data in this compilation were selected for use in a model for reactions between a generic waste injection site comprised of an immobile mineral/organic carbon solid

phase; an aqueous liquid phase; and, simple organic wastes which have been injected into this subsurface environment. The reactions and associated data considered are limited to sorption of the waste constituents onto the organic carbon (humus) portion of a mineral/soil solid phase (13-25); biotransformation by microorganisms, principally bacterial (4-13,24-32); solubility (35,36,38); Henry's constants (39-41); and, hydrolysis (34,35,37). In an undisturbed subsurface environment, the number of microorganisms is expected to be substantially less than that at the surface and in surficial soils (5,30,31). Injection of the wastes could cause an increase in both the population of microorganisms and the organic content of the soil. However, the established soil microflora are likely to resist introduction of new microorganisms (43,p.22). There is less organic material bound to the minerals (43,46), so that adsorption onto mineral surfaces becomes important (25,p.319) when considering retardation mechanisms. Lastly, at these depths both the ambient temperature (45) and pressure will differ from that prevailing at the surface; both are higher, and have a major effect on homogeneous and heterogeneous reactions involving injected organic material. The model for the injection process is depicted in Figure 1.

Nonchemical removal of the various waste constituents from a containment site by diffusion, or migration by movement of groundwater are excluded from this report. Sequential reactions of an organic waste such as biodegradation of a product formed from hydrolysis are not considered; only the initial reactions are covered. Calculations on the fate of organic wastes in the deep subsurface is outside the scope of this report.

Additional information on reactions of organic materials in the soil/water environment near the land surface can be obtained from References 5 and 15.

1.1 Organic Wastes Constituents

The individual organic constituents in this grouping were obtained from the Proposed Grouping of Listed U and P Waste Codes (49). The result is a tabulation for about 300 organic chemicals representing both liquid and solid forms of the pure chemicals. Some

Ideal Injection Well and Site

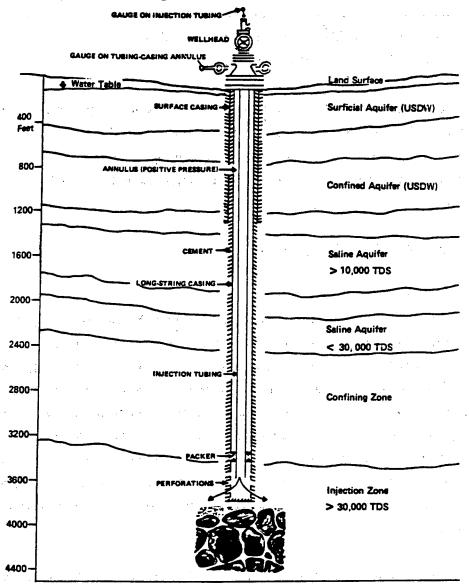


Figure 1. A model of the injection environment (57). At land surface: organic carbon (OC) is 0.06 to 2% w (assumed bound to minerals, see Refs. 15 and 20); microorganisms (Ref. 43) are 10^6 to 10^7 bacteria per gram of soil (Ref. 5), other microorganisms such as 1,000,000 to 20,000,000 actinomycetes; 5000 to 900,000 fungi; 1,000 to 100,000 yeasts; 1,000 to 500,000 algae; 1,000 to 500,000 protozoa. Temperature is variable. In the deep subsurface environment, such as 4,000 to 7,000 feet; it can be assumed: organic carbon <0.06%; microorganisms (principally bacterial) << 10^6 cells per gram; temperatures exceed about 35 °C. However, aerobic bacteria have been found beneath the subsurface at the Savannah River Laboratory (58).

of these chemicals are identified as constituents in hazardous wastes from both specific and non-specific sources. These sources have an industry and EPA hazardous waste number (49) which is included in this tabulation.

Table 1-1 lists selected organic wastes that have been injected into deep geologic formations (2,3,13). These wastes from industrial operations were classified by Donaldson into 15 categories of water-soluble oils, and two classes of water-insoluble oils (2). More recent information is found in the comprehensive report of Reference 57.

1.2 Deep Subsurface Soil/Water Environment

A description of selected injection wells was tabulated by Donaldson up to about the year 1971 (2). The listing covered geologic formations at depths ranging from 1500 feet to 7200 feet (300 to 2400 meters). These formations were mostly comprised of three types of materials: unconsolidated sands, sandstone and a vugular carbonate (limestone and dolomite). The disposal sites were selected because they were porous, permeable, and had a large areal extent. Types of formations and their geological properties are given in Tables 1-2 and 1-3. It should be noted that injection of the wastes displaces the natural pore waters from these wells (2, p.34). The properties of the formation soils are important in containing the organic wastes, for example sorption capacity for individual organic wastes.

Unconsolidated sands in the Frio formation of Harris County, Texas are medium to fine-grained, loosely cemented and calcareous (50). The components of the sands varied with depth; however, quartz and smectite (montmorillonite) clays and feldspars were usually present between 5709 - 6290 feet depths. The clay content measured as the methylene blue capacity varied from 2 to 48 me/100 g (50).

Most sandstones are comprised of a small number of minerals; an average sandstone has 66.8% SiO₂ (primarily quartz); 11.5% feldspar; 11.1% carbonate minerals; 6.6% micas and clays; 1.8% iron oxides; and, 2% other minerals (13,p.16). Limestone is mostly CaCO₃ (calcite), but can contain a significant amount of noncarbonate constituents such as clay minerals (13,p.15).

Clay minerals are important because they represent a highly reactive and large surface area (13,25,55): about 500 m²/g for montmorillonite (55). This reactivity is manifested by a large capacity for adsorbing many polar organic substances (51,p.288). For example, recently Charlesworth (23) studied the interaction of clay minerals with organic nitrogen compounds and found up to 70% adsorption on the basis of total nitrogen, by nine common clays under laboratory conditions. See Table 1-4. Clays also catalyze redox and other reactions such as cracking of hydrocarbons (55).

Experimental studies uniformly show a direct correlation between the fractional organic content of a specific soil or sediment, and the soil/water partition coefficient. These studies are limited to surface samples which were obtained within about 9 meters (30 feet) beneath the surface. Organic matter due to decaying plant and animal life (humus) concentrates in this surface layer: "the concentration decreases abruptly with depth so that the deep subsurface is almost wholly mineral matter in most soils" (43,p.4). Adsorption of organic wastes and bacteria onto mineral surfaces may be more important in the deep subsurface than in shallow soil because of the dearth of organic matter. However, this accumulation at the mineral/water interface might be more reversible than adsorption at an organic carbon/water interface (52,p.2). The result will be an excess of immobile organic wastes at the mineral/water interface due to weak van der Waals or electrostatic adsorption (53). Also, any gases evolved will accumulate at this interface; the gas/water distribution is assumed to follow Henry's law. The water/soil interface represents a source of nutrients for bacteria (52,p.2).

Because oxygen is absent (30,31,54), bacterial action is anaerobic (e.g.,methanogenic). Bacteria have been found in water samples obtained from wells, and in lithologic cores; their presence in groundwater is well documented (54,p.299). See Table 1-5. Microorganisms are not evenly distributed in soils; for example, in a sandy soil "the majority of bacteria are associated with a minor fraction of the soil- organic matter..." (25,p.319).

2.0 GROUPINGS OF WASTES

Organic industrial wastes injected into the deep subsurface environment up to about the year 1970 are complicated mixtures of pure organic chemicals and inorganic salts in

water (2,13). The pH varied from 2 to highly alkaline (8% NaOH); the content of inorganic salt such as NaCl varied, as did the total amount of organic waste material. See Table 1-1. It is recognized that reactions which determine the fate of the organic content should be viewed from the perspective of one substance reacting competitively with other constituents in the waste. For example, sorption might involve adsorption and desorption of one or more organic compounds at the solid/water interface; hydrolysis rates can be affected by ionic strength and catalytic metal ions. However, in this report we have decided to focus on individual organic substances in tabulating data on reactions which determine the fate in the deep subsurface.

Organic chemicals can be organized into a number of different categories; most often this is done by functional group. In this report, we have chosen to follow the organization given in Attachment 2 of the Proposed Grouping of Listed U and P Waste Codes (47). This selection was made to ensure compatibility with that of the EPA document, and the Part 261 of the Code of Federal Regulations (49). These identify the substances as acute (P) and toxic (U) hazardous wastes (49). See the printout of the tables of data in the Appendix.

For each pure organic substance in the Appendix, the Grouping category is given, followed by the name of the organic and the date identifying when the data were tabulated. Other commonly used names for the organics are given, together with the Chemical Abstracts Services (CAS) number. For purposes of identification, we also include the Hazardous Waste Number, and an Industry/EPA Generic No. The Industry/EPA Generic No. is not, however, related to the source of the waste injected into the subsurface disposal system. The empirical chemical formula is also given. Under the heading "Description", there is a listing of selected physical properties of the pure substance, and a reference to source of the data. The physical properties are intended to provide information on the organic chemical to be used in predicting its form in the deep subsurface, e.g., gaseous, soluble in water, or solid. The general heading Retention properties is defined as the chemical and biochemical reactions which determine the fate of the organic within the injection zone. These values are all at 25°C and about 1 bar pressure.

The data in the tabulation of the Appendix are the database for laboratory conditions which are to be used in predicting the fate of organic wastes disposed to the deep subsurface by injection wells.

3.0 METHODOLOGY FOR ASSESSMENT OF FATE

The fate of an organic substance injected into the deep subsoil depends mainly on the transformation of the original material into an initial product. An example is chemical transformation by hydrolysis of a bromoalkane to form an alcohol (34). This alcohol product may then undergo further chemical or physicochemical reactions; however, these subsequent reactions are not considered in this report. Where possible, each of the reactions treated here (biotransformation, hydrolysis) is given a half-life $(t_{1/2})$ so that a numerical value can be calculated and assigned to the "fate" process. An exception is sorption which is considered to be kinetically fast, and therefore as an equilibrium process. In the absence of experimental data obtained at depth, correlations using these surface soil data are suggested for filling gaps.

3.1 Equilibrium Sorption Correlation

Sorption of organic substances onto an organic matrix such as humus materials is an important retardation process in the topsoil. The sorption is kinetically slow (17,18), but virtually irreversible (17). In topsoil, the result of this adsorption includes removal of the organic substance from the aqueous phase with a probable resulting change in reactivity of the sorbed species, e.g., biotransformation rate may change (16).

Sorption of organic substances onto organic carbon (e.g.,humus) bound to minerals has been studied extensively (14-23). Selected data obtained from soils and sediments are shown in Table 3-1. Karickhoff developed a semi-empirical estimation of sorption of hydrophobic organic substances onto soils from aqueous media (14). A key relation is the adsorption onto the organic carbon (OC) fraction of natural soils or sediments. The relationship between the concentration of sorbed material and the concentration of pollutant in the water phase for dilute solutions at equilibrium is given by eq 1.

$$\Gamma = K_p C \tag{1}$$

 Γ = concentration of organic waste sorbed onto OC

C = solution concentration of the sorbent

 K_p = sorption partition coefficient for the heterogeneous process

A definition of a dilute solution is an aqueous phase concentration less than 10⁻⁵ M in pollutant concentration, or below one-half its solubility (14). The dependence of sorption on the OC content of the soil is given by eq 2

$$K_{p} = K_{OC} OC \tag{2}$$

where OC is the weight fraction of organic carbon in a sediment or soil (16); K_{OC} is a fugacity coefficient ratio (16). Table 3-2 lists selected values of OC for differing sediments (15) and soils (20).

Such studies have shown a linear correlation between sorption of hydrophobic organic chemicals and the octanol/water partition coefficients (15,21).

3.1.1 Octanol/Water Partition Coefficient

Values of octanol/water partition coefficient are obtained from laboratory measurements using purified and carefully characterized chemicals. The relationship between these laboratory data and partitioning of an organic waste between soil and water is a linear correlation between the $\log K_{oc}$ and $\log K_{ow}$. For a water and sediment system, Karickhoff et al. found with a correlation coefficient, $r^2 = 1.00$ (15)

$$\log K_{oc} = (1.00) \log K_{ow} - 0.21$$
 (3)

Chiou et al. (21) developed a correlation applicable to the partition equilibrium of non-ionic organic compounds distributed between organic matter bound to soil and the water phase, eq 4 with $r^2 = 0.989$:

$$\log K_{om} = (0.904)\log K_{ow} - 0.779 \tag{4}$$

where, for eq 3 and eq 4

$$K_{oc}$$
 = a sorption coefficient on an organic carbon (oc) basis, equal to $\frac{K_p}{oc}$ (15)

 $K_{ow} = \text{octanol/water partition coefficient}$, often determined by shaking known volume of octanol containing small amount of solute with water, then measuring solute concentration in octanol (C^O) and water (C^W); $K_{OW} = C^O/C^W$.

 K_{om} = solute partition coefficient between soil organic phase and water. The soil organic phase is Woodburn silt loam soil: 1.9% organic matter, 68% silt, 21% clay, 9% sand, cation exchange capacity of 14 meq/100 g soil.

We have used eq 4 to calculate values of K_{om} for selected missing data in the printout in Section 7.0. In the absence of laboratory data on K_{ow} , the magnitude of K_{oc} can be estimated from the solubility, for example, for selected hydrophobic hazardous wastes (21), using eq 5. See Table 3-3.

$$\log K_{om} = -0.729 \log S + 0.001 \tag{5}$$

where S = moles/L for supercooled liquids and $r^2 = 0.996$ (21).

However, Southworth and Keller (20) found that for soils with low OC, e.g., 0.05%, correlations such as eq 5 predict lower sorption for polar substances such as phenol, because of interactions between the organic pollutant and the mineral surfaces. This is probably because sorption onto a mineral surface is less likely to be totally irreversible than sorption onto an organic carbon surface such as humus bound to the mineral.

In summary, the partitioning of hydrophobic organic wastes between water and the organic content of a soil can be correlated to laboratory measurements of the octanol/water partition coefficient. The value of K_{om} is a measure of sorption of the organic waste; and, eq 4 has been used in this work to estimate some values of K_{om} . It is suggested that eq 4 be used to calculate the magnitude of K_{om} in the deep subsurface, after adding terms or otherwise modifying the equation to account for increased

temperature and pressure in the deep subsurface.

3.2 Biofilm Model Kinetics

Bouwer and McCarty (46) described a model for the biotransformation of trace organics in subsurface ground waters according to biofilm processes. An ideal biofilm is composed of a homogeneous matrix of bacteria and extracellular polymers that bind the bacteria together, and to the subsurface organic matter, which in turn is bound to the rock. The model can be interpreted for a subsurface environment generally characterized by low substrate and nutrient concentrations, and high specific surface area. Bacteria are likely to be attached to and therefore be immobile on the deep subsurface soils and minerals. The population of native bacteria is much smaller than in a surface soil. Webster et al. (1985) found 10⁶ to 10⁷ cells per gram of subsurface soil, at depths from 2 to 9 meters. The number of cells is undoubtedly far less in the deep subsurface, from 300 to 2500 m deep, in an injection environment.

Bouwer and McCarty (46) give eq 6 as an expression for the degradation half-life, $t_{1/2}$, of the substrate:

$$t_{1/2} = \frac{\ln 2}{[(kX)/K_s]} \tag{6}$$

where k = first order reaction maximum rate constant, sec⁻¹, of substrate utilization by bacteria and X = first order reaction of organisms such as bacteria capable of degrading the organic contaminant; $K_s = first$ monod half-maximum-rate concentration.

Table 3-4 lists half-lives based on microorganism concentrations, under both aerobic and methanogenic conditions. Note that degradation in the absence of oxygen is tenfold less than under aerobic conditions. Also, the rate of biotransformation decreases by tenfold, with a tenfold decrease in organism concentration. Therefore, half-lives of biotransformation processes may be increased due to the fewer number of organisms available in the deep subsurface.

In predicting the fate of organic wastes in the deep subsurface, the half-life of organic wastes in this environment might be estimated by assuming that: (a) $t_{1/2}$ can be divided by a factor of 10 for laboratory measured aerobic biotransformations; and, by 100 for anaerobic. (b) It might also be assumed that the microorganism concentration is 0.001 mg/L or less, so that laboratory data obtained at surface conditions could be divided by 10^5 .

With regard to temperature effects on biotransformation reactions, Fuller and Warrick (43,p.25) note that "microbial decomposition of organic materials occurs most rapidly in the mesophilic (25 to 35 °C) range, higher and lower temperatures favor some organisms,... Thus, biodegradation will occur over a wide range of temperature, although it takes place more slowly near the extremes of 0 °C and 65 °C...". In the absence of other information on the kinds of microorganisms down to 5000 feet it might be assumed, as a first approximation, that data at 25 °C can be used to estimate biotransformation rates in the deep subsurface under anaerobic conditions.

3.3 Hydrolysis Half-Life in the Deep Subsurface

Hydrolysis is a chemical process in which an organic hazardous waste constituent reacts with water to form a different substance. The detailed mechanism may involve intermediate products; however, the rate of hydrolysis of a substance RX usually follows the kinetic equation (33)

$$\frac{-d\left[RX\right]}{dt} = k_h \left[RX\right] \tag{7}$$

$$\ln[RX] = -k_h t \tag{7A}$$

so that when one-half of substance RX has reacted in time $t_{1/2}$,

$$t_{1/2} = \frac{\ln 2}{k_h} \tag{7B}$$

for the simplified reaction $RX + H_2O = ROH + HX$, where X is a functional group such as Cl^- . In eq 7, k_h is a pseudo first order reaction rate constant. The reaction rate is a function of a number of parameters such as pH, temperature, ionic strength and concentration of metal ions which might catalyze the reaction. Examples of catalytic metals are ions of Cu, Co, Mn and Mg (33).

As noted, the rate of hydrolysis for organic wastes depends on a number of parameters, mainly pH, temperature, concentration of organic pollutant, functional group of the organic wastes, and pressure (33). Mabey and Mill (33) have summarized critically evaluated hydrolysis rates and half-lives at pH 7, 25 °C, atmospheric pressure, and zero ionic strength. However, in the deep subsurface, the temperature is higher and therefore the hydrolysis rates should increase. The enhanced hydrolysis rates due to temperature of the deep subsurface e.g., 55 °C, can be calculated from the equations given by Mabey and Mill (33), for example. Effects of increased pressure, catalytic effects and differences in hydrolysis rates due to ionic strength are difficult to predict; these may be negligible compared with the larger effects of increased temperature, as a first approximation.

Hydrolysis half-life is estimated according to the usual expression

$$t_{1/2} = \frac{0.693}{k_h} \tag{8}$$

In eq 8, k_h is related to pH by the following equation used by Mabey and Mill to calculate k_h

$$k_h = k_B [OH^-] + k_A [H^+] + k_N$$
 (9)

where $K_W = [OH^-][H^+]$, and k_B , k_A , and k_N are second-order rate constants for acid and base catalyzed and neutral processes, respectively (33). The constant, K_W is the ion product of water. This is equal to -13.995 at 25 °C and -13.275 at 50 °C (56). When the pH is fixed, eq 8 describes the half-life of the hydrolysis reaction. Table 3-5 summarizes selected $t_{1/2}$ values for hazardous organic wastes taken from the larger tabulation in Mabey and Mill (33).

4.0 SUMMARY AND CONCLUSIONS

The principal homogeneous and heterogeneous chemical reactions controlling the fate of organic wastes in the near surface soil/water environment include sorption, biodegradation and chemical degradation such as hydrolysis (33). For example, Sutton and Barker (42) injected about 2100 liters of an aqueous mixture containing 70 to 100 ppm of dissolved n-butyric acid, phenol, o-chlorophenol and dimethyl phthalate into a soil. Injection depth was below the water table for this site, at 5.5 to 6.25 meters below the ground surface. For these organic substances, biotransformation appeared to be more important than sorption for a period of 204 days.

In summary, a model is proposed to depict the fate of organic waste substances disposed to a deep subsurface environment by injection wells. The model assumes that the principal reactions are sorption onto organic carbon, accumulation at the soil/water interface according to a Henry's law isotherm, biotransformation by native bacteria, and hydrolysis. A tabulation of values for use in predicting the retention of organic wastes is given in the Appendix. In the absence of experimental data, correlations based on laboratory measurements are suggested with appropriate modifications to reflect a soil/water environment down to 2500 m. Sorption of organic nonpolar compounds at 25°C is estimated from the relation for the equilibrium partitioning between soil organic matter and an aqueous phase by eq 4 (21), assuming that organic substances sorb independently of one another. In the absence of organic matter, accumulation of organic wastes at the mineral/water interface is reflected by Henry's constant. When laboratory measurements, or other data, are lacking the half-life of a trace organic can be estimated from Table 3-4 (46), for biotransformations. Bouwer and McCarty based their model, in part, on "large-scale subsurface contamination with low active organic concentrations and

slow ground-water movement...." (46). Under these conditions, biotransformation rates are expected to be slow by comparison with those at the surface. The effects of temperature in the deep subsurface environment on microbial action differ, depending on the kind of bacteria in this region (43,p.24). Hydrolysis half-lives from the evaluation by Mabey and Mill, are probably higher at, e.g., 55 °C, a temperature based on a geothermal gradient of 20 °C/km (45), and an average disposal depth of 5000 feet (1.5 km).

5.0 RECOMMENDATIONS FOR ADDITIONAL RESEARCH

After reviewing the current data on fate of organic wastes in the deep subsurface environment in consideration of the information required, some specific recommendations are appropriate. The optimum future laboratory measurements and predictive activities should include:

- 1. Development of a predictive model for the fate of organic wastes in the deep subsurface environment which takes into account the important reactions of specific organic substances. The computerized model should account for the following major chemical and physicochemical reactions: (a) rate of initial biotransformation of organic wastes, organized according to functional groups; (b) rate of bioconversion of organic wastes into carbon dioxide and water; (c) rate of waste sorption and equilbrium partition coefficient for the organic carbon portion of subsurface soils, for each functional group; (d) rate of adsorption and desorption of the organic materials onto minerals; (e) rate of hydrolysis by functional group in both pore water and in the liquid injected. In this way, the hydrolysis can be sorted into reactions for both "new" and "aged" disposal wells; (f) effects of temperature on the biokinetics, sorption and hydrolysis reactions, at depth.
- 2. Experiments which measure biodegradation, sorption and hydrolysis under laboratory conditions which closely simulate generic deep subsurface environments. Selected experiments with typical organic wastes would be used to improve the theoretical model, and might uncover additional important reactions of organic compounds in an abiotic environment.
- 3. Development of instrumentation to identify subsurface bacteria, and to monitor their

growth as a function of waste organic nutrients in the laboratory. An excellent approach appears to be a laser induced fluorescence instrument coupled with selected materials which give a "fingerprint" typical of specific bacteria (44).

- 4. Samples of soil at the depths of disposal sites. The soil would be analyzed for its organic carbon content. Knowledge of the OC content would help in development of the predictive model.
- 5. Samples of bacteria from the disposal well. Biotransformation may be the ultimate fate of the organic wastes; therefore, it is imperative that the native microorganisms be identified and their rates of biodegradation measured. In addition, bacteria introduced by injection of the wastes should be identified by sampling the disposal site.
- 6. Assessment of any advantage of organic wastes disposal in very deep wells, e.g., below 10,000 feet (3000 m). Computer modeling with an appropriate code would be used. The temperatures (e.g., 85 °C) and pressures in very deep wells may be sufficient to thermally decompose the hazardous organic substances into harmless degradation products. One related study is the recent publication by Takach et al. (45) on the stability of natural gas to 29,000 feet, which utilizes thermodynamic calculation of equilibrium compositions.

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Description of waste	Depth of well (ft)	Description of formation	Ref.
Aldehydes			
Chlorinated aldehydes (1.9%); acetic acid (1.3%); acetaldehyde (0.3%); pH = 5	5,400	Unconsolidated sands ranging in thickness from 100 to 200 ft, separated by thin shale beds $\phi = 30-40\%$, k=0.5-2.3 darcys	2
Formaldehyde (0.4%); acetaldehyde (0.4%); butyonaldehyde (0.3%); crotonaldehyde (0.2%); methyl alcohol (0.2%); butyl alcohol (0.14%); acrolein (0.02%); formic acid (0.01%);	4,500	Unconsolidated sand $\phi = 30\text{-}40\%$, k=1.0-3.0 darcys	2
acetic acid (0.01%) Acetaldehyde (750 ppm); propanol (100 ppm); acetone (350 ppm); acrolein (100 ppm); butanol (150 ppm); methyl ethyl ketone (300 ppm); vinyl methanol ketone (600 ppm); cyclobutanone (600 ppm); dioxane (100 ppm); NaCl (5,000 ppm); NaHCO ₈ (1,000 ppm); pH = 7.3;	4,000	Sandstone and detrital granite $\phi=25\%$, k = 500 - 1,000 md	2
Alcohols			
Hexanol (0.10%); amyl alcohol (0.06%); methanol (0.02%); dibasic acids (0.82%); hexamethylene imine (0.14); cyclohexane (0.02%); dodecane (0.01%); NaNO ₈ (1.3%); NaHCO ₈ (0.35%); unidentified organics (2.40%); pH = 9.1	3,300	Loosely consolidated sand - 260 ft thick, $\phi = 30-35\%$, $k = 1.0$ - 2.0 darcys	2

Description of waste	Depth of well (ft)	Description of formation	Ref.
Polyvinyl alcohol (300 ppm); methanol (1,400 ppm); acetic acid (400 ppm); hydroquinone (100 ppm); silicones (23 ppm); vinyl acetate, (1,200 ppm); methyl acetate (800 ppm); acetaldehyde (500 ppm); ethyl acetate (50 ppm); sodium acetate (2,300 ppm); pH = 5	7,000	Sandstone interbedded with shale.	2
Glycols			
Ethylene glycol (1.0%);			
K ₂ CO ₈ (1.0%);	5,000	Unconsolidated	2
Ethylene glycol (500 ppm); propylene dichloride (125 ppm); propylene glycol (11,000 ppm); acetone (40 ppm); ethylene dichloride (1,200 ppm); dichloroisoproyl ether (90 ppm); dichloroethyl ether (40 ppm); NaCl (5%); Na ₂ SO ₄ (0.05%); Na ₂ CO ₈ (0.01%);	6,000	Unconsolidated sand	2
Phenols Phenols (6.2%); cyanides (1.2%); ammonia (0.3%); thiocyanides (0.1%); NaCl (0.3%);	1,800	Sandsone - $\phi = 15\%$, $k = 200 \text{ md}$	2

Description of waste	Depth of well (ft)	Description of formation	Ref.
Acids			
Acetic acid (0.5%);	3,500	Unconsolidated sand -	2
chloroacetic acid (0.4%);	•	$\phi = 27 - 35\%$	
acetaldehyde (0.1%);		k= 0.5 - 2 darcys	
acetaldol (0.1%);			
Butanol (0.05%);			
crotonaldehyde (0.05%);			
NaCl (0.5%);			
pH = 2.5;		• .	
COD = 29,400 ppm;			
BOD = 9,900 ppm			
Adipic acid (8%);	6 000	Sandstone - $\phi = 28\%$,	•
caprolactum (1%);	6,000	Sandstone - $\psi = 28\%$, k = 0.2-1.0 darcy	2
butanol (0.5%);		k = 0.2-1.0 darcy	
benzene (0.2%);			
cyclohexane (0.3%);			
NaCl (2%);			
1401 (270),			
Organic acids	5,200	Loosely consolidated sand	2
(dibasic - principally adipic) (2.6%)			
nitric acid (0.8%)			
Acetic acid (3,000 ppm);	6,000	Sandstone - $\phi=30\%$	2
paratoluic acid (100 ppm);	,	k = 500 md	_
benzoic acid (50 ppm);			
xylene (100 ppm);			
methyl ethyl ketone (50 ppm);			
CaCl ₂ (200 ppm);			
MgSO ₄ (200 ppm);			
pH = 3.0			
C ₂ -C ₆ organic acids (6,700 ppm);	7,000	Sandstone	2
C ₁ -C ₄ alcohols (2,000 ppm);	7,000	$\phi = 27\%, k = 300 \text{ md}$	2
C_1 - C_4 alcohols (2,000 ppm); C_1 - C_8 ketones (1,400 ppm);		ψ 21/0, K = 300 Hd	
propylene glycols (400 ppm);		•	
propylene glycols (400 ppm); propyl esters (2,000 ppm);			
propylene oxide (3,000 ppm);			
		·	
C_6 - C_6 hydrocarbons (2,000 ppm); Na ₂ SO ₄ (2,000 ppm);			•
•			
NaCl (500 ppm);		•	

Description of waste	Depth of well (ft)	Description of formation	Ref.
Acetic acid (160 ppm);	7,200	Sandstone containing beds	2
maleic acid (100 ppm);		of clay $\phi = 26\%$,	
<pre>xylene sulfonic acid (50 ppm);</pre>		k=200 md	•
benzoic acid (50 ppm);		en e	*
butyl alcohol (950);			
styrene (100 ppm);			
isoproponal (100 ppm);			
phthalic acid (300 ppm);			
methyl ethyl ketone (250 ppm);			
NaOH (650 ppm);			
H_2SO_4 (500 ppm);			
Na ₂ CO ₈ (350 ppm);			
Nitriles			
Glycol nitrile (0.5%);	6,500	Sandstone - 400 ft thick,	2
iminodiacetonitrile (0.3%);		$\phi = 28\%$, k = 200 md	
nitrilotriacetonitrile (0.2%);	•		
other organic nitriles (1.0%);			
methanol (0.1%);		•	
methyl formate (0.1%);			
formaldehyde (0.3%);			
formic acid(0.5%);		•	
H ₂ SO ₄ (1.7%);			
(NH ₄) ₂ SO ₄ (4.3%);			
Disulfides			
	•		
Ethyl disulfide (30%);	2,200	Unconsolidated sand with	2
propyl disulfide (40%),	•	thin clay beds 100 ft thick	
butyl disulfide (25%);			
mercaptans (5%);		•	
water insoluble oil			
Ethyl disulfide (4%);	1,500	Limestone - vugular	2
propyl disulfide (6%);			
butyl disulfide (14%);			
pentyl disulfide (14%);			
hexyl disulfide (3%);			
NaOH (8%);			

Table 1-2 Selected geologic formations used for organic waste disposal (2).

Type of Formation	Formation	Location	Depth (ft)
Unconsolidated sand	Nacatoch	Louisiana	1,000
	Cockfield		3,000
	Catahoula	Texas	5,000
	Glorieta		1,300
	Frio	· · ·	7,500
Sandstone	Eutaw	Florida	3,500
·	Mount Simon	Illinois	4,000
	Tar Springs	Indiana	2,300
	Bethel		2,800
	Eau Claire		4,000
·	Mount Simon		5,500
	Sylvania	Michigan	1,000
	Yeso	New Mexico	1,000
	Burgoon	Pennsylvania	1,000
	Oriskany		5,500
	"Granite Wash"	Texas	5,000
	Greta		4,500
Vugular carbonate	Lake City	Florida	1,800
_	Cedar Valley	Illinois	2,500
	Arbuckle	Kansas	4,000
	Virginian		3,200
	St. Peter	Kentucky	1,000
	Dundee	Michigan	4,000
	Arbuckle	Oklahoma	2,000
	Bass Islands	Pennsylvania	1,600
	San Andres	Texas	5,000

Table 1-3 Properties of geologic material from three types of underground injection formations (2).

Unconsolidated Sand			Sandstone			Vugular Limestor		
Depth (ft)	Φ (%)	k (md)	Depth (ft)	Φ (%)	k (md)	Depth (ft)	Φ (%)	k (md)
4,442	35.4	452	2,466	13.7	505	5,021	17.5	12.3
4,458	37.2	3,185	2,467	14.0	417	5,021	15.2	5.3
4,512	35.9	958	2,468	15.1	482	5,022	21.0	5.8
4,527	32.8	4,200	2,469	16.6	289	5,022	22.0	4.6
4,540	30.0	943	2,470	16.7	3 30	5,022	16.8	58.9
4,625	31.1	2,980	2,471	11.4	197	5,022	19.0	83.8
4,636	32.5	4,750	2,472	13.2	165	5,023	12.9	4.2
4,674	31.3	3,930	2,473	14.6	413	5,023	14.5	1.7
4,682	29.8	2,560	2,474	17.2	229	5,023	19.5	5.5
4,696	29.6	2,020	2,475	18.1	147	5,023	19.5	4.6
4,706	35.3	1,880	2,476	16.6	962	5,033	16.5	1.1
4,716	31.2	5,800	2,477	17.5	885	5,033	16.0	7.5
4,877	25.2	12	2,478	12.1	124	5,034	12.9	0.7
4,898	31.1	5,120	2,479	7.1	3	5,034	15.5	4.6
4,911	32.6	4,210	2,480	15.8	665	5,246	18.4	0.7
4,922	28.6	2,500	2,481	14.8	367	5,246	17.9	0.6
4,940	34.5	988	2,482	12.9	780	5,247	17.3	0.5
4,951	33.9	1,830	2,483	15.2	845	5,247	16.7	0.6
4,962	33.1	2,270	2,484	16.0	862	5,247	18.6	0.6
4,970	34.0	1,626	2,485	14.7	450	5,247	18.2	0.5

Table 1-4a Adsorption of neutral and basic organonitrogen compounds by clay/oil shale mixtures (5:1 weight ratios) (23).

Clay	% Total N adsorbed	, , , , , , , , , , , , , , , , , , , ,		Ion exchange capacity meq/100g
Montmorillonite (Mont)	50	92	29	99
K-Mont	40	84	27	
Na-Mont	55	92	26	
Ca-Mont	55	97	34	
Mg-Mont	70	>99	30	
Cu-Mont	7 0	>99	38	
Halloysite	40	83	18	26
Kaolinite '	3 0	65	11	13
Muscovite	10	20	2	2

Table 1-4b Adsorption of selected neutral and basic nitrogen compounds by ion-exchanged montmorillonite (23).

Organic		Exc	hangeabl	e cation	
Nitrogen Compound	K ⁺	Na ⁺	Ca ⁺⁺	Mg ⁺⁺	Cu ⁺⁺
n-Butyronitrile	25	40	30	30	15
Pyridine	16	8	2	<1	<1
2,6-Dimethylpyridine	24	10	7	1	<1
2,5-Dimethylpyrrole	3	3	. <1	<1	<1
Quinoline	24	21	9	5	<1
Indole	23	26	12	12	<1
Stearonitrile	40	60	5 5	45	3 0

Table 1-5 Examples of microorganisms present in subsoils near the surface.

Material or Microorganism	Organic Chemical Reacted	Ref.
Arthrobacter		5
Pseudomonas		5
Enterobacter		5
Methanogenic Aquifer	•	6
Material	1,1-Dichloroethylene	6
	trans-1,2-Dichloroetheylene	6
	cis-1,2-Dichloroethylene	6
	Trichloroethylene	6
	1,2-Dibromoethane	6
	Benzene	6
	Toluene	6
	Ethylbenzene	6
	o-Xylene	6
Organic Sediment from	·	7
the Everglades	1,1-Dichloroethene	7
	cis-1,2-Dichloroethene	7
	trans-1,2-Dichloroethene	7
Estuarine Water	Phenol	8
	p-Chlorophenol	8
	2,4-Dichlorophenol	8
	2,4,5-Trichlorophenol	8
	Pentachlorophenol	8 .
Alluvial Groundwater	Alkylpyridines	9
Sea Water	Naphthalene; Toluene	10
Soil Samples,	-	11
30 to 457 cm	Trichloroethylene; 1,2-Dichloroethylene	11
Interstitial Waters		12
of a Sediment		12
0-41 cm. depth	Polychlorinated Biphenyls (PCBs)	12

Table 3-1 Sediments and soils--physical properties (16)

Location	% O.C.	pН	CEC me. 100 g	Silt	Clay	Swelling Clay
Missouri River, Stanton, N. Dak.	2.07	7.79	23.7	41.8	55.2	40.1
Missouri River, Lake Oahe, N. Dak.	2.28	7.44	19.0	35.4	31.0	25.7
Missouri River, Big Bend Lake, S. Dak.	0.72	7.83	33.0	31.2	68.6	60.8
Missouri River, Onawa, Iowa	0.15	8.32	3.7	10.7	6.8	6.1
Loess sample, Turin, Iowa	0.11	8.34	12.4	75.6	17.4	16.3
Ohio River, Ceredo, W. Va.	3.04	6.90	11.86	27.1	52.6	-
Soil eroded hillside, Ceredo, W. Va	0.48	4.54	18.9	34.4	63.6	13.8
Ohio River, Leavenworth, Ind.	0.95	7.79	11.3	48.7	35.7	10.1
Mississippi River, Columbus, Ky.	0.66	7.76	15.4	25.8	39.5	29.7
Soil, Fern Clyffe State Park, Ill.	1.30	5.5	8.5	71.4	28.6	21.9
Illinois River, Lorenzo, Ill.	1.88	7.60	8.3	42.7	7.1	2.8
Illinois River, Lacon, Ill.	1.67	7.55	8.5	52.7	21.2	14.8
Confluence of Illi- nois and Sanga- mon Rivers	2.38	6.70	31.2	13.6	69.1	57.6
Mississippi River, McClure, Ill.	1.48	7.75	20.9	55.4	42.9	37.1
Small stream, Watkinsville, Ga.	1.21	6.35	3.7	13.9	18.6	2.0

Table 3-2a Pyrene and methoxychlor sorption coefficients for linear portion of isotherms (15).

Sediment size	Organic	Pyrene			chlor		
fraction	carbon	K_p	(r ²)	K_{oc} (×10 ^{-δ})	K_p	(r ²)	Koc (×10-5)
Hickory Hill							
Sand	0.13	42	(0.85)	0.32	53	(0.85)	0.41
Coarse silt	3.27	3000	(0.95)	0.92	2600	(0.97)	0.80
Medium silt	1.98	2500	(0.95)	1.3	1800	(0.91)	0.91
Fine silt	1.34	1500	(0.90)	1.1	1400	(0.97)	1.0
Clay	1.20	1400	(0.75)	1.2	1100	(0.99)	0.92
$Doe\ Run$			` ,			,	
Sand	0.086	9.4	(0.97)	0.11	8.3	(0.98)	0.097
Coarse silt	2.78	2100	(0.92)	0.76	2200	(0.93)	0.80
Medium silt	2.34	3000	(0.80)	1.3	1700	(0.94)	0.73
Fine silt	2.89	3600	(0.88)	1.2	2300	(0.95)	0.80
Clay	3.29	3800	(0.90)	1.2	2400	(0.98)	0.73
Oconee River			, ,			, ,	•
Sand	0.57	68	(0.38)	0.12	95	(0.93)	0.17
Coarse silt	2.92	3200	(0.99)	1.1	2500	(0.97)	0.86
Medium silt	1.99	2300	(0.96)	1.2	2000	(0.93)	1.0
Fine silt	2.26	2500	(0.99)	1.1	2100	(0.96)	0.93
Clay*			, ,			` ,	

^{*} The clay portion of this sediment was allowed to age in suspension and degraded substantially. Therefore, no sorption was done on this fraction.

Soil OC normalized sorption coefficients (K_{OC}) (Ref. 20) Table 3-2b

	Sorption coefficient (K_{OC})								
Compound	. Apison*		Fullerton ^b			Dormont ^c			
Ketones									
Acetophenone 2-Acetonaphthone Benzophenone 4-Acetylbiphenyl 9-Acetylanthracene	185 1200 580 2400 8400	± ± ± ±	8 ^d 100 100 200 1200	270 410 530 710 1170	± ± ± ±	80 80 130 50 230	105 950 440 1900 1750	± ± ± ±	4 20 30 120 160
Alcohols									
Benzyl Alcohol Sec-Phenethylalcohol 1-Naphthalenemethanol 4-Biphenylmethanol 9-Anthracenemethanol	<5 37 165 700 8000	± ± ±	2 12 35 3300	<5 <5 <5 105 1500	± ±	2 50	<5 52 270 495 2700	± ± ±	4 14 70 480
Phenols	[
Phenol p-Cresol 3,5-Dimethylphenol 5-Indanol 2,3,5-Trimethylphenol 1-Naphthol	55 3420 460 11,500 5700 369,000	* * * * * *	25 970 90 1900 2750 128,000	710 3350 1400 10,000 6400 >600,000	± ± ± ± ±	75 580 400 2470 570 300,000	7 115 190 240 260 1280	± ± ± ± ±	7 19 40 40 4 340
Nonplar compounds									
Naphthalene Biphenyl Phenanthrene 1,4-Dichlorobenzene 1,2,4-Trichlorobenzene	1000 3300 5900 665 2100	± ± ± ±	80 260 420 105 360	960 900 1400 850 1300	± ± ± ±	50 60 200 270 610	400 870 5800 280 885	± ± ± ±	52 65 375 105 160

^a0.1% OC. ^b0.06% OC. ^c1.2% OC. ^dMean ± S.E., n = 3-6.

Table 3-3 Water solubilities (S), molar volumes (\overline{V}), octanol-water partition coefficients (K_{ow}), and soil organic matter-water distribution coefficients (K_{om}) of selected organic solutes (21).

Compound	log S, mol/L	V, L/mol	$\log{(\mathrm{S}\overline{\mathrm{V}})}$	log K _{om}	log K _{ow}
benzene	-1.64	0.0894	-2.69	1.26	2.13
anisole	-1.85	0.109	-2.82	1.30	2.11
Chlorobenzene	-2.36	0.102	-3.35	1.68	2.84
ethylbenzene	-2.84	0.123	-3.75	1.98	3.15
1,2-dichlorobenzene	-2.98	0.113	-3.9 8	2.27	3.38
1,3-dichlorobenzene	-3.04	0.114	-3.9 8	2.23	3.38
1,4-dichlorobenzene	(-3.03)	0.118	-3.96	2.20	3.39
1,2,4-trichlorobenzene	-3.57	0.125	-4.47	2.70	4.02
2-PCB	(-4.57)	0.174	-5.33	3.23	4.51
2,2'-PCB	(-5.08)	0.189	-5.57	3.68	4.80
2,4'-PCB	(-5.28)	0.189	-5.97	3.89 /	5.10
2,4,4'-PCB	(-5.98)	0.204	-6.67	4.38	5.62

^{*}The listed solubilities are the 20-25 °C values. The numbers in parentheses are the supercooled liquid solute solubilities. For 2-, 2,2', 2,4'-, and 2,4,4'-PCB, the calculations were based on their solid solubilities of 3760, 717, 637, and 115 μ g/L at 20 °C, respectively and an assumption of 13.5 cal/(mol K) or their entropies of fusion. The molar volumes of PCBs are estimated by using the densities of liquid Aroclor mixtures that have approximately the same chlorine atoms as the individual PCBs.

Table 3-4 Half-lives of biotransformation modeled with first-order batch kinetics as a function of active organism concentration (46).

	anism ntration	Degradatio	on half-life, days
mg/L	No./ml*	Aerobic respiration [†]	Methanogenesis [§]
10.	10 ⁷	0.014	0.14
1.	10^{6}	0.14	1.39
0.1	10^{5}	1.39	13.9
0.01	10^{4}	13.9	139.
0.001	10^{3}	139.	1,390.

Organism dry weight taken as 10^{-12} g/cell † k/K_s = 5 l/mg cells-day § k/K_s = 0.5 l/mg cells-day

Table 3-5 Hydrolysis half-life of selected organic substances.

Data are from Mabey and Mill (33), and (34, 37) for pH=7,

25 °C, and zero ionic strength.

organic substance	t _{1/2}
CH ₂ Cl ₂	704 yr
CH ₂ Br Cl	44 yr
CH ₂ Br ₂	183 yr
CH Cl ₃	3500 yr
CH Br Cl ₂	137 yr
CH Br ₂ Cl	274 yr
CH Br ₃	686 yr
CHI Cl ₂	275 yr
CHFICI	1.0 yr
CCl ₄	7000 yr (1 ppm),
	7 yr (1000 ppm)
CH ₃ F	30 yr
CH ₃ Cl	0.93 yr
CH ₃ Br	2 0 day
CH ₃ I	110 day
$C_2 H_5 Cl$	38 day
$C_2 H_{\delta} Br$	3 0 day
$C_2 H_6 I$	49 day
$(CH_3)_2$ CH Cl	38 day
(CH ₃) ₂ CH Br	2.1 day
$(CH_3)_2$ CHI	2.9 day
CH ₃ CH ₂ CH ₂ Br	2 6 day
$(CH_3)_3$ CF	50 day
$(CH_3)_3$ CCl	23 sec
CH ₂ =CH CH ₂ Cl	69 day
CH ₂ =CH CH ₂ Br	12 hr
CH ₂ =CH CH ₂ I	2.0 day
C ₆ H ₅ CH ₂ Cl	15 hr
p-CH ₃ C ₆ H ₄ CH ₂ Cl	0.43 hr
C ₆ H ₅ CH Cl ₂	0.1 hr
C ₆ H ₅ C Cl ₃	19 sec
C ₆ H ₆ CH ₂ Br	1.32 hr
p-CH ₃ C ₆ H ₄ CH ₂ Br	4.3 min
CH ₃ COOC ₂ H ₅	2.0 yr
CH ₃ CONH ₂ (acetamide)	3950 yr

7.0 APPENDIX

The organic substances tabulated in this Appendix are included because they are listed in the Groupings of wastes, in Reference 47. Other organic compounds are included if they are likely to be found in a subsurface environment, or if they provide useful data such as sorption measurements. The intent of this tabulation is to provide values of half-lives for hydrolysis and biotransformation reactions, and equilibrium sorption values at the soil/water interface. Data are largely at 25 °C and atmospheric pressures. These data then can provide the basis for predicting the rates and equilibrium values in the deep subsurface, after modification of the appropriate mathematical equations. There are numerous gaps in the tables, representing a lack of experimental values. An especially large gap is represented by rates of biological transformations. Our calculations have been limited to sorption coefficients, estimated from octanol/water partition coefficients using eq 4.

The organic wastes are organized into the following major groupings:

- 7.1 Halogenated Solvents
- 7.2 Phthalates
- 7.3 Halogenated Pesticides
- 7.4 Polynuclear Aromatics
- 7.5 Oxygenated
- 7.6 Polymerizables
- 7.7 Phenolics
- 7.8 Nonhalogenated Solvents
- 7.9 Reactive (Non-Cyanide)
- 7.10 Nonhalogenated Pesticides
- 7.11 Organo-Nitrogens
- 7.12 Organo-Sulfurs
- 7.13 Halogenated N.O.S.
- 7.14 Pharmaceuticals
- 7.15 Dyes

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CONVERSION TABLE

- 1 Pa (Pascal) = 10^{-5} bar = 0.00750062 torr
- 1 atm = 760 torr
- 1 torr = 1 mm Hg
- 1 kPa = 1000 Pa
- $1~\mathrm{mg/L} = 1~\mathrm{g/m}^3$
- $1 \text{ kPa m}^3/\text{mol} = .0009869 \text{ atmos. m}^3/\text{mol}$
- 1 kPa = 7.50062 torr
- 1 m = 3.279 ft

ATOMIC WEIGHTS OF SELECTED ELEMENTS

Oxygen: 15.9994

Hydrogen: 1.0080

Nitrogen: 14.0067

Sulfur: 32.064

Carbon: 12.0112

Phosphorus: 30.9738

FORMAT FOR PROPERTY VALUES OF WASTE ORGANICS

For each pure organic substance in the Appendix, the Grouping category is given, followed by the name of the organic and the date identifying when the data were tabulated. Other commonly used names for the organics are given, together with the Chemical Abstracts Services (CAS) number. For purposes of identification, we also include the Hazardous Waste Number, and an Industry/EPA Generic No. The Industry/EPA Generic No. is not, however, related to the source of the waste injected into the subsurface disposal system. The empirical chemical formula is also given. Under the heading "Description", there is a listing of selected physical properties of the pure substance, and a reference to source of the data. The physical properties are intended to provide information on the organic chemical to be used in predicting its form in the deep subsurface, e.g., gaseous, soluble in water, or solid. The general heading Retention properties is defined as the chemical and biochemical reactions which determine the fate of the organic within the injection zone. These values are all at 25°C and about 1 bar pressure.

The data in the tabulation of the Appendix are the database for laboratory conditions which are to be used in predicting the fate of organic wastes disposed to the deep subsurface by injection wells.

Other Names: CAS No.: Hazardous Waste No.: Industry/EPA Generic No.: Formula: C7H7CIN2S Description: Physical properties: Ref. Molecular weight: Melting point (C): Boiling point (C): Density (g/cc, 20 C): Vapor pressure (torr, 25 C): Henry's law constant (atm m**3 mole**-1, 25 C): Solubility in water (mg/L, 25 C): Retention properties: Ref. log (octanol/water) partition (25 C): Partition coefficient, soil/water: Hydrolysis rate in water (t1/2, month**-1, 25 C): Biodegradation rate in water (t1/2, month**-1):

7.1 Halogenated Solvents

HALOGENATED SOLVENTS bis(Chloromethyl) ether

```
Other Names: sym-Dichloromethyl ether; Oxybis[chloromethane]; BCME;
             Chloro(chloromethoxy) methane; Chloromethyl ether;
             Dichlorodimethyl ether; alpha-alhpa'-Dichlorodimethyl ether;
             Monochloromethyl ether
Haz Waste #: P016
                          Ind/EPA Gen #:
             (CH2-C1)20: C2H4C120
Formula:
Description: Colorless liquid; suffocating odor; strong irritant to eyes,
             respiratory tract; listed as a carcinogen by the EPA
PHYSICAL PROPÈRTIES: - - - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                               114.97 2
Boiling point (C):
                                                               106.00
                                                                      2
Density (g/cc, 20C):
                                                               1.3150 2
Solubility in water (mg/L, 25C):
                      Decomposed by water into HCI and formaldehyde 2
RETENTION PROPERTIES:
log (octanol/water) partition (25C):
                                                               -0.38 3
Partition coefficient, soil/water:
                                                               0.0750 17
Hydrolysis rate, water (1/month, 25C):
                                     Very rapid, to form HCl and HCHO 3
Biodegradation rate in water (1/month):
                                                        Not important 3
                                  HALOGENATED SOLVENTS
August 1987
CAS #:
         108-90-7
                                  Chlorobenzene
Other Names: Monochlorobenzene; Benzene Chloride
Haz Waste #: U037
                    Ind/EPA Gen #: F002,K015,K105,K085
Formula:
             C6H5C1
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                               112.56 1
Melting point (C):
                                                               -45.34 1
Boiling point (C):
                                                               131.70 1
Density (g/cc, 20C):
                                                               1.1012 1
Vapor pressure (torr, 250):
                                                              12.0000 5
Henry's law const (atm m**3/mol, 25C):
                                                          0.39300E-02 3
Solubility in water (mg/L, 25C):
RETENTION PROPERTIES: - - - - - - -
                                                               488.00 3
                                                              - - - Ref.
log (octanol/water) partition (25C):
                                                                 2.84 3
Partition coefficient, soil/water:
                                                            61.4000 17
```

Biodegradation rate in water (1/month): Bioaccumulated and biodegraded 3

HALOGENATED SOLVENTS Chloroform

Other Names: Trichloromethane; Formyl trichloride; Freon 20; R 20; Methane trichloride; Methenyl trichloride; Methyl

trichloride; NCI-CO2686; Trichloroform

Haz Waste #: UO44 Ind/EPA Gen #:

Formula: CHC13

Description: Highly refractive, non-flammable, heavy, very volatile,

sweet tasting liquid (Ref. 1); banned by the FDA from use in drug, cosmetic and food packaging products in 1976; listed

as a carcinogen by the EPA (Ref. 2)

PHYSICAL PROPERTIES:		Ref.
Molecular weight:	119.39	2
Melting point (C):	-63.50	2
Boiling point (C):	61.70	3
Density (g/cc, 20C): 1.484; 1.48069 specific gravity,	25 C (Ref. 2)	1
Vapor pressure (torr, 25C):	150.00	3
Henry's law const (atm m**3/mol, 25C):	0.33900E-02	3
Solubility in water (mg/L, 25C):	8200.00	3
RETENTION PROPERTIES:		Ref.
log (octanol/water) partition (25C):	1.97	3
Hydrolysis rate, water (1/month, 25C):	450.00	4
Biodegradation rate in water (1/month):	Significant	3

August 1987

CAS #: 95-50-1

HALOGENATED SOLVENTS 1,2-Dichlorobenzene

Other Names: o-Dichlorobenzene; Orthodichlorobenzene

Haz Waste #: U070 Ind/EPA Gen #: F002,K042,K085,K105

Formula: 1,2-C6H4Cl2

Description: Liquid; can cause injury to liver, kidneys; high

concentrations cause CNS depression (Ref. 2)

concentrations cause one depression (i.e., 2)		
PHYSICAL PROPERTIES:		Ref.
Molecular weight:	147.00	1
Melting point (C):	-16.97	1
Boiling point (C):	180.40	1
Density (g/cc, 20C):	1.3022	1
Vapor pressure (torr, 25C):	1.5000	3
Henry's law const (atm m**3/mol, 25C):	0.19400E-02	3
Solubility in water (mg/L, 25C):	145.00	3
RETENTION PROPERTIES:		Ref.
log (octanol/water) partition (25C):	3.38	3
Partition coefficient, soil/water:	189.00	17
Biodegradation rate in water (1/month):	Bioaccumulated	3

August 1987

HALOGENATED SOLVENTS 1,3-Dichlorobenzene

CAS #: 541-73-1

Other Names: m-Dichlorobenzene

Ind/EPA Gen #: K085,K105 Haz Waste #: U071

Formula: 1,3-C6H4Cl2

Formula: 1,3-C6H4C12 PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - - - Ref. Molecular weight: 147.00 1 Melting point (C): -24.76 1 Boiling point (C): 173.00 1 Density (g/cc, 20C): 1.2828 1 Vapor pressure (torr, 25C): 2.2800 3 Henry's law const (atm m**3/mol, 25C): 0.26300E-02 3 Solubility in water (mg/L, 25C): 123.00 3
RETENTION PROPERTIES: ---- Ref. log (octanol/water) partition (25C):

Partition coefficient, soil/water:

Biodegradation rate in water (1/month):

3.38 3
189.00 17
Bioaccumulates 3 3.38 3

HALOGENATED SOLVENTS

August 1987 CAS #: 764-41-0

1,4-Dichloro-2-butene

Haz Waste #: U074 Ind/EPA Gen #: Formula: CICH2CH=CHCH2CI; C4H6CI2

Description: Highly toxic by inhalation; suspected carcinogen (Ref. 13)

Molecular weight:

125.00 13

August 1987

CAS #: 75-71-8

HALOGENATED SOLVENTS

Dichlorodifluoromethane

Other Names: Freon-12

Haz Waste #: U075

Ind/EPA Gen #:

Formula: CC12F2

Description: Almost odorless gas; can react violent with Al or Mg

(Ref. 13)

Molecular weight: 121.00 13 Melting point (C): -155.00 13

Boiling point (C):

-29.80 13

HALOGENATED SOLVENTS 1,1-Dichloroethane

·	
Other Names: Ethylidene dichloride; Ethylidene	e chloride
Haz Waste #: U076 Ind/EPA Gen #:	
Formula: C12CHCH3; C2H4C12 Description: Dily liquid; odor and taste as of	f chloroform: parcotic in
high concentrations (Ref. 2)	i cirrororiii, narcoore iii
PHYSICAL PROPERTIES:	Ref.
Molecular weight:	98.96 3
Melting point (C):	-97.00 3 57.30 3
Boiling point (C): Density (g/cc, 20C):	(25 C) 1.1680 2
Vapor pressure (torr, 25C):	180.00 3
Henry's law const (atm m**3/mol, 25C):	0.54500E-02 3
Solubility in water (mg/L, 25C):	(20 C) 5500.00 3
RETENTION PROPERTIES:	Ref. 1.79 3
Partition coefficient, soil/water:	6.9000 17
Biodegradation rate in water (1/month):	Degradation significant 3
August 1987 HALOGENATE	O SOLVENTS
CAŠ #: 107-06-2 1,2-Dichlor	
Other Names: Ethylene dichloride; Borer sol; Independent of the Destrux of Borer-sol; sym-Dichlorous Dichlorous on; alpha, beta-Dichlorous Dichlorous of the Dichlorous of the Dichloride; NIC-C00511; Educhloride of the Dutch liquid of the Dutch li	oethane; Di-chlor-mulsion; loroethane; EDC; ENT 1656; ide; Freon 150; thylene chloride;
	•
Description: Heavy liquid; pleasant odor; sweet	
Description: Heavy liquid; pleasant odor; sweet irritating to the respiratory tra	et taste; vapors are act and conjunctiva; listed
irritating to the respiratory tra as a carcinogen by the EPA (Ref.	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transcription as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2) Ref.
irritating to the respiratory transactions as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Boiling point (C): Density (g/cc, 20C): Vapor pressure (torr, 25C): Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): RETENTION PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)
irritating to the respiratory transaction as a carcinogen by the EPA (Ref. PHYSICAL PROPERTIES:	et taste; vapors are act and conjunctiva; listed 2)

August 1987
CAS #: 75-35-4

HALOGENATED SOLVENTS
1,1-Dichloroethylene

Other Names: 1,1-Dichloroethene; Vinylidene chloride; asym-Dichloroethylene

Haz Waste #: U078 Ind/EPA Gen #: Formula: CH2=CC12; C2H2C12

Description: Liquid; mild, sweet odor resembling that of chloroform;

irritant to skin, mucous membranes; narcotic in high concentrations; caused kidney and liver damage in

experimental animals (Ref. 2)

experimental animate (non 2)		
PHYSICAL PROPERTIES:		
Molecular weight:	96.94 5	
Melting point (C):	-122.10 5	
Boiling point (C):	. 37.00 5	
Vapor pressure (torr, 25C):	591.00 3	
Henry's law const (atm m**3/mol, 25C):	0.0150 3	
Solubility in water (mg/L, 25C):	400.00 3	
RETENTION PROPERTIES:	Ref.	
log (octanol/water) partition (25C):	1.48 3	
Partition coefficient, soil/water:	3.6200 17	
Biodegradation rate in water (1/month):	Significant 3	

August 1987 HALOGENATED SOLVENTS
CAS #: 156-59-2 cis-1,2-Dichloroethylene

Other Names: cis-1,2-Dichloroethene

Haz Waste #: U079 Ind/EPA Gen #:

Formula: cis-CHC1=CHC1

PHYSICAL PROPERTIES: - - - -- - - Ref. Molecular weight: 96.95 1 Melting point (C): -80.50 5 Boiling point (C): 60.20 1 Density (g/cc, 20C): (15 C) 1.2917 1 Vapor pressure (torr, 25C): 27.4600 5 Henry's law const (atm m**3/mol, 25C): 0.75100E-02 5 Solubility in water (mg/L, 25C): 3500.00 5

HALOGENATED SOLVENTS trans-1,2-Dichloroethylene

Other Names: 1,2-Dichloroethene; Acetylene Dichloride; alpha, beta-Dichloroethylene; sym-Dichloroethaz Waste #: U079 Ind/EPA Gen #: Formula: trans-CHCI=CHCI Description: Colorless, mobile liquid, sweet odor PHYSICAL PROPERTIES:	+ thylene 96.95 -49.44 47.70 (10 C) 1.2631 200.00 0.53200E-02 600.00	1 3 3 3 Ref. 3 17
August 1987 CAS #: 78-87-5 Other Names: Propylene dichloride Haz Waste #: U083 Formula: CH2CICHCICH3 HALOGENATED SOLVENTS 1,2-Dichloropropane		
PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Boiling point (C): Vapor pressure (torr, 25C): Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): RETENTION PROPERTIES:	113.00 -100.00 96.80 (20 C) 42.0000 0.28200E-02 (20 C) 2700.00 	3 3 3 3 Ref. 3

August 1987 CAS #: 542-75-6

Density (g/cc, 20C):

HALOGENATED SOLVENTS 1,3-Dichloropropene

Other Names: 1,3-Dichloropropylene Haz Waste #: U084 Ind/EPA Gen #: cis and trans-CICH2CH=CHCI Formula: - - - Ref. Molecular weight: 111.00 3 104. (cis); 112. (trans) 3 Boiling point (C): (20 C) 25.0000 3 Vapor pressure (torr, 25C): 0.35500E-02 3 Henry's law const (atm m**3/mol, 25C): HALOGENATED SOLVENTS August 1987 75-69-4 Fluorotrichloromethane CAS #: Other Names: Trichlorofluoromethane; Freon-II; Freon 11; Frigen 11; Trichloromonofluoromethane; Arcton 9 Haz Waste #: U121,U229 Ind/EPA Gen #: F002 Formula: CC13F Description: Liquid at temperatures below 23.7; faint ethereal odor; non-flammable; decomposes into harmful materials by flames or high heat; may be narcotic in high concentrations; use as an aerosol propellant regulated by government (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. 137.38 2 Molecular weight: Melting point (C): -111.00 3 Boiling point (C): 23.80 3 Density (g/cc, 20C): (25 C, gas, air=1) 5.0400 2 Vapor pressure (torr, 25C): (20 C) 667.00 3 Henry's law const (atm m**3/mol, 25C): 0.0583 3 Solubility in water (mg/L, 25C): 1100.00 3 RETENTION PROPERTIES: ---- Ref. log (octanol/water) partition (25C): Partition coefficient, soil/water: 2.53 3 32.2000 17 Biodegradation rate in water (1/month): Not significant 3 HALOGENATED SOLVENTS August 1987 CAS #: 630-20-6 1,1,1,2-Tetrachloroethane Ind/EPA Gen #: Haz Waste #: U208 Formula: C13CCH2C1; C2H2C14 Description: Liquid PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref. Molecular weight: 168.00 13 Boiling point (C): 135.10 13

(25 C) 1.5420 13

HALOGENATED SOLVENTS 1,1,2,2-Tetrachloroethane

Other Names: Acetylene tetrachloride; Tetrachloroethane; Cellon; sym-Tetrachloroethane; Bonoform Haz Waste #: U209 Ind/EPA Gen #: CHC12CHC12; C2H2C14 Formula: Description: Nonflammable, heavy, mobile liquid; sweetish, suffocating odor similar to chloroform; has the highest solvent power of the chlorinated hydrocarbons; powerful narcotic; liver poison (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - -- - - Ref. Molecular weight: 167.86 2 Melting point (C): -36.00 3 Boiling point (C): 146.00 3 Vapor pressure (torr, 25C): (20 C) 5.0000 3 Henry's law const (atm m**3/mol, 25C): 0.43200E-03 3 (20 C) 2900.00 3 Solubility in water (mg/L, 25C): RETENTION PROPERTIES: - - - - - -- - - - Ref. log (octanol/water) partition (25C): 2.56 3 Partition coefficient, soil/water: 34.3000 17 Biodegradation rate in water (1/month): Not significant 3 HALOGENATED SOLVENTS August 1987 CAS #: 127-18-4 Tetrachloroethylene Other Names: Perchloroethylene; Ethylene Tetrachloride; Nema; Tetrachloroethene; Tetracap; Tetropil; Perclene; Ankilostine; Didakene Haz Waste #: U210 Ind/EPA Gen #: F001,F002,K030 CC12=CC12 Formula: Description: Colorless, nonflammable liquid, ethereal odor - - - Ref. Molecular weight: 165.85 Melting point (C): -22.70 3 Boiling point (C): 121.20 1 Density (g/cc, 20C): 1.6230 2 Vapor pressure (torr, 25C): 14.0000 Henry's law const (atm m**3/mol, 25C): 0.0287 3 Solubility in water (mg/L, 25C): 150.00 to RETENTION PROPERTIES: ------200.00 3 --- Ref. log (octanol/water) partition (25C): 2.88 5 Partition coefficient, soil/water: 66.8000 17 Hydrolysis rate, water (1/month, 25C):

Biodegradation rate in water (1/month):

0.73 4

Significant 5

HALOGENATED SOLVENTS Carbon tetrachloride

Other Names: Tetrachloromethane; Perchloromethane; Necatorina; Benzinoform; Carbon chloride; Carbona; ENT 4705; Fasciolin; Flukoids; Freon 10; Halon 104; Methane tetrachloride; Necatorina: Tetrachlorocarbon; Tetrafinol; Tetraform; Tetrasol; Univerm; Vermoestricid Ind/EPA Gen #: F001,K016 Haz Waste #: U211 Formula: CC14 Description: Colorless, clear, nonflammable heavy liquid (Ref. 1); poisoning by inhalation, ingestion or skin absorption; can be fatal; listed as a carcinogen by the EPA (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 153.84 2 -23.00 2 Melting point (C): Boiling point (C): 76.70 2 Density (g/cc, 20C): 1.589 (25 C); specific gravity 1.59472 (20 C) (Ref. 1) Vapor pressure (torr, 25C): 90.0000 3 Henry's law const (atm m**3/mol, 25C): 0.30200E-02 3 Solubility in water (mg/L, 25C): 785.00 3 RETENTION PROPERTIES: - - - Ref. log (octanol/water) partition (25C): 2.64 3 Partition coefficient, soil/water: 40.5000 17 Hydrolysis rate, water (1/month, 25C): Hydrolysis not likely 3 Biodegradation rate in water (1/month): Significant 3 August 1987 HALOGENATED SOLVENTS CAS #: 75-25-2 Bromoform Other Names: Tribromomethane Ind/EPA Gen #: Haz Waste #: U225 Formula: CHBr3 Description: Heavy liquid; chloroform odor; sweetish taste; abuse may lead to habituation or addiction (Ref. 2) - - - Ref. Molecular weight: 252.80 3 Melting point (C): 8.30 3 Vapor pressure (torr, 25C): (34 C) 10.0000 3 Henry's law const (atm m**3/mol, 25C): 0.53200E-03 3 Solubility in water (mg/L, 25C): (15 C) 3010.00 3 - **- - - - - - - - - - - - - - - R**ef. RETENTION PROPERTIES: - - - - - -2.30 3 log (octanol/water) partition (25C): Partition coefficient, soil/water: 19.9000 17 Biodegradation rate in water (1/month): Not significant 3

HALOGENATED SOLVENTS 1,1,1-Trichloroethane

Other Names: Methylchloroform; Chlorothene Haz Waste #: U226	
Molecular weight: Melting point (C): Boiling point (C): Density (g/cc, 20C):	133.42 2 -30.40 3 74.00 1 1.3249 1
Vapor pressure (torr, 25C): Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): (RETENTION PROPERTIES:	(20 C) 96.0000 3 0.49200E-02 3 (20 C) 480.00 to 4400.00 3
Partition coefficient, soil/water: Hydrolysis rate, water (1/month, 25C): Biodegradation rate in water (1/month):	9.2000 17 180.00 4 Significant 3
August 1987 HALOGENATED CAS #: 79-00-5 1,1,2-Trichl	
Other Names: Vinyl trichloride Haz Waste #: U227	
PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Boiling point (C): Density (g/cc, 20C): Vapor pressure (torr, 25C):	
Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): RETENTION PROPERTIES: log (octanol/water) partition (25C): Partition coefficient, soil/water: Biodegradation rate in water (1/month):	0.84600E-03 3 (20 C) 4500.00 3

August 1987 CAS #: 79-01-6

Biodegradation rate in water (1/month):

HALOGENATED SOLVENTS Trichloroethylene

Significant 3

Other Names: 1,1,2-Trichloroethene; Ethylene Trichloride; Ethinyl Trichlroide; Tri-clene; Trielene; Trichloran; Trichloren; Algylen; Trimar; Triline; Tri; Triethylene; Westrosol; Chlorylen; Gemalgene; Germalgene Ind/EPA Gen #: F001,F002,K030 Haz Waste #: U228 CHC1=CC12 Formula: Description: Nonflammable liquid; odor similar to chloroform; dissolves most fixed and volatile oils (Ref. 1); slowly decomposed by light in the presence of moisture; moderate exposures can cause symptoms similar to alcohol inebriation: higher concentrations have a narcotic effect (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - -Molecular weight: 131.39 1 Melting point (C): -73.00 3 Boiling point (C): 86.70 2 Density (g/cc, 20C): 1.4649 1 Vapor pressure (torr, 25C): 57.9000 Henry's law const (atm m**3/mol, 25C): 0.0117 3 Solubility in water (mg/L, 25C): 1100.00 3 RETENTION PROPERTIES: - - - -- - - Ref. log (octanol/water) partition (25C): 2.29 3 Partition coefficient, soil/water: 19.5000 17 Hydrolysis rate, water (1/month, 25C): 0.89 4

7.2 Phthalates

August 1987 **PHTHALATES** CAS #: 117-81-7 bis-(2-ethylhexyl) phthalate Other Names: DEHP Ind/EPA Gen #: Haz Waste #: U028 C6H4 (C00CH2CH (C2H5) C4H9) 2 Formula: PHYSICAL PROPERTIES: - - - - - - -Ref. Molecular weight: 391.00 3 Melting point (C): -50.00 3 Boiling point (C): 387.00 3 Vapor pressure (torr, 25C): 0.20000E-06 3 Solubility in water (mg/L, 25C): 0.4000 3 RETENTION PROPERTIES: - - - - - -- - - Ref. log (octanol/water) partition (250): (approx.) 8.73 3 Partition coefficient, soil/water: 0.12000E+08 17 Biodegradation rate in water (1/month): Significant 3 **PHTHALATES** August 1987 CAS #: 84-74-2 di-n-Butyl phthalate Other Names: DBP; 1,2-Benzenedicarboxylic acid dibutyl ester; n-Butyl phthalate; Phthalic acid dibutyl ester Ind/EPA Gen #: Haz Waste #: U069 C6H4(C00C4H9)2; C16H22O4 Formula: Description: Oily liquid; insect repellent for the impregnation of clothing (Ref. 2) PHYSICAL PROPERTIES: - - - - - - -- - - Ref. Molecular weight: 278.34 2

log (octanol/water) partition (25C): (approx.) 5.20 3

Melting point (C):

Boiling point (C):

Vapor pressure (torr, 25C):

Solubility in water (mg/L, 25C):

RETENTION PROPERTIES: - - - - -

Biodegradation rate in water (1/month):

-35.00 3

340.00 3

0.1000 3

13.0000 3

- - - Ref.

Rapid 3

PHTHALATES Diethyl phthalate

Other Names: DEP; 1,2-Benzenedicarboxylic aci Phthalic acid ethyl ester; Ethyl Haz Waste #: U088 Ind/EPA Gen #: Formula: C6H4(C00C2H5)2 Description: Colorless, practically odorless, disagreeable taste; irritating t in high concentrations, narcotic	phthalate; Palatinol A oily liquid; bitter, o mucous membranes, and,
PHYSICAL PROPERTIES:	-` Ref.
Molecular weight:	222.20 3
Melting point (C):	-40.50 3
Boiling point (C):	298.00 3
Density (g/cc, 20C):	(14 C) 1.2320 2
Vapor pressure (torr, 25C):	0.0500 3
Henry's law const (atm m**3/mol, 25C):	0.84600E-06 3
Solubility in water (mg/L, 25C):	896.00 3
RETENTION PROPERTIES:	Ref.
log (octanol/water) partition (25C):	3.22 3
Biodegradation rate in water (1/month):	Rapid 3

August 1987 CAS #: 131-11-3 PHTHALATES

Dimethyl phthalate

Other Names: DMP; 1,2-Benzenedicarboxylic acid, dimethyl ester Haz Waste #: U102 Ind/EPA Gen #: C6H4 (C00CH3) 2 Formula: PHYSICAL PROPERTIES: - - - -Ref. Molecular weight: 194.20 3 Melting point (C): 5.50 3 Boiling point (C): 282.00 3 Vapor pressure (torr, 25C): ₹ ⟨0.01 0.32400E-06 3 Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): 4320.00 3 RETENTION PROPERTIES: - - - - -- - - Ref. log (octanol/water) partition (25C): 2.12 3 Biodegradation rate in water (1/month): Rapid 3

PHTHALATES August 1987 CAS #: 117-84-0 d-n-Octyl phthalate Other Names: DOP Ind/EPA Gen #: Haz Waste #: U107 PHYSICAL PROPERTIES: - - - - - - -- - - - Ref. Molecular weight: 391.00 3 Melting point (C): -25.00 3 Boiling point (C): 220.00 3 Vapor pressure (torr, 25C): (0.2 3 Solubility in water (mg/L, 25C): 3.0000 3 RETENTION PROPERTIES: - - - - - -- - - - Ref. log (octanol/water) partition (25C): (approx.) 9.20 3 Biodegradation rate in water (1/month): Significant 3 August 1987 PHTHALATES CAS #: 85-44-9 Phthalic anhydride Other Names: 1,2-Benzenedicarboxylic acid anhydride; 1,3-Isobenzofurandione Haz Waste #: U190 Ind/EPA Gen #: KO23,KO24,KO93,KO94 Formula: C8H403 - - - - - Ref. Molecular weight: 148.11 2 Melting point (C): (sublimes) 130.80 2 Boiling point (C): 295.00 2

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

1.5300 2

Soluble in 162 parts water 2

7.3 Halogenated Pesticides

August 1987 CAS #: 309-00-2

HALOGENATED PESTICIDES Aldrin Pesticides

Other Names: 1,2,3,4,10,10-Hexachloro-1,4,4a,5,8,8a-hexahydro-1,4-endoexo-5,8-dimethanonaphthalene; HHDN; compd 118; Octalene Ind/EPA Gen #: Haz Waste #: P004 Formula: C12H8C16 Description: Crystals PHYSICAL PROPERTIES: - - - - - - - - -- - - Ref. Molecular weight: 364.93 2 Melting point (C): 104.00 2 Vapor pressure (torr, 250): 0.60000E-05 3 Henry's law const (atm m**3/mol, 25C): 0.49600E-03 3 Solubility in water (mg/L, 25C): 0.0170 to 0.1800 3 RETENTION PROPERTIES: - - - - - -- - - - Ref. log (octanol/water) partition (25C): 5.17 3 Partition coefficient, soil/water: 7847.00 17 Hydrolysis rate, water (1/month, 250): Slow hydrolysis 3 Biodegradation rate in water (1/month): Biotransforms to dieldrin 3

August 1987

HALOGENATED PESTICIDES

CAS #: 94-75-7

2,4-D Pesticides

Other Names: (2,4-Dichlorophenoxy)acetic acid; Hedonal; Trinoxal

Formula: C6H3Cl2OCH2COOH; C8H6Cl2O3

Description: Crystals; causes irritation to eyes, G.I. disturbances

(Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Molecular weight: 221.04 1
Melting point (C): 138.00 1

Boiling point (C): (0.4) 160.00 1

HALOGENATED PESTICIDES Dieldrin Pesticides

6,7-epoxy-1,4,4a,5,6,7,8,8a- dimethanonphthalene; Compoun HEOD; ENT 16225; Octalox Haz Waste #: PO37 Ind/EPA Gen #: Formula: C12H8C160 Description: A stereoisomer of endrin (PO	d 497; Insecticide no. 497;
PHYSICAL PROPERTIES:	Ref.
Molecular weight:	380.93 2
Melting point (C):	176.00 to 177.00 2 (20 C) 0.17800E-06 3
Vapor pressure (torr, 25C): Henry's law'const (atm m**3/mol, 25C):	(20 C) 0.17800E-06 3 0.58000E-04 3
Solubility in water (mg/L, 25C):	0.2000 3
RETENTION PROPERTIES:	Ref.
Hydrolysis rate, water (1/month, 25C):	Hydrolysis slow 3
Biodegradation rate in water (1/month):	Bioaccumulation moderate 3
	NATED PESTICIDES Endosulfan Pesticides
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5 methano-2,4,3-benzodioxathie	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene-
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5 methano-2,4,3-benzodioxathie 1,4,5,6,7,7-Hexachloro-5-nor sulfite; 1,2,3,4,7,7-Hexachl 5,6-bisoxymethylene sulfite; Haz Waste #: PO50 Ind/EPA Gen #: Formula: C9H6C16O3S Description: Brown crystals	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5 methano-2,4,3-benzodioxathie 1,4,5,6,7,7-Hexachloro-5-nor sulfite; 1,2,3,4,7,7-Hexachl 5,6-bisoxymethylene sulfite; Haz Waste #: PO50 Ind/EPA Gen #: Formula: C9H6Cl6O3S Description: Brown crystals PHYSICAL PROPERTIES:	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex
CAS #: 115-29-7 alpha- Other Names: 6,7,8,9,10,10-Hexachloro-1,5	Endosulfan Pesticides ,5a,6,9,9a-hexahydro-6,9- pin 3-oxide; Malix; Thiodan; bornene-2,3-dimethanol cyclic orobicyclo[2,2,1]-2-heptene- Chlorthiepin; Thionex

August 1987

HALOGENATED PESTICIDES

72-20-8 CAS #: Endrin Pesticides Other Names: 3,4,5,6,9,9-Hexachloro-1a,2,2a,3,6,6a,7,7a-octahydro-2,7:3,6-dimethanonaphth[2,3-b]oxirene; Mendrin; Nendrin; 1,2,3,4,10,10-Hexachloro-6,7-epoxy-1,4,4a,5,6,7,8,8aoctahydro-endo, endo-1, 4:5,8-dimethanonaphthalene; Hexadrin; Compound 269; Experimental insecticide no. 269; ENT 17251 Ind/EPA Gen #: D012 Haz Waste #: P051 Formula: C12H8C160 Description: Crystals; toxic to fish; manufacture and use has been discontinued in the U.S. (Ref. 2) - - - - - - - - Ref. Molecular weight: 380.93 2 Decomposes 245 (235, Ref. 3) Melting point (C): 2 Vapor pressure (torr, 25C): 0.20000E-06 3 Henry's law const (atm m**3/mol, 25C): 0.50000E-06 3 Solubility in water (mg/L, 25C): 0.2600 3 RETENTION PROPERTIES: - - - - -- - - Ref. log (octanol/water) partition (25C):

Partition coefficient, soil/water:

Hydrolysis rate, water (1/month, 25C):

Biodegradation rate in water (1/month):

Bioaccumulation important 3 August 1987 HALOGENATED PESTICIDES CAS #: 640-19-7 Fluoroacetamide pesticides Other Names: 2-Fluoroacetamide; Fluoroacetic acid amide; 1081; Fussol; Monofluroracetamide; Fluorakil 100 Haz Waste #: P057 Ind/EPA Gen #: Formula: CH2FC0NH2 Description: Crystals; sublimes on heating PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 77.06 2 Solubility in water (mg/L, 25C): Freely soluble 2 HALOGENATED PESTICIDES August 1987

CAS #: 62-74-8

Na Fluoracetic acid pesticides

Other Names: Fluoroacetic acid, sodium salt; Compound 1080; 1080; Fratol Haz Waste #: P058 Ind/EPA Gen #:

CH2FC00Na; C2FH2Na02 Formula:

Description: Fine white powder; extremely toxic

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.

Molecular weight:

100.02 17

Solubility in water (mg/L, 25C):

Soluble 2

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HALOGENATED PESTICIDES
August 1987
         76-44-8
CAS #:
                                Heptachlor Pesticides
Other Names: 1,4,5,6,7,8,8-Heptachloro-3a,4,7,7a-tetrahydro-4,7-methano-
            1H-indane; Velsicol 104; Drimox; Drinox; Heptamul;
            1,4,5,6,7,8,8-Heptachloro-3a,4,7,7a-tetrahydro-4,7-
            methanoindene; E3314
Haz Waste #: P059
                         Ind/EPA Gen #:
Formula:
            C10H5C17
Description: Crystals; the EPA has cancelled registration of pesticides
            containing this compound with the exception its use through
            subsurface ground insertion for termite control and the
            dipping of roots or tops of non-food plants (FR vol. 40,
            p. 28850, 7/9/75); poisoning may occur by ingestion.
            inhalation, skin contamination; stimulates CNS (Ref. 2)
----Ref.
Molecular weight:
                                                            373.35
                                                                   2
Melting point (C):
                                               95.00 to
                                                            96.00
Vapor pressure (torr, 250):
                                                       0.30000E-03
Henry's law const (atm m**3/mol, 25C):
                                                       0.14800E-02 3
0.1800 3
                                                           - - - - Ref.
log (octanol/water) partition (25C):
                                                             4.40 3
Partition coefficient, soil/water:
                                                           1580.00 17
Hydrolysis rate, water (1/month, 25C):
Biodegradation rate in water (1/month):
                                                Hydrolysis rapid 3
                                                Will bioaccumulate 3
August 1987
                                HALOGENATED PESTICIDES
CAS #:
        465-73-6
                                Isodrin pesticides
Other Names: Compound 711
Haz Waste #: P060
                         Ind/EPA Gen #:
Formula:
          C12H8C16
```

Description: endo, endo-isomer of Aldrin (P004); cyrstals

PHYSICAL PROPERTIES: - - - - - - -

Molecular weight:

Melting point (C):

- Ref.

364.93 2

242.00

240.00 to

August 1987 CAS #: 8001-35-2

HALOGENATED PESTICIDES Toxaphene Pesticides

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Other Names: Camphene polychlorinates (67-69% CI); Alltox; Geniphene;
            Motox; Penphene; Toxakil; Camphechlor; Synthetic 3956;
            Chlorinated camphene; Polychlorocamphene; Phenacide;
            Phenatox; Strobane-T; Alltex; Anatox; Camphochlor;
            Camphoclor; Chem-phene; Chlorocamphene; Compound 3956;
            Crestoxo; Cristoxo; Clor chem T-590; ENT 9735; Estonox;
            Fasco-terpene; Gy-phene; Hercules toxaphene; Hercules 3956;
             Kamfochlor; M 5055; Melipax; NCI-C00259; PCHK; Phenacide;
            Octachlorocamphene; Polychlorcamphene; Phenatox; Toxadust;
             Toxakil; Toxon 63; Toxyphen; Vertac 90%; (others)
                          Ind/EPA Gen #: D015,K041,K098
Haz Waste #: P123
Formula:
            ca. C10H10C18 (C10H8C110, C10H(18-n)Cln, C10H(16-n)Cln)
Description: Yellow, waxy solid; pleasant piney odor. A very complex,
            but reproducible mixture of at least 177 C10 polychloro
             derivatives. Can cause mild irritation of, and be
            absorbed through, skin. CNS stimulant. Listed as a
            carginogen by the EPA. (Ref. 2); available as a wettable
             powder, emulsifiable concentrate, dust, granule, bait, oil
             and emulsion; can be extremely persistent in soil receiving
             direct application (Ref. 12)
PHYSICAL PROPERTIES: - - - - - - -
                                                     (average) 414.00 3
Molecular weight:
Melting point (C):
                                                 65.00 to
                                                               90.00
Boiling point (C):
                                                     Decomposes >120
Vapor pressure (torr, 25C):
                                              0.2000 to
                                                             0.4000
Henry's law const (atm m**3/mol, 25C):
                                                         0.48900E-02
Solubility in water (mg/L, 25C):
                                              0.5000 to
                                                              3.0000 3
RETENTION PROPERTIES:
                                                            - - - Ref.
log (octanol/water) partition (25C):
                                                                3.30 3
Partition coefficient, soil/water:
                                                              160.00 17
Hydrolysis rate, water (1/month, 25C):

Slow hydrolysis
Biodegradation rate in water (1/month): Biodegrades and bioaccumulates
                                  HALOGENATED PESTICIDES
August 1987
CAS #: 5103-71-9
                                  Chlordane Pesticides
Other Names: 1,2,4,5,6,7,8,8-Octachloro-3a,4,7,7a-tetrahydro-4,7-methano-
             indane; trans-Chlordane
Haz Waste #: U036
                          Ind/EPA Gen #: K032,K033,K034,K097
Formula:
             C10H4C18
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             406.00 3
Melting point (C):
                                                103.00 to
                                                            105.00
Boiling point (C):
                                                  (at 2 torr) 175.00 3
Vapor pressure (torr, 25C):
                                                         0.10000E-04
Henry's law const (atm m**3/mol, 25C):
                                                         0.48000E-04
                                                                     3
RETENTION PROPERTIES:
                                            ----- Ref.
log (octanol/water) partition (25C):
                                                                2.78 3
Partition coefficient, soil/water:
                                                             54.2000 17
Biodegradation rate in water (1/month): Bioaccumulation important 3
```

	ALOGENATED PESTICIDES	
CAS #: 510-15-6	Chlorobenzilate pesticides	
Other Names: Benzeneacetic acid, 4-c alpha-hydroxy, ethyl es Haz Waste #: U038 Ind/EPA (Formula: (C6H4C1)2C(OH) (C0OC2H5) Description: Viscous liquid PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Vapor pressure (torr, 25C): Solubility in water (mg/L, 25C):	ster Gen #:	
	HALOGENATED PESTICIDES o-Chloro-m-cresol pesticides	
Other Names: 4-Chloro-m-cresol; 4-Chloro-3-methylphenol; Chlorocresol; 3-Methyl-4-chlorophenol; Parachlorometacresol; 6-Chloro-m-cresol; 6-Chloro-3-hydroxytoluene; 2-Chloro-5-hydroxytoluene		
Haz Waste #: U039 Ind/EPA (Formula: C7H7CIO	en ⋕:	
Description: Dimorphous crystals; usually a phenolic odor; aqueous solution turns yellow on exposure to light and air (Ref. 2) PHYSICAL PROPERTIES: Ref.		
Molecular weight:	142.58 2	
Melting point (C):	55.5 and 66. (ligroin crystals) 2	
Boiling point (C):	235.00 2 (20 C) 3850.00 3	
Solubility in water (mg/L, 25C): RETENTION PROPERTIES:	(20 C) 3850.00 3	
log (octanol/water) partition (25C):		
Partition coefficient, soil/water:	77.2000 17	
Biodegradation rate in water (1/mont	th): Readily biodegraded 3	
	•	

HALOGENATED PESTICIDES 2-Chlorophenol Pesticides

Other Names: o-Chlorophenol Haz Waste #: U048 Ind/EPA Gen #: Formula: C6H50C1 Description: Liquid PHYSICAL PROPERTIES: - - - -Molecular weight: 128.60 3 9.30 2 Melting point (C): Boiling point (C): 175.00 2 Density (g/cc, 20C): 1.2573 2 (20 C) 2.2000 3 Vapor pressure (torr, 25C): Henry's law const (atm m**3/mol, 25C): 0.82800E-05 log (octanol/water) partition (25C):

Biodegradation rate in water (1/month):

Significant biodegradation 3 HALOGENATED PESTICIDES August 1987 CAS #: 72-54-8 4,4'-DDD Pesticides Other Names: 1,1-Dichloro-2,2-bis(p-chlorophenyl)ethane; TDE; DDD; 1,1'-(2,2-Dichloroethylidene)bis[4-chlorobenzene]; p,p'-DDD; Tetrachlorodiphenylethane; p,p'-TDE; Rhothane; Dichlorodiphenyldichloroethane Haz Waste #: U060 Ind/EPA Gen #: Formula: C14H10C14 Description: Crystals; slightly irritating to skin; symptoms similar to DDT (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref. Molecular weight: 320.05 2 109.00 to 2 Melting point (C): 110.00 (30 C) 0.10200E-05 3 Vapor pressure (torr, 25C): log (octanol/water) partition (25C): 5.99 3 log (octanol/water) partition (25C): 5.99 3 Biodegradation rate in water (1/month): Important biotransformation 3

HALOGENATED PESTICIDES August 1987 CAS #: 789-02-6 4.4'-DDT Pesticides Other Names: p,p'-DDT Haz Waste #: U061 Ind/EPA Gen #: Formula: C14H9C15 PHYSICAL PROPERTIES: - - - - - - -- - - Ref. Molecular weight: 354.50 Melting point (C): 108.50 to 109.00 2 Boiling point (C): 185.00 Vapor pressure (torr, 250): (20 C) 0.19000E-06 3 Solubility in water (mg/L, 25C): $\langle 1.5 \text{ to } 5.5 3 \rangle$ - - - Ref. RETENTION PROPERTIES: log (octanol/water) partition (25C): 3.98 3 Hydrolysis rate, water (1/month, 25C): Hydrolysis may be important 3 Biodegradation rate in water (1/month): Both biotransformation and bioaccumulation are important 3 HALOGENATED PESTICIDES August 1987 CAS #: 2303-16-4 Diallate pesticides Other Names: Carbamothioic acid bis(1-methylethyl)-S-(2,3-dichloro-2propenyl) ester Haz Waste #: U062 Ind/EPA Gen #: ((CH3)2CH)2NCOSCH2CCI=CHCI Formula: Description: Brown liquid PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -- - - Ref. Molecular weight: 270.24 2 Boiling point (C): (9) 150.00 2 (ppm) 40.0000 2 Solubility in water (mg/L, 25C): August 1987 HALOGENATED PESTICIDES CAS #: 96-12-8 1,2-Dibromo-3-chloropropane pesticides Other Names: BBC 12; 3-Chloro-1,2-dibromopropane; DBCP; Fumagon; 1-Chloro-2.3-dibromopropane: Dibromochloropropane: Fumazone; NCI-C00500; Nemabrom; Nemafume; Nemagon; Nemanax; Nemapaz; Nemaset; Nematocide; Nematox; Nemazon; OS 1897:

0xy DBCP; SD 1897 Ind/EPA Gen #: Haz Waste #: U066 C3H5Br2CI Formula: Description: Dark amber to dark brown liquid; moderately volatile and degraded in moist soil; a carcinogen (Ref. 12) - - - Ref. Molecular weight: 236.36 2 Boiling point (C): 196.00 2 Density (g/cc, 20C): 2.0930 2 Vapor pressure (torr, 25C): (mmHg, 21 C) 0.8000 2 Solubility in water (mg/L, 25C): Slightly 2

August 1987

HALOGENATED PESTICIDES CAS #: 106-46-7 TALUGENATED PESTICION TALUGENATE PESTIC

Other Names: p-Dichlorobenzene; Paracide; PDB; Para-zene; UO72 Ind/EPA Gen #: K085,K105 Di-chloricide; Paramoth Haz Waste #: U072 Formula: PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 147.00 1 Melting point (C): 53.04 1 Boiling point (C): 174.10 1 (liquid) 1.2475 1 Density (g/cc, 20C): Vapor pressure (torr, 25C): 1.1800 3 Henry's law const (atm m**3/mol, 25C): 0.28800E-02 3 Solubility in water (mg/L, 25C): 79.0000 3
RETENTION PROPERTIES: - - - - - - - - - - - Ref. log (octanol/water) partition (25C):

Partition coefficient, soil/water:

Biodegradation rate in water (1/month):

3.39 3

193.00 17

Bioaccumulates 3 August 1987 HALOGENATED PESTICIDES CAS #: 120-83-2 2,4-Dichlorophenol Pesticides Other Names: 2.4-DCP Ind/EPA Gen #: Haz Waste #: U081 Formula: C6H4C120 PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref. Molecular weight: 163.00 3 Melting point (C): 45.00 3 Boiling point (C): 210.00 3 Vapor pressure (torr, 25C): 0.1200 3 Henry's law'const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): RETENTION PROPERTIES: --------0.66600E-05 3 4500.00 3 ---- Ref. log (octanol/water) partition (25C):
Hydrolysis rate, water (1/month, 25C):
Biodegradation rate in water (1/month):

2.75 3
Hydrolysis not important 3
Rapid biodegradation 3 HALOGENATED PESTICIDES August 1987 CAS #: 87-65-0 2,6-Dichlorophenol pesticides Haz Waste #: U082 Ind/EPA Gen #: KO43 Formula: C6H4C120 Description: Needles PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 163.00 13 163.00 13 66.00 to 67.00 13 219.00 to 220.00 13 Melting point (C): Boiling point (C):

HALOGENATED PESTICIDES August 1987 Dimethylcarbamoyl chloride pesticides CAS #: 79-44-7 Other Names: N.N-Dimethylcarbamyl chloride; DDC; DMCC; TL 389; Chloroformic acid dimethylamide; (Dimethylamino)carbonyl chloride; Dimethylcarbamic acid chloride; Dimethylcarbamic chloride; N,N-Dimethylcarbamidoyl chloride Ind/EPA Gen #: Haz Waste #: U097 C3H6C1N0 Formula: Description: Colorless liquid; a carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - - - - - - - - - - - -Molecular weight: RETENTION PROPERTIES: ------- - - - - Ref. Hydrolysis rate, water (1/month, 25C): Hydrolyzes rapidly 12 August 1987 HALOGENATED PESTICIDES Hexachlorobenzene CAS #: 118-74-1 Other Names: Perchlorobenzene; Anticarie; Bunt-cure; Bunt-no-more; Julin's carbon chloride (Ref. 2) Ind/EPA Gen #: Haz Waste #: U127 Formula: C6C16 PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - Ref. Molecular weight: 284.80 2 228.70 1 Melting point (C): Boiling point (C): 319.30 1 Density (g/cc, 20C): (kg/L, liquid) 1.5960 1 Vapor pressure (torr, 25C): (kPa) 0.34400E-03 3 Henry's law const (atm m**3/mol, 25C): (kPa m**3/mol, 20 C) 0.50000E-02 Solubility in water (mg/L, 25C): (q/m**3) 0.50000E-02 3HALOGENATED PESTICIDES August 1987 CAS #: 87-68-3 Hexachlorobutadiene Pesticides Other Names: HCBD Haz Waste #: U128 Ind/EPA Gen #: Formula: CC12=CC1-CC1-CC12 PHYSICAL PROPERTIES: - - - - - - - - - - - -Molecular weight: 260.80 3 Melting point (C): -21.00 3 Boiling point (C): 215.00 3 Vapor pressure (torr, 25C): 0.1500 3 Henry's law const (atm m**3/mol, 25C): 0.0103

log (octanol/water) partition (25C):

Partition coefficient, soil/water:

2.0000 3 - - - Ref.

3.74 3

400.00 17

HALOGENATED PESTICIDES Lindane pesticides

```
Other Names: 1,2,3,4,5,6-Hexachlorocyclohexane; gamma-HCH; gamma-Benzene
            hexachloride; gamma Hexachlor; ENT 7796; Aparsin; Aphtiria;
            gamma-BHC; Gammalin; Gamene; Gamiso; Gammexane; Gexane;
            Jacutin; Kwell; Lindafor; Lindatox; Lorexane; Quellada;
            Streunex; Tri-6; Viton; Aalindan; Aficide; Agrisol G-20;
            Agrocide; Agronexit; Ameisenatod; Ben-hex; Bentox 10;
            Celanex; Detmol-extrakt; Detox 25; Hortex; Isotox; Forlin;
            Lindagam; Lindagrain; gamma-Lindane; Lindex; Omnitox;
            Tap 85; Nexol-E; Nexit; Linvur; Gamacid; Gamaphex;
            Gammahexa; Gammahexane; Gammalin; (many others)
Haz Waste #: U129
                         Ind/EPA Gen #: DO13
            C6H6C16
Formula:
Description: Slightly musty odor; crystals; poisoning may occur by
            ingestion, inhalation or percutaneous absorption; this
            and its isomers (alpha-HCH, beta-HCH and technical HCH) have
            been listed as carcinogens by the EPA (Ref.2); crystals are
            colorless; CAS numbers for isomers: alpha--319-84-6,
            beta--319-85-7, technical--608-73-1 (Ref. 12)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             290.85 2
Melting point (C):
                                                            112.50 2
                                  (mmHg, 20 C) 0.94000E-05 2
Vapor pressure (torr, 25C):
Solubility in water (mg/L, 25C):
RETENTION PROPERTIES: -----
                                                            7.8000 9
                                                            - - - Ref.
log (octanol/water) partition (250):
                                                              3.72 16
Partition coefficient, soil/water:
                                                             348.00 17
August 1987
                                 HALOGENATED PESTICIDES
CAS #:
        77-47-4
                                 Hexachlorocyclopentadiene Pesticides
Other Names: HCCPD; 1,2,3,4,5,5-Hexachloro-1,3-cyclopentadiene
Haz Waste #: U130 Ind/EPA Gen #: K034
Formula:
          C5C16
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                            272.80 3
Melting point (C):
                                                             -9.90 3
Boiling point (C):
                                                            239.00 3
Vapor pressure (torr, 25C):
                                                            0.0810 3
Henry's law!const (atm m**3/mol, 25C):
                                                            0.0164 3
1.8000 3
                                                        ---- Ref.
log (octanol/water) partition (25C):
                                                              3.99 3
rog (octanol/water) partition (25C):

Hydrolysis rate, water (1/month, 25C):

Hydrolysis could occur 3
Biodegradation rate in water (1/month):
                                                   Bioaccumulated 3
```

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HALOGENATED PESTICIDES
August 1987
         70-30-4
CAS #:
                                Hexachlorophene pesticides
Other Names: 2,2'-Methylenebis[3,4,6-trichlorophenol]; G-11; AT-7;
            2,2'-Dihydroxy-3,3',5,5',6,6'-hexachlorodiphenylmethane;
            Bis(3,5,6-trichloro-2-hydroxyphenyl)methane; Bilevon;
            Dermadex; Exofene; Gamophen; Hexosan; pHisohex; Surgi-Cen;
            Surofene
Haz Waste #: U132
                         Ind/EPA Gen #:
            (C6HC130H)2CH2; C13H6C1602
Formula:
Description: Crystals; forms salts with alkalies and alkaline earths;
            excessive dosage to animals results in symptoms of
            neurotoxicity; use regulated by FDA (Ref. 2)
Molecular weight:
                                                            406.92
Melting point (C):
                                               164.00 to
                                                            165.00 2
Solubility in water (mg/L, 25C):
                                                  Almost insoluble 2
                                 HALOGENATED PESTICIDES
August 1987
CAS #:
        143-50-0
                                 Kepone Pesticides
Other Names: Chlordecone; Decachlorooctahydro-1,3,4-metheno-2H-
            cyclobuta[cd]pentalen-2-one; GC-1189; CIBA 8514; Clordecone;
            Compound 1189; Decano-4-one; Decachloro-1,3,4-metheno-2H-
            cyclobuta(cd)pentalen-2-one; Decachloroketone; ENT-16391;
            1,1a,3,3a,4,5,5a,5b,6-Decachlorooctahydro-1,3,4-metheno-
            2H-cyclobuta(cd)pentalen-2-one; General Chemicals 1189;
            Decachlorotetracyclodecanone; Merex; NCI-C00191;
            Decachlorooctahydrokepone-2-one
Haz Waste #: U142
                         Ind/EPA Gen #:
            C10C | 100
Formula:
Description: Crystals; has caused tumors in factory workers exposed to
            this substance; listed as a carcinogen by the EPA (Ref. 2)
            no longer manufactured or used in the U.S. (Ref. 12)
- Ref.
Molecular weight:
                                                            490.68 2
Boiling point (C):
                                                  Sublimes at 350 C 3
```

Solubility in water (mg/L, 25C):

(100 C) 4000.00 3

HALOGENATED PESTICIDES August 1987 CAS #: 82-68-8 Pentachloronitrobenzene pesticides Other Names: Quintozene; PCNB; Terrachlor; PKhNB; Avicol; Botrilex; Brassicol; Folosan; Terraclor; Tilcarex; Tritisan Ind/EPA Gen #: Haz Waste #: U185 Formula: C6C15N02 Description: Fine needles, platelets PHYSICAL PROPERTIES: - - - - - - - -- - - Ref. Molecular weight: 295.36 2 Melting point (C): 144.00 Boiling point (C): (some decomposition) 328.00 2 Density (g/cc, 20C): (25 C) 1.7180 2 Solubility in water (mg/L, 25C): Practically insoluble 2 HALOGENATED PESTICIDES August 1987 CAS #: 58-90-2 2,3,4,6-Tetrachlorophenol pesticides Haz Waste #: U212 Ind/EPA Gen #: Formula: C6H2C140 Description: Needles; highly toxic orally (Ref. 13) PHYSICAL PROPERTIES: - - - - -- - - - Ref. Molecular weight: 232.00 13 Melting point (C): 70.00 13 Boiling point (C): (15 mm Hg) 150.00 13

August 1987 HALOGENATED PESTICIDES
CAS #: 95-95-4 2,4,5-Trichlorophenol pesticides

Other Names: Collunosol
Haz Waste #: U230 Ind/EPA Gen #:
Formula: C6H3Cl30
PHYSICAL PROPERTIES: - - - - - - - - - - Ref.

 Molecular weight:
 197.46 1

 Melting point (C):
 (sublimes) 67.00 1

 Boiling point (C):
 253.00 1

August 1987

Molecular weight:

Melting point (C):

Solubility in water (mg/L, 25C):

HALOGENATED PESTICIDES

88-06-2 CAS #:

2,4,6-Trichlorophenol Pesticides Other Names: Dowicide 2S; Omal Ind/EPA Gen #: Haz Waste #: U231 Formula: C6H3C130 Description: Crystals; strong phenolic odor; volatile with steam; forms a sodium salt (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - - Ref. Molecular weight: 197.46 2 Melting point (C): 69.00 2 Boiling point (C): 246.00 2 1.4901 2 Density (g/cc, 20C): Vapor pressure (torr, 25C): 1.0000 3 Henry's law const (atm m**3/mol, 25C): 0.72000E-05 3 Solubility in water (mg/L, 25C):
RETENTION PROPERTIES: -----800.00 3 - - - - Ref. log (octanol/water) partition (25C): 3.38 3 Partition coefficient, soil/water: 189.00 17 Biodegradation rate in water (1/month): Significant 3 HALOGENATED PESTICIDES August 1987 CAS #: 93-76-5 2,4,5-Trichlorophenoxy acetic acid pest. Other Names: 2,4,5-T; 2-(2,4,5-Trichlorophenoxy)propionic acid; Fenoprop; 2,4,5-TC; Miller Nu Set (Hormone Spray) Haz Waste #: U232 Ind/EPA Gen #: KO42 C8H5C1303 Formula: Description: Crystals; irritating to eyes, skin, mucous membranes (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - - Ref. Molecular weight: 255.49 1 Melting point (C): 153.00 1 Density (g/cc, 20C): 1.8000 1 Solubility in water (mg/L, 25C): (mg/kg, 30 C) 238.00 1 August 1987 HALOGENATED PESTICIDES CAS #: 93-72-1 Silvex pesticides Other Names: 2-(2,4,5-Trichlorophenoxy) propionic acid; 2,4,5-TC Formula: C9H7C1303 PHYSICAL PROPERTIES: - - - - - - - - - - - - - -

- - - - Ref.

269.53 1

181.60 1

(%) 0.0140 1

August 1987

HALOGENATED PESTICIDES Pentachlorophenol Pesticides

CAS #: 87-86-5

Other Names: Penta; PCP; Penchlorol; Santophen 20 Ind/EPA Gen #: KOO1 Haz Waste #: U242

C6HC150; C6C150H Formula:

Description: Needle-like crystals; very pungent odor when hot; sublimes

in needles; forms a sodium salt; ingestion causes increase

then decrease in respiration, blood pressure, urinary output, etc. leading to death; Causes lung, liver, kidney damage; contact dermatitis; may be absorbed through skin

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref. Molecular weight: 266.35 2 2 Melting point (C): 190.00 to 191.00 Boiling point (C): 309.00 to 310.00 2 Density (g/cc, 20C): 1.9780 Vapor pressure (torr, 25C): (20 C) 0.11000E-03 Henry's law const (atm m**3/mol, 25C): 0.88200E-05 3 Solubility in water (mg/L, 25C): RETENTION PROPERTIES: - - - - - - -14.0000 3 - - - - Ref. log (octanol/water) partition (25C): 5.01 3 5624.00 17 Partition coefficient, soil/water: Biodegradation rate in water (1/month): Significant 3

August 1987

HALOGENATED PESTICIDES

CAS #: 1888-71-7

Hexachloropropene pesticides

Other Names: 1,1,2,3,3,3-Hexachloro-1-propene; Hexachloropropylene

Haz Waste #: U243 Ind/EPA Gen #:

C12C=CC1CC13; C3C16 Formula:

Description: Highly toxic by inhalation; irritant (Ref. 13)

- - - - Ref. Molecular weight: 249.00 13 Boiling point (C): 209.00 to 210.00 13

Density (q/cc, 20C):

1.7600 13

August 1987

HALOGENATED PESTICIDES

CAS #: 72-43-5

Methoxychlor

Other Names: 1,1'-(2,2,2-Trichlorethylidene)-bis[4-methyoxybenzene];

1,1,1-Trichloro-2,2-bis(p-methoxypehnyl)ethane; Marlate;

2,2-Di-p-anisyl-1,1,1-trichloroethane; DMDT; Methoxy-DDT

Ind/EPA Gen #: D014 Haz Waste #: U247

Formula: C16H15C1302

Description: Dimorphic crystals

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -

Molecular weight:

Melting point (C):

78. to 78.2 or 86. to 88. (dimorphic) 2

Solubility in water (mg/L, 25C): Practically insoluble 2 7.4 Polynuclear Aromatics

August 1987 CAS #: 53-96-3

POLYNUCLEAR AROMATICS 2-Acetylaminofluorene

Other Names: N-2-Fluorenylacetamide; N-9H-Fluoren-2-ylacet-amide; AAF; 2-FAA; 2-AAF; 2-Acetamidofluorene; 2-Acetaminofluorene; N-Acetyl-2-aminofluorene; 2-(Acetylamino)fluorene; FAA; N-Fluoren-2-ylacetamide; 2-Fluorenylacetamide Haz Waste #: U005 Ind/EPA Gen #: C15H13N0 Formula: Description: Crystals; listed as a carcinogen by the EPA PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -- - - Ref. 223.26 2 Molecular weight: 194.00 2 Melting point (C): Insoluble 2 Solubility in water (mg/L, 25C): POLYNUCLEAR AROMATICS August 1987 CAS #: 225-51-4 Benz(c)acridine Other Names: 1.2-Benzacridine; alpha-Chrysidine; alpha-Naphthacridine; alpha-Phenonaphthacridine; 2,1-Naphthacridine Haz Waste #: U016 Ind/EPA Gen #: Formula: C17H11N Description: Air pollutant; brilliant yellow needles; experimental carcinogen (Ref. 13) PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref. Molecular weight: 229.00 13 Melting point (C): 108.00 13 POLYNUCLEAR AROMATICS August 1987 CAS #: 56-55-3 Benz[a]anthracene Other Names: 1,2-Benzanthracene; 2,3-Benzphenanthrene; Tetraphene; BA; Benzanthrene; Naphthanthracene; 1,2-Benz(a)anthracene; Benzanthracene; 1,2-Benzanthrene; Benzo(a)anthracene; Benzo(a) pehnanthrene; Benzo(b) pehnanthrene; Benzoanthracene; 1,2-Benzoanthracene; 2,3-Benzopehnanthrene Haz Waste #: U018 Ind/EPA Gen #: Formula: C18H12 Description: Plates; greenish-yellow fluorescence; listed as a carcinogen by the EPA (Ref. 2); no reported commercial use or application (Ref. 12) PHYSICAL PROPERTIES: -----Ref. 228.30 3 Molecular weight: Melting point (C): 157.00 3 155.00 to Boiling point (C): Sublimes 3 Vapor pressure (torr, 25C): 0.50000E-08 3 Solubility in water (mg/L, 25C): 0.0140 3 RETENTION PROPERTIES: - - - Ref. log (octanol/water) partition (25C): 5.61 3 0.19608E+05 17 Partition coefficient, soil/water: Biodegradation rate in water (1/month): Slow 3

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August 1987
                                   POLYNUCLEAR AROMATICS
CAS #:
         50-32-8
                                   Benzo[a]pyrene
Other Names: 3,4-Benzpyrene; (formerly 1,2-Benzpyrene); B(a)P; BP;
             Benz(a)pyrene; 3,4-Benz(a)pyrene; Benzo(d,e,f)chrysene;
             .6,7-Benzopyrene; 3,4-Benzypyrene; 3,4-BP
                           Ind/EPA Gen #:
Haz Waste #: U022
             C20H12
Formula:
Description: Yellowish plates; crystals may be monoclinic or
             orthorhombic; listed as a carcinogen by the EPA (Ref. 2)
             combustion byproduct; a constituent of creosote (Ref. 12)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                                 252.30 3
Melting point (C):
                                                                 179.00 3
Boiling point (C):
                                                   310.00 to
                                                                 312.00 3
Vapor pressure (torr, 25C):
                                                           0.50000E-08 3
Henry's law const (atm m**3/mol, 25C):
                                                                 0.0126 3
Solubility in water (mg/L, 25C): 0.38000E-02 3 RETENTION PROPERTIES: - - - - - - - - - - - Ref.
log (octanol/water) partition (25C):
Biodegradation rate in water (1/month):
                                                                   6.04 3
                                                            Biodegraded 3
August 1987
                                 POLYNUCLEAR AROMATICS
CAS #: 218-01-9
                                   Chrysene
Other Names: Benzo(a)phenanthrene
```

Haz Waste #: U050 Ind/EPA Gen #: Formula: C18H12 PHYSICAL PROPERTIES: - - - - - - -- - - Ref. Molecular weight: 228.30 3 Melting point (C): 256.00 3 Boiling point (C): 448.00 3 - - - - - Ref. log (octanol/water) partition (25C): 5.61 3 Partition coefficient, soil/water: 0.19600E+05 17 Biodegradation rate in water (1/month): Bioaccumulated 3

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POLYNUCLEAR AROMATICS
August 1987
CAS #:
         53-70-3
                                  Dibenz[a,h]anthracene
Other Names: 1,2:5,6-Dibenzanthracene; DB(a,h)A; 1,2,5,6-DBA;
             1,2:5,6-Benzanthracene; Dibenzo(a,h)anthracene
Haz Waste #: U063
                          Ind/EPA Gen #:
            C22H14
Formula:
Description: Plates, leaflets; crystals may be monoclinic or
            orthorhombic; listed as a carcinogen by the EPA (Ref. 2);
            no reported commercial application; a product of incomplete
            combustion (Ref. 12)
PHYSICAL PROPERTIES: - - - - - - - -
                                                             - - - - Ref.
Molecular weight:
                                                              278.40
                                                                     3
Melting point (C):
                                                              270.00
                                                                     3
Boiling point (C):
                                                            Sublimes 3
Vapor pressure (torr, 250):
                                                (approx.) 0.10000E-09
Solubility in water (mg/L, 25C):
                                                         0.50000E-02 3
RETENTION PROPERTIES:
                                                           - - - - Ref.
log (octanol/water) partition (25C):
                                                                5.93 3
Biodegradation rate in water (1/month):
                                                     Degraded slowly 3
                                  POLYNUCLEAR AROMATICS
August 1987
CAS #:
        189-55-9
                                  Dibenzo(a,i)pyrene
Other Names: DB(a,i)P; 1,2,7,8-Dibenzopyrene; Debenz(a,i)pyrene;
             3,4:9,10-Dibenzopyrene; 3,4:9,10-Dibenzpyrene
            Benzo[rst]pentaphene
Haz Waste #: U064
                          Ind/EPA Gen #:
Formula:
            C24H14
Description: Not commercially produced, but found as a combustion
             product; a carcinogen (Ref. 12)
PHYSICAL PROPERTIES: - - - - - - - - -
                                                              302.37 17
Molecular weight:
Solubility in water (mg/L, 25C):
                                                           Insoluble 12
August 1987
                                  POLYNUCLEAR AROMATICS
CAS #:
         57-97-6
                                  7,12-Dimethylbenz[a]anthracene
Other Names: 9,10-Dimethyl-1,2-benzanthracene; 1,4-Dimethyl-2,3-
             benzphenanthrene
Haz Waste #: U094
                          Ind/EPA Gen #:
Formula:
             C20H16
Description: Plates, leaflets; faint greenish-yellow tinge
- - - - Ref.
Molecular weight:
                                                              256.33 2
Melting point (C):
                                                122.00 to
                                                              123.00 2
```

Solubility in water (mg/L, 25C):

Insoluble 2

POLYNUCLEAR AROMATICS

Fluoranthene

Other Names: Benzo(j,k)fluorene Haz Waste #: U120 Ind/EPA Gen #: Formula: C16H10 PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref. Molecular weight: 202.30 3 111.00 3 log (octanol/water) partition (25C): Partition coefficient, soil/water: 5.33 3 0.11000E+05 17 Biodegradation rate in water (1/month): Significant 3 POLYNUCLEAR AROMATICS August 1987 CAS #: 193-39-5 Indeno (1,2,3-c,d) pyrene Other Names: 2,3-o-Phenylenepyrene; IP; 1,10-(o-Phenylene)pyrene; 1,10-(1,2-Phenylene)pyrene; o-Phenylenepyrene; 2,3-Phenylenepyrene Ind/EPA Gen #: Haz Waste #: U137 Formula: C22H12 Description: A carcinogen; not commercially manufactured; a byproduct of combustion (Ref. 12) - - - Ref. Molecular weight: 276.30 3 Melting point (C): 162.00 to 164.00 3

 Melting point (C):
 162.00 to 164.00 3

 Vapor pressure (torr, 25C):
 (approx.) 0.10000E-09 3

 Solubility in water (mg/L, 25C):
 0.6200 3

 RETENTION PROPERTIES:
 - - - - - - - - - Ref.

 log (octanol/water) partition (25C): 7.66 3 Biodegradation rate in water (1/month): Slow 3 August 1987 POLYNUCLEAR AROMATICS CAS #: 56-49-5 3-Methylcholanthrene Other Names: 1,2-Dihydro-3-methyl-benz[j]aceanthrylene; 3-MECA; 3-MC; 20-Methylcholanthrene Haz Waste #: U157 Ind/EPA Gen #: Formula: C21H16 Description: Pale yellow, slender prisms PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 268.34 2 Melting point (C): 179.00 to 180.00 2 Boiling point (C): (80 mm Hg) 280.00 2 Density (g/cc, 20C): 1.2800 2 Solubility in water (mg/L, 25C): Insoluble 2

August 1987 CAS #: 91-20-3 POLYNUCLEAR AROMATICS Naphthalene

Other Names: Naphthene; Naphthalin; Tar camphor Haz Waste #: U165 Ind/EPA Gen #: K023,K024

Formula: C10H8

Description: Monoclinic prismatic plates; also sold as white scales,

powder, balls or cakes; odor of moth balls; volatilizes appreciably at room temperature; sublimes appreciably at temperatures above melting point; poisoning may occur by ingestion of large doses, inhalation or skin absorption; symptoms range from nausea and vomiting to convulsions and

coma (Ref. 2)

coma (nor. 2)		
PHYSICAL PROPERTIES:		Ref.
Molecular weight:	128.20	3
Melting point (C):	80.60	3
Boiling point (C):	(sublimes) 218.00	3
Vapor pressure (torr, 25C):	0.0492	3
Henry's law const (atm m**3/mol, 25C):	0.36000E-03	3
Solubility in water (mg/L, 25C):	34.4000	3
RETENTION PROPERTIES:		Ref.
log (octanol/water) partition (25C):	3.37	3
Partition coefficient, soil/water:	185.00	17
Biodegradation rate in water (1/month):	Rapid	3

7.5 Oxygenated

OXYGENATED August 1987 CAS #: 460-19-5 Cyanogen Other Names: Ethanedinitrile; Dicyan; Oxalic acid dinitrile Haz Waste #: PO31 Ind/EPA Gen #: Formula: C2N2 Description: Highly poisonous gas; almond-like odor; toxic effects similar to those of hydrogen cyanide (Ref. 2) - - - Ref. Molecular weight: 52.04 2 -27.90 2 Melting point (C): -21.17 Boiling point (C): 2 (-21.17 C) 0.9537 2 Density (g/cc, 20C): Solubility in water (mg/L, 25C): 4 vols of gas in 1 vol of water 2 August 1987 OXYGENATED CAS #: 506-77-4 Cyanogen chloride Haz Waste #: P033 Ind/EPA Gen #: Formula: CNCI Description: Liquid; vapors are highly irritant and very poisonous; toxic effects similar to those of hydrogen cyanide (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 61.48 2 Melting point (C): (solidifies) -6.00 2 13.80 2 Boiling point (C): Density (g/cc, 20C): 1.1860 2 Soluble 2 Solubility in water (mg/L, 25C): OXYGENATED August 1987 CAS #: 624-83-9 Isocyanic acid methyl ester Other Names: Methyl isocyanate; Isocyanatomethane Haz Waste #: P064 Ind/EPA Gen #: CH3NCO; C2H3NO Formula: Description: Liquid with powerful odor; highly toxic by inhalation and absorption; lacrimator; extremely flammable (Ref. 13) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.

Molecular weight:

Boiling point (C):

Density (g/cc, 20C):

57.00 13

45.00 13

0.9600 13

43.00 to

August 1987 OXYGENATED CAS #: 107-12-0 Propanenitrile Other Names: Propionitrile; Ethyl cyanide Haz Waste #: P101 Ind/EPA Gen #: CH3CH2CN; C3H5N Formula: Description: Liquid; poisonous when heated to decomposition or on contact with acids; pleasant, ethereal, sweetish odor (Ref. 2) PHYSICAL PROPERTIES: - -- - Ref. Molecular weight: 55.08 2 Melting point (C): -91.80 2 Boiling point (C): 97.20 2 Density (g/cc, 20C): 0.7818 2 Solubility in water (mg/L, 25C): 11.9 g/100 g at 40 C 2 August 1987 **OXYGENATED** CAS #: 75-07-0 Aceta I dehyde Other Names: Ethanal; Ethylaldehyde; "Aldehyde"; Acetic aldehyde Haz Waste #: U001 Ind/EPA Gen #: K009,K010 Formula: CH3CH0; C2H40 Description: Flammable liquid; characteristic, pungent odor; general narcotic action; large doses may cause death by respiratory paralysis; symptoms of chronic intoxication resemble those of chronic alcoholism; irritating to mucous membranes (Ref. 2) PHYSICAL PROPERTIES: - - - - -- - - Ref. Molecular weight: 44.05 3 Melting point (C): -124.00 3 Boiling point (C): 21.00 3 Density (g/cc, 20C):

Vapor pressure (torr, 25C):

Henry's law const (atm m**3/mol, 25C):

Solubility in water (mg/L, 25C):

RETENTION PROPERTIES: - - - - -

Biodegradation rate in water (1/month):

(16 C) 0.7880

0.32400E-04 3

Miscible 3

2

740.00 3

--- Ref.

High 3

August 1987 CAS #: 98-86-2

OXYGENATED Acetophenone

Other Names: 1-Phenylethanone; Phenyl methyl ketone; Acetylbenzene; Hypnone Ind/EPA Gen #: Haz Waste #: U004 Formula: C6H5COCH3; C8H8O Description: Liquid; forms laminar crystals at low temperatures PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref. 120.15 2 Molecular weight: 20.50 2 Melting point (C): Boiling point (C): 202.00 2 (15 C) 1.0330 2 Density (g/cc, 20C): Solubility in water (mg/L, 25C): Slightly soluble 2 **OXYGENATED** August 1987 CAS #: 98-82-8 Cumene Other Names: (1-Methylethyl)benzene; Cumol; Isopropylbenzene Haz Waste #: U055 Ind/EPA Gen #: K022 Formula: C9H12 Description: Colorless liquid; narcotic in high concentrations (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 120.19 2 Boiling point (C): 153.00 2 152.00 to Density (g/cc, 20C): 0.8620 2 Solubility in water (mg/L, 25C): Insoluble 2 **OXYGENATED** August 1987 CAS #: 1464-53-5 Diepoxybutane Other Names: 2,2'-Bioxirane; 1,2:3,4-Diepoxybutane Haz Waste #: U085 Ind/EPA Gen #: Formula: C4H6O2

Vapor pressure (torr, 25C):

OXYGENATED

Formal dehyde Other Names: Methanal; Oxymethylene; Oxomethane; Methylene Oxide; Formic aldehyde; Methyl aldehyde; BFV; Fannoform; Formalin; Formaldehyde solution; Formalith; Formol; Fyde; Hoch; Ivalon; Karsan; NCI-CO2799; Oxymethylene Ind/EPA Gen #: HCHO; CH20 Description: Flammable, colorless gas at ordinary temperatures; pungent suffocating odor; very reactive; polymerizes easily; very irritating to mucous membranes; listed as a carcinogen by the EPA (Ref. 2); most frequently marketed as a 37-56 percent aqueous solution stabilized with 10-15 percent methanol to prevent polymerization (Ref. 12) PHYSICAL PROPERTIES: - - - - - - -- - - - Ref. Molecular weight: 30.00 3 Melting point (C): -92.00 3 -118.00 to Boiling point (C): -21.00 to -19.00 Density (g/cc, 20C): (air=1.00) 1.0670 2 Vapor pressure (torr, 25C): 10.0000 3 Solubility in water (mg/L, 25C): Very soluble, up to 55% 2 RETENTION PROPERTIES: ----- Ref. Biodegradation rate in water (1/month): High 3 **OXYGENATED** August 1987 CAS #: 64-18-6 Formic acid Other Names: Methanoic acid; Formylic acid Haz Waste #: U123 Ind/EPA Gen #: HC00H; CH202 Formula: Description: Colorless liquid; pungent odor; strong reducing agent; dangerously caustic to skin (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - -- - - - - Ref. Molecular weight: 46.03 3 Melting point (C): 8.40 3 Boiling point (C): 100.50 Density (g/cc, 20C): 1.2200 2

40.0000 3

```
OXYGENATED
August 1987
        765-34-4
                                 Glycidaldehyde
CAS #:
Other Names: Epihydrinaldehyde; 2,3-Epoxy-1-propanal; Glycidal;
            Epihydrine aldehyde; 2,3-Epoxypropionaldehyde;
            Oxirane-carboxaldehyde
                          Ind/EPA Gen #:
Haz Waste #: U126
Formula:
            C3H402
Description: Colorless liquid
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                              72.10 14
                                                             113.00 14
Boiling point (C):
                                                             1.1403 14
Density (g/cc, 20C):
                                 DXYGENATED
August 1987
CAS #:
        123-63-7
                                 Paraldehyde
Other Names: Paracetaldehyde
                          Ind/EPA Gen #:
Haz Waste #: U182
Formula:
            C6H1203
Description: Liquid; characteristic aromatic odor; warm, but disagreeable
            taste; abuse may lead to habituation or addiction (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             132.16 2
Melting point (C):
                                                              12.00 2
Boiling point (C):
                                                             124.00 2
Density (g/cc, 20C):
                                                       (25 C) 0.9940 2
Solubility in water (mg/L, 25C): Soluble in eight parts water
                                  OXYGENATED
August 1987
CAS #: 106-51-4
                                 p-Quinone
Other Names: Quinone; 2,5-Cyclohexadiene-1,4-dione; 1,4-Benzoquinone;
            1,4-Cyclohexadienedione
Haz Waste #: U197
                      Ind/EPA Gen #:
Formula:
            C6H402
Description: Yellow monoclinic prisms; penetrating odor resembling
            chlorine; irritating vapors; can cause dermatitis; vapors
            can damage eyes (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             108.09 2
Melting point (C):
                                                   (sublimes) 115.70 2
                                                             1.3180 2
Density (g/cc, 20C):
Solubility in water (mg/L, 25C):
                                                   Slightly soluble 2
```

August 1987 CAS #: 636-21-5 **OXYGENATED**

o-Toluidine hydrochloride

Other Names: 2-Amino-1-methylbenzene hydrochloride; NCI-CO2335;

1-Amino-2-methylbenzene hydrochloride; o-Aminotoluene

hydrochloride; 2-Aminotoluene hydrochloride;

1-Methyl-2-aminobenzene hydrochloride;

2-Methyl-1-aminobenzene hydrochloride; o-Methylaniline

hydrochloride; o-Methylaniline hydrochloride;

o-Methylbenzenamine hydrochloride; 2-Methylbenzenamine hydrochloride; 2-Toluidine hydrochloride; o-Tolylamine

hydrochloride

Haz Waste #: U222

Ind/EPA Gen #:

Formula:

C7H9N.HCI

Description: Listed as a carcinogen by the EPA (Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.

Molecular weight:

143.62 17

7.6 Polymerizables

POLYMERIZABLE Acrolein

```
Other Names: 2-Propenal; Allyl aldehyde; Acrylic aldehyde; Acrylaldehyde;
            Acraldehyde; Aqualin
Haz Waste #: P003
                         Ind/EPA Gen #:
Formula:
            CH2=CH-CH0
Description: Flammable liquid; pungent odor; irritates eyes and mucosa;
            unstable (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             56.06 2
                                                            -87.70 3
Melting point (C):
Boiling point (C):
                                                             52.50 3
Density (g/cc, 20C):
                                                            0.8389 2
Vapor pressure (torr, 25C):
                                                            220.00 3
Henry's law const (atm m**3/mol, 250):
                                                       0.77000E-04 3
Solubility in water (mg/L, 25C):
                                                    (?) 0.40000E+06 3
RETENTION PROPERTIES:
                                                   log (octanol/water) partition (25C):
                                                             -0.09 3
Partition coefficient, soil/water:
                                                            0.1400 17
Biodegradation rate in water (1/month):
                                                             Rapid 3
                                 POLYMERIZABLE
August 1987
CAS #:
      107-18-6
                                 Allyl alcohol
Other Names: 2-Propen-1-ol; 1-Propenol-3; Vinyl carbinol
Haz Waste #: P005
                         Ind/EPA Gen #:
Formula:
            CH2=CH-CH20H
Description: Colorless liquid; pungent, mustard-like odor; irritating
            to the eyes; causes severe irritation of mucous membranes
            (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                            58.08
                                                                   3
Melting point (C):
                                                            -50.00
                                                                   3
Boiling point (C):
                                                            97.00 3
                                                96.00 to
Density (g/cc, 20C):
                                                            0.8540 2
Vapor pressure (torr, 25C):
                                                           23.8000 3
Henry's law const (atm m**3/mol, 25C): (15 C) 0.10000E-05 3
Solubility in water (mg/L, 25C):
RETENTION PROPERTIES: ------
                                                          Miscible 3
                                                         - - - - Ref.
```

Biodegradation rate in water (1/month):

Degraded 3

August 1987 POLYMERIZABLE CAS #: 151-56-4 Ethylenimine Other Names: Aziridine; azacyclopropane; dimethtylenimine Haz Waste #: P054 Ind/EPA Gen #: Formula: C2H5N Description: Liquid; intense odor of ammonia; poisonous; strongly alkaline; strongly irritating to eyes, skin, mucous membranes PHYSICAL PROPERTIES: - -- - - Ref. 43.07 2: Molecular weight: 57.00 2 Boiling point (C): 56.00 to Density (g/cc, 20C): (24 C) 0.8321 Solubility in water (mg/L, 25C): Miscible 2 **POLYMERIZABLE** August 1987 CAS #: 75-55-8 2-Methylaziridine Other Names: Propyleneimine Haz Waste #: P067 Ind/EPA Gen #: Formula: C3H7N Description: Fuming oil; highly toxic; suspected human carcinogen; may polymerize explosively in presence of acids (Ref. 13) - - - - Ref. Molecular weight: 57.00 13 Boiling point (C): 66.00 to 67.00 13 Density (g/cc, 20C): (16 C) 0.8120 13 **POLYMERIZABLE** August 1987 CAS #: 75-86-5 2-Methyllactonitrile Other Names: Acetone cyanohydrin; alpha-Hydroxyisobutyronitrile; USAF RH-8; 2-Hydroxy-2-methyl-propanenitrile Haz Waste #: P069 Ind/EPA Gen #: Formula: C4H7N0

Description: Poisonous PHYSICAL PROPERTIES: - - -

Molecular weight:

Melting point (C):

Boiling point (C):

Density (g/cc, 20C):

- - Ref.

85.12 14

-20.00 14

(23 mm Hg) 82.00 14

(19 C) 0.9320 14

```
August 1987
                                  POLYMERIZABLE
        107-19-7
CAS #:
                                  Propargyl alcohol
Other Names: 2-Propyn-1-ol
Haz Waste #: P102
                          Ind/EPA Gen #:
            HC=-CCH20H; C3H40
Formula:
Description: Moderately volatile liquid; mild geranium odor; irritating
            to skin and mucous membranes (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - -
                                                              - - - Ref.
Molecular weight:
                                                               56.06 2
                                                 -52.00 to
                                                              -48.00 2
Melting point (C):
                                                              115.00 2
Boiling point (C):
                                                 114.00 to
                                                              0.9715 2
Density (g/cc, 20C):
                                                            Miscible 2
Solubility in water (mg/L, 25C):
                                  POLYMERIZABLE
August 1987
CAS #:
         79-06-1
                                  Acrylamide
Other Names: Propenamide
Haz Waste #: U007
                          Ind/EPA Gen #:
            CH2=CHC0NH2; C3H5N0
Formula:
Description: Monomer--flake-like crystals; highly toxic and irritant;
            causes CNS paralysis; can be absorbed through unbroken skin
PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.
Molecular weight:
                                                               71.08 2
Melting point (C):
                                                               84.50 2
Boiling point (C):
                                                    (25 mm Hg) 125.00
                                                                     2
Density (g/cc, 20C):
                                                       (30 C) 1.1220 2
Solubility in water (mg/L, 25C):
              215.5 q/kg (30 C); polymer--soluble and insoluble forms 2
August 1987
                                  POLYMERIZABLE
CAS #:
        79-10-7
                                  Acrylic acid
Other Names: 2-Propenoic acid; Vinylformic acid
Haz Waste #: U008
                          Ind/EPA Gen #:
            CH2=CHC02H; C3H402
Description: Corrosive liquid; acrid odor and fumes; strong irritant;
             polymerizes readily in the presence of oxygen (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                               72.06 2
Melting point (C):
                                                               14.00
                                                                     2
```

Boiling point (C):

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

141.00 2

Miscible 2

(16 C) 1.0621 2

POLYMERIZABLE Acrylonitrile

Other Names: Vinyl cyanamide; Propenonitrile; ABS; Absaf Abson; Acrylon; B 32; Bakelite; Blendex; Car Cevian V; Cyanoethylene; Cycolac; Dow 500; EPA 3530; EPB 3570; Formid; Forsan; Fortyler Goodyear 600; Hycar 1441; Kralastic; Kralon Lastilac; Lustran; Lustropak; Lustrum; Novoc 2-Propenenitrile; Ravikral; RExene; Ronfalin Stylac; Terluran; Tyoylac; VCN; Ventox; Viny Acritet; (many other commercial names) Haz Waste #: U009	rbacryl; ENT 54; ne; Fumigrain; ; Lacqran; dur; n; Sicoflex; yl cyanide; as a carcinogen
PHYSICAL PROPERTIES: Molecular weight:	Ref. 53.06 3
Melting point (C):	-82.00 3
Boiling point (C):	78.50 3
Vapor pressure (torr, 25C):	100.00 3
Henry's law const (atm m**3/mol, 25C):	0.66600E-04 3
Solubility in water (mg/L, 25C):	0.73500E+05 3
RETENTION PROPERTIES:	Ref.
log (octanol/water) partition (25C):	-0.14 3
Partition coefficient, soil/water:	0.1200 17
Biodegradation rate in water (1/month):	Significant 3
August 1987 POLYMERIZABLE CAS #: 123-73-9 Crotonaldehyde	
Other Names: 2-Butenal Haz Waste #: U053 Ind/EPA Gen #: Formula: CH3CH=CHCHO	, ,
PHYSICAL PROPERTIES:	Ref.
Molecular weight:	70.09 3
Melting point (C):	-74.00 3
Boiling point (C): 104.00	
Henry's law const (atm m**3/mol, 25C):	0.14000E-04 3
Solubility in water (mg/L, 25C):	0.18000E+06 3
RETENTION PROPERTIES:	Ref.
Biodegradation rate in water (1/month):	Rapid 3

POLYMERIZABLE August 1987 CAS #: 140-88-5 Ethyl acrylate Other Names: 2-Propenoic acid ethyl ester; Acrylic acid ethyl ester Haz Waste #: U113 Ind/EPA Gen #: CH2=CHC00CH2CH3; C5H802 Formula: Description: Monomer--liquid; acrid, penetrating odor retained by clothing; lacrimator; easily polymerized (Ref. 2) Polymer--transparent, elastic substance; practically no odor; little adhesive power; resists usual solvents (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - -- - - - Ref. Molecular weight: 100.11 2 Melting point (C): -72.00 2 99.40 2 Boiling point (C): 2 g/100 mL at 20 C 2 Solubility in water (mg/L, 25C): August 1987 POLYMERIZABLE CAS #: 75-21-8 Ethylene oxide Other Names: Oxirane; Anprolene Haz Waste #: U115 Ind/EPA Gen #: Formula: C2H40 Description: Colorless, flammable gas; explosive; highly irritating to eyes, mucous membranes (Ref. 2) PHYSICAL PROPERTIES: - - - - - - -- - - Ref. Molecular weight: 44.05 2 Melting point (C): -111.00 2 Boiling point (C): 10.70 2 0.8694 2 Density (g/cc, 20C): Solubility in water (mg/L, 25C): Soluble 2 RETENTION PROPERTIES: - - - - -- - - Ref. Hydrolysis rate, water (1/month, 25C): 0.40 15 August 1987 POLYMERIZABLE CAS #: 97-63-2 Ethyl methacrylate Other Names: Ethylmethyl acrylate; Ethyl-alpha-methyl acrylate; Ethyl-2-methylacrylate; Ethyl-2-methyl-2-propenoate Haz Waste #: U118 Ind/EPA Gen #: Formula: C6H1002 Description: Liquid PHYSICAL PROPERTIES: - -- - Ref. Molecular weight: 114.16 14

Melting point (C):

Boiling point (C):

Density (g/cc, 20C):

< -75. C 14

(25 C) 0.9110 14

119.00 14

POLYMERIZABLE August 1987 CAS #: 108-31-6 Maleic anhydride Other Names: 2,5-Furandione; cis-Butenedionic anhydride; Toxilic anhydride Haz Waste #: U147 Ind/EPA Gen #: Formula: C4H2O3 Description: Orthorhombic needles; powerful irritant; causes burns; inhalation can cause pulmonary edema; in Malathion (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - Ref. 98.06 2 Molecular weight: 52.80 2 Melting point (C): Boiling point (C): 202.00 2 1.4800 2 Density (g/cc, 20C): Solubility in water (mg/L, 25C): Soluble forming maleic acid 2 POLYMERIZABLE August 1987 CAS #: 109-77-3 Malononitrile Other Names: Propanedinitrile; Methylene cyanide; Dicyanomethane; Cvanoacetonitrile Haz Waste #: U149 Ind/EPA Gen #: Formula: CH2(CN)2; C3H2N2 Description: Colorless solid PHYSICAL PROPERTIES: - - - - - - - - - - - - -Molecular weight: 66.06 2 Melting point (C): 32.00 2 Boiling point (C): 218.00 to 219.00 2 Density (g/cc, 20C): 1.1910 2 Soluble 2 Solubility in water (mg/L, 25C): POLYMERIZABLE August 1987 CAS #: 126-98-7 Methacrylonitrile Other Names: 2-Methyl-2-propenenitrile; Isopropenylnitrile; alpha-Methylacrylonitrile; Isopropene cyanide Haz Waste #: U152 Ind/EPA Gen #: CH2=C(CH3)C=N; C4H5NDescription: Liquid; lacrimator; inisidious poison; delayed skin reaction (Ref. 2) PHYSICAL PROPERTIES: - - - - -- - - - Ref. Molecular weight: 67.09 2 -35.80 2 Melting point (C): Boiling point (C): 90.30 2 Density (g/cc, 20C): 2 0.8001

Solubility in water (mg/L, 25C):

2.57 wt % at 20 C 2

POLYMERIZABLE August 1987 CAS #: 80-62-6 Methyl methacrylate Other Names: 2-Methylpropenoic acid methyl ester Haz Waste #: U162 Ind/EPA Gen #: CH2=C (CH3) - C00CH3 Formula: Description: Polymerizes easily forming a clear plastic known as Lucite, Plexiglas, Perspex (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. 100.10 3 Molecular weight: -48.00 3 Melting point (C): Boiling point (C): 100.00 to 101.00 3 Vapor pressure (torr, 25C): 37.0000 3 Solubility in water (mg/L, 25C): Slightly 3 - - - - Ref. Biodegradation rate in water (1/month): Rapid 3 August 1987 **POLYMERIZABLE** CAS #: 504-60-9 1,3-Pentadiene Other Names: Piperylene Haz Waste #: U186 Ind/EPA Gen #: H3CCH=CHCH=CH2; C5H8 Formula: Description: Liquid PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 68.00 13 Boiling point (C): 42.00 13 August 1987 POLYMERIZABLE CAS #: 95-80-7 Toluene diamine Other Names: Diaminotoluene; 2,4-Diaminotoluene; 5-Amino-o-toluidine; 3-Amino-p-toluidine; Benzofur MT; C.I. Oxidation Base 20; C.I. 76035; Developer B; 2,4-Diamino-1-methylbenzene; 1,3-Diamino-4-methylbenzene; 2,3-Diamino-1-toluene; 2,4-Diaminotoluol; Eucanine GB; Fouramine; Fourrine M; MTD; Nako TMT; NCI-C02302; Pelagol grey J; Fourrine 94; 4-Methyl-1,3-benzenediamine; 4-Methyl-m-phenylenediamine Haz Waste #: U221 Ind/EPA Gen #: Formula: C7H10N2 Description: Used in the synthesis of polyurethanes; crystal; used in a 4:1 mixture with 2-6 Diaminotoluene to make Toluene Diisocyanate; a carcinogen (Ref. 12)

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Molecular weight:

Solubility in water (mg/L, 25C):

122.17 17

Soluble 12

August 1987 CAS #: 584-84-9 POLYMERIZABLE Toluene diisocyanate

Other Names: Toluene 2,4-diisocyanate; 2,4-Diisocyanatotoluene; 2,4-tolylene diisocyanate; TDI; Nacconate 100 Haz Waste #: U223 Ind/EPA Gen #: KO27 C9H6N202 Formula: Description: Liquid; sharp pungent odor; causes skin irritation, allergic eczema, also bronchial asthma (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - -

Ref. Molecular weight: 174.15 2 Melting point (C): 21.50 2 19.50 to Boiling point (C): 251.00 2 Density (g/cc, 20C): 1.2244

Solubility in water (mg/L, 25C):

Reacts with water with evolution of CO2 2

August 1987

POLYMERIZABLE Urethane

CAS #: 51-79-6

Other Names: Urethan; Carbamic acid ethyl ester; Ethyl carbamate;

Ethyl urethan

Haz Waste #: U238 Ind/EPA Gen #:

NH2C00C2H5 Formula:

Description: Crystals; cooling saline taste

PHYSICAL PROPERTIES: - - - - - - - -Ref. Molecular weight: 89.09 2 Melting point (C): 2 48.00 to 50.00 Boiling point (C): 182.00 to 184.00 2 Density (g/cc, 20C): 1.1000 2 Solubility in water (mg/L, 25C): 1 q/0.5 mL 2

7.7 Phenolics

```
PHENOLICS
August 1987
CAS #:
          51-28-5
                                   2,4-Dinitrophenol
Other Names: Aldifen; alpha-Dinitrophenol
Haz Waste #: P048,U104
                           Ind/EPA Gen #:
Formula:
             C6H3(N02)2(OH)
Description: Yellow'orthorhombic crystals; highly toxic; readily
             absorbed through intact skin; vapors absorbed through
             respiratory tract (Ref. 2)
PHYSICAL PROPERTIES: - - - - - -
                                                                 - - - Ref.
Molecular weight:
                                                                184.11 2
Melting point (C):
                                                                114.00 3
Boiling point (C):
                                                              Sublimes 3
                                                                1.6830 2
Density (g/cc, 20C):
Solubility in water (mg/L, 25C):
                                                        (18 C) 5600.00 3
RETENTION PROPERTIES:
                                                              - - - Ref.
log (octanol/water) partition (25C):
                                                                  1.53 3
Biodegradation rate in water (1/month):
                                                           Significant 3
                                   PHENOLICS
August 1987
CAS #: 8001-58-9
                                   Creosote
Other Names: Coal creosote; Brick oil; Coal tar creosote; Coal tar oil;
             Creosote from coal tar; Creosote oil; Creosotum; Heavy oil;
             Cresylic creosote; Liquid pitch oil; Naphthalene oil;
             Tar oil; Wash oil
                           Ind/EPA Gen #:
Haz Waste #: U051
Formula:
Description: A distillate of coal tar consisting of liquid and solid
             aromatic hydrocarbons, tar acids and tar bases; translucent
             brown to black, oily liquid; characteristic sharp odor;
             (Ref. 2); a carcinogen (Ref. 12)
PHYSICAL PROPERTIES: - - - - - - - - - - -
```

Density (q/cc, 20C):

Solubility in water (mg/L, 25C):

(38 C, typical value) 1.0600 2

Practically insoluble 2

Significant 3

Biodegradation rate in water (1/month):

PHENOLICS August 1987 CAS #: 100-02-7 4-Nitrophenol Other Names: p-Nitrophenol Ind/EPA Gen #: Haz Waste #: U170 Formula: C6H4 (NO2) 0H Description: Colorless to slightly yellow, odorless crystals; sweetish then burning taste (Ref. 2) - - - - - - - - - - - - - Ref. PHYSICAL PROPERTIES: - - - - - - -Molecular weight: 139.10 3 (sublimes) 115.00 3 Melting point (C): Boiling point (C): (decomposes) 279.00 3 Vapor pressure (torr, 25C): 2.2000 3 Solubility in water (mg/L, 25C): 0.16000E+05 3 ----Ref. RETENTION PROPERTIES: - - - - - - - - - log (octanol/water) partition (250): 1.91 3 Biodegradation rate in water (1/month): Significant 3 August 1987 PHENOLICS CAS #: 108-95-2 Phenol Other Names: Carbolic acid Haz Waste #: U188 Ind/EPA Gen #: K022 Formula: C6H5OH PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - - Ref. 94.11 3 Molecular weight: 40.90 3 Melting point (C): Boiling point (C): 182.00 3 Henry's law const (atm m**3/mol, 25C): 0.13000E-06 3 0.93000E+05 3 - - - - Ref. log (octanol/water) partition (25C): Biodegradation rate in water (1/month): 1.46 3 Significant 3 **PHENOLICS** August 1987 CAS #: 108-46-3 Resorcinol Other Names: 1,3-Dihydroxybenzene Haz Waste #: U201 Ind/EPA Gen #: Formula: C6H4(OH)2 PHYSICAL PROPERTIES: - - - - - - - -Molecular weight: 110.10 3 Melting point (C): 276.00 to 280.00 3 Boiling point (C): 281.00 3 Vapor pressure (torr, 25C): (138 C) 5.0000 3

Rapid 3

Biodegradation rate in water (1/month):

7.8 Nonhalogenated Solvents

August 1987 CAS #: 75-15-0

NONHALOGENATED SOLVENTS

Carbon disulfide

Other Names: Carbon bisulfide; Dithiocarbonic anhydride Ind/EPA Gen #: F005 Haz Waste #: P022 Formula: S=C=S Description: Highly refractive, mobile, very flammable liquid; poisonous; acute fire and explosion hazard; poisoning may be caused by inhalation, ingestion and skin absorption (Ref. 2) - - - Ref. Molecular weight: 76.14 3 -111.00 3 Melting point (C): Boiling point (C): 46.30 3 (vapor--2.67 (air=1)) 1.2632 2 Density (g/cc, 20C): Vapor pressure (torr, 25C): 360.00 3 Henry's law const (atm m**3/mol, 25C): 0.0133 3 Solubility in water (mg/L, 25C): (20 C) 2940.00 3 NONHALOGENATED SOLVENTS August 1987 CAS #: 107-15-3 Ethylene diamine Other Names: 1,2-Ethanediamine Haz Waste #: P053 Ind/EPA Gen #: H2NCH2CH2NH2 Formula: - - - Ref. Molecular weight: 60.11 3 Melting point (C): 8.50 3 116.00 3 Boiling point (C): Vapor pressure (torr, 25C): (21 C) 10.0000 3 Solubility in water (mg/L, 25C): Miscible 3 RETENTION PROPERTIES: - - - -· - - - Ref. Biodegradation rate in water (1/month): Quantitative 3 August 1987 NONHALOGENATED SOLVENTS CAS #: 57-55-6 Propylene glycol Other Names: 1,2-Propanediol; Methyl glycol; 1,2-Dihydroxypropane Haz Waste #: P100 Ind/EPA Gen #: Formula: CH3CH0HCH20H; C3H802 Description: dl-form--hygroscopic, viscous liquid; slightly acrid taste; tents to oxidize at high temperatures; also I and d forms (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - -- - - Ref. 76.09 2 Molecular weight: Melting point (C): -59.00 2 Boiling point (C): dl-form--188.2; I form--88.-90. (12 mm Hg); d-form--94-96. Density (g/cc, 20C): dl-form--1.036; d-form--1.04 Solubility in water (mg/L, 25C): Miscible 2

NONHALOGENATED SOLVENTS August 1987 CAS #: 67-64-1 Acetone Other Names: 2-Propanone; Dimethyl ketone; beta-Keto-propane; Pyroacetic ether Ind/EPA Gen #: F003,K022 Haz Waste #: U002 CH3COCH3; C3H60 Formula: Description: Volatile, highly flammable liquid; pungent, sweetish taste; characteristic odor; serious poisoning rare; keep away from fire and plastics (solvent for plastics) (Ref. 2) - - - Ref. Molecular weight: 58.08 2 Melting point (C): -94.00 2 Boiling point (C): 56.50 2 Density (g/cc, 20C): (25 C) 0.7880 2 Solubility in water (mg/L, 25C): Miscible 2 NONHALOGENATED SOLVENTS August 1987 CAS #: 75-05-8 Acetonitrile Other Names: Methyl cyanide; Cyanomethane; Ethanenitrile Haz Waste #: U003 Ind/EPA Gen #: CH3CN; C2H3N Formula: Description: Liquid; ether-like odor; poisonous (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - -- - - Ref. Molecular weight: 41.05 2 Melting point (C): -45.00 2 Boiling point (C): 81.60 2 Density (g/cc, 20C): (15 C) 0.7875 2

Solubility in water (mg/L, 25C):

Miscible 2

August 1987

NONHALOGENATED SOLVENTS Aniline

CAS #: 62-53-3

Other Names: Aminobenzene; Benzenamine; Aniline oil; Phenylamine;

Aminophen; Kyanol

Ind/EPA Gen #: K083,K103,K104 Haz Waste #: U012

Formula: C6H5NH2; C6H7N

Description: Oily liquid; colorless when freshly distilled, darkens on

exposure to light or air; poinonous; characteristic odor and burning taste; cumbustible; volatile with steam; intoxication may occur from inhalation, ingestion or

cutaneous absorption; many salts (Ref. 2)

PHYSICAL PROPERTIES:				Ref.
Molecular weight:			93.10	3
Melting point (C):			-6.00	3
Boiling point (C):			184.00	3
Density (g/cc, 20C):			1.0220	2
Vapor pressure (torr, 25C):			0.3000	3
Solubility in water (mg/L, 25C):		•	0.34000E+05	
RETENTION PROPERTIES:	- -			Ref.
Biodegradation rate in water (1/month):			Rapid	3

August 1987

NONHALOGENATED SOLVENTS

71-43-2 CAS #:

Benzene

Other Names: Cyclohexatriene; Benzol; Benzin; Benzine; Benzole; Phene;

Benzolene; Bicarburet of hydrogen; Carbon oil; Pyrob; Coal naphtha; Mineral naphtha; Motor benzol; Pyrobenzol; NCI-C55276; Nitration benzene; Phenyl hydride; Pyrobenzole

Haz Waste #: U019 Ind/EPA Gen #: KO25

Formula: C6H6

Description: Clear, colorless, volatile, highly flammable liquid; a

major raw material of the chemical industry; listed as a

carcinogen by the EPA (Ref. 2, 12)	
PHYSICAL PROPERTIES:	Ref.
Molecular weight: 78.11	2
Melting point (C): 5.53	2
Boiling point (C): 80.10	2
Density (g/cc, 20C): 0.8787	2
Vapor pressure (torr, 25C): 95.2000	3
Henry's law const (atm m**3/mol, 25C): 0.55500E-02	5
Solubility in water $(mg/L, 25C)$: 0.17800E+12	
RETENTION PROPERTIES:	Ref.
log (octanol/water) partition (25C): 2.13	3
Biodegradation rate in water (1/month): Biodegraded	3

NONHALOGENATED SOLVENTS

CAS #: 71-36-3 n-Butyl alcohol Other Names: 1-Butanol; Butyl alcohol; Propyl carbinol Haz Waste #: U031 Ind/EPA Gen #: F003 CH3CH2CH2CH2OH; C4H1OO Description: Highly refractive liquid; vapors irritate and cause cough - - - Ref. Molecular weight: 74.12 2 Melting point (C): -90.00 2 118.00 2 Boiling point (C): 117.00 to Density (g/cc, 20C): 0.8100 2 Solubility in water (mg/L, 25C): 9.1 ml/100 ml 2 NONHALOGENATED SOLVENTS August 1987 CAS #: 110-82-7 Cyclohexane Other Names: Hexahydrobenzene Haz Waste #: U056 Ind/EPA Gen #: Formula: C6H12 PHYSICAL PROPERTIES: - - - - - - - - - - - -- - Ref. Molecular weight: 84.16 3 Melting point (C): 6.30 3 Boiling point (C): 81.00 3 Vapor pressure (torr, 25C): 77.0000 3 Henry's lawwconst (atm m**3/mol, 25C): 0.1600 3 Solubility in water (mg/L, 25C): (20 C) 55.0000 3 August 1987 NONHALOGENATED SOLVENTS CAS #: 108-94-1 Cyclohexanone Other Names: Ketohexamethylene; Pimelic ketone; Hytrol O; Anone; Nadone Haz Waste #: U057 Ind/EPA Gen #: F003 Formula: C6H100 Description: Oily liquid; odor reminiscent of peppermint and acetone; vapor harmful (Ref. 2)

PHYSICAL PROPERTIES: - - - - - -Molecular weight: 98.14 2 Melting point (C): -32.102Boiling point (C): 155.60 2 Density (g/cc, 20C): 0.9478 2 Solubility in water (mg/L, 25C): 87 g/L at 20 C 2

```
August 1987
                                   NONHALOGENATED SOLVENTS
CAS #:
         123-91-1
                                   1,4-Dioxane
Other Names: 1,4-Diethylene dioxide; Dioxane; Di(ethylene oxide);
             Diethylene dioxide; Diethylene ether; Diethylene oxide;
             Diokan; 1,4-Dioxacyclohexane; Dioxan; 1,4-Dioxan;
             Dioxyethylene ether; Glycol ethylene ether; NCI-CO3689;
             NE 220; Tetrahydro-p-dioxin; Tetrahydro-1,4-dioxin
                           Ind/EPA Gen #:
Haz Waste #: U108
             C4H802
Formula:
Description: Flammable liquid; faint, pleasant odor; vapor harmful; may
             cause CNS depression; listed as a carcinogen by the EPA
             (Ref. 2)
PHYSICAL PROPERTIES: - -
Molecular weight:
                                                                 88.10 2
Melting point (C):
                                                                 11.80 2
Boiling point (C):
                                                                101.10 2
Density (g/cc, 20C):
                                                                1.0329 2
Solubility in water (mg/L, 25C):
                                                    Infinitely soluble 12
August 1987
                                   NONHALOGENATED SOLVENTS
CAS #: 141-78-6
                                   Ethyl acetate
Other Names: Acetic acid ethyl ester; Acetic ether; Vinegar naphtha
Haz Waste #: U112
                           Ind/EPA Gen #: F003
             CH3C00C2H5; C4H802
```

Description: Clear, volatile, flammable liquid; characteristic fruity odor; pleasant taste when diluted (Ref. 2)

- - - Ref.

2

88.10 2

77.00 2

24.00 15

0.9020 2

1 mL/10 mL 2

- - - - Ref.

-83.00

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -

RETENTION PROPERTIES: ------

Molecular weight:

Melting point (C):

Boiling point (C):

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

Hydrolysis rate, water (1/month, 25C):

August 1987 CAS #: 60-29-7

Boiling point (C):

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

NONHALOGENATED SOLVENTS

Ethyl ether Other Names: 1,1'-Oxybisethene; Ethoxyethane; Ether; Diethyl ether; Ethyl oxide; Diethyl oxide; Sulfuric ether; Anesthetic ether Ind/EPA Gen #: F003 Haz Waste #: U117 C2H50C2H5; C4H100 Formula: Description: Mobile, very volatile, highly flammalbe liquid; vapor heavier than air; characteristic, sweetish, pungent odor; burning taste; tends to form explosive peroxides under the influence of air and light; mildly irritating to skin, mucous membranes; inhalation of high concentrations may cause unconsciousness and death due to respiratory paralysis; explosion hazard (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - -Molecular weight: 74.12 2 Melting point (C): -116.3 (stable crystals); -123.3 (metastable crystals) Boiling point (C): 34.60 2 Density (g/cc, 20C): 0.7134 2 Vapor pressure (torr, 25C): (mm Hg at 20 C) 439.80 2 Solubility in water (mg/L, 25C): 6.05% (w/w)NONHALOGENATED SOLVENTS August 1987 CAS #: 110-00-9 Furan Other Names: Furfuran; Oxole; Tetrole; Divinylene oxide Haz Waste #: U124 Ind/EPA Gen #: Formula: C4H40 Description: Liquid; varpors are narcotic; can be absorbed through skin (Ref. 2) PHYSICAL PROPERTIES: - - ----- Ref. Molecular weight: 68.07

(758 mm Hg) 32.00

(19.4 C) 0.9371

Insoluble 2

2

2

August 1987 CAS #: 98-01-1

NONHALOGENATED SOLVENTS Furfural

Other Names: Furfuraldehyde; 2-Furancarboxaldehyde; 2-Furaldehyde; Pyromucic aldehyde; Artificial oil of ants; "Furfurol" Haz Waste #: U125 Ind/EPA Gen #: C40CH0: C5H402 Formula: Description: Colorless oily liquid; peculiar odor; turns yellow to brown on exposure to air and light and resinifies; irritates mucous membranes and acts on CNS; causes lacrimation, inflammation of eyes, irritation of throat, headache PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. 96.09 3 Molecular weight: -38.70 3 Melting point (C): Boiling point (C): 162.00 3 Density (g/cc, 20C): (25 C) 1.1563 2 Vapor pressure (torr, 25C): (approx.) 1.6500 3 Solubility in water (mg/L, 25C): (? C) 0.91000E+05 3 RETENTION PROPERTIES: - - - - - - -Biodegradation rate in water (1/month): Quantitative 3 NONHALOGENATED SOLVENTS August 1987 CAS #: 78-83-1 Isobutanol Other Names: 2-Methyl-1-propanol; Isobutyl alcohol Haz Waste #: U140 Ind/EPA Gen #: F005 (H3C) 2CHCH20H; C4H100 Formula: Description: Colorless, refractive liquid; flammable; moderately toxic; eye irritant (Ref. 13) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 74.00 13 -108.00 13 Melting point (C): Boiling point (C): 108.10 13 Solubility in water (mg/L, 25C): Sparingly soluble 13 August 1987 NONHALOGENATED SOLVENTS CAS #: 67-56-1 Methanol Other Names: Methyl alcohol; Carbinol; Wood spirit; Wood alcohol Ind/EPA Gen #: F003 Haz Waste #: U154 Formula: CH30H; CH40 Description: Flammable, poisonous, mobile liquid; slight alcoholic odor; poisoning may occur from ingestion, inhalation or percutaneous absorption (Ref. 2) - - - - - - - - Ref. Molecular weight: 32.04 2 Melting point (C): -97.80 2 Boiling point (C): 64.70 2 Density (g/cc, 20C): (vapor density=1.11; air=1) 0.7915 2 Solubility in water (mg/L, 25C): Miscible 2

August 1987 CAS #: 78-93-3

NONHALOGENATED SOLVENTS Methyl ethyl ketone

August 1987 CAS #: 108-10-1 NONHALOGENATED SOLVENTS
Methyl isobutyl ketone

Other Names: Isopropylacetone; 4-Methyl-2-pentanone; Hexone

Haz Waste #: U161 Ind/EPA Gen #: F003

Formula: CH3COCH2CH(CH3)2; C6H12O

Description: Colorless liquid; faint, ketonic and camphor odor

August 1987

NONHALOGENATED SOLVENTS

CAS #: 79-46-9 2-Nitropropane

Haz Waste #: U171 Ind/EPA Gen #: Formula: CH3CH(NO2)CH3; C3H7NO2

Description: Liquid

 PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.

 Molecular weight:
 89.09 2

 Melting point (C):
 -93.00 2

 Boiling point (C):
 120.30 2

 Density (g/cc, 20C):
 (25 C) 0.9821 2

 Solubility in water (mg/L, 25C):
 Slightly soluble 2

NONHALOGENATED SOLVENTS August 1987 109-06-8 CAS #: 2-Picoline Other Names: alpha-Picoline; 2-Methylpyridine Haz Waste #: U191 Ind/EPA Gen #: Formula: C6H7N Description: Colorless liquid; strong, unpleasant odor; irritating to respiratory tract (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - -- - - Ref. 93.12 2 Molecular weight: Melting point (C): -70.002Boiling point (C): 128.00 to 129.00 2 Density (q/cc, 20C): (15 C) 0.9500 2 Solubility in water (mg/L, 25C): Freely soluble 2 August 1987 NONHALOGENATED SOLVENTS CAS #: 110-86-1 Pyridine Haz Waste #: U196 Ind/EPA Gen #: F005 Formula: C5H5N Description: Flammable, colorless liquid; characteristic disagreeable odor; sharp taste; may cause CNS depression, irritation of skin and respiratory tract (Ref. 2) ---- Ref. PHYSICAL PROPERTIES: - - - - - - - - - - -Molecular weight: 79.10 2 Melting point (C): -42.00 2 Boiling point (C): 115.00 to 116.00 2 Density (g/cc, 20C): (25 C) 0.9780 2 Solubility in water (mg/L, 25C): Miscible 2 NONHALOGENATED SOLVENTS August 1987 CAS #: 109-99-9 Tetrahydrofuran Other Names: Diethylene oxide; Tetramethylene oxide Haz Waste #: U213 Ind/EPA Gen #: Formula: C4H80 Description: Liquid; ether-like odor; irritating to skin, eyes, mucous membranes; narcotic in high concentrations (Ref. 2) - - Ref. Molecular weight: 72.10 2 Melting point (C): -108.50 2 Boiling point (C): 66.00 2

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

0.8892 2

Miscible 2

August 1987 NONHALOGENATED SOLVENTS CAS #: 108-88-3 Toluene

Other Names: Toluol Ind/EPA Gen #: F005,K036 Haz Waste #: U220 Formula: C6H5CH3 PHYSICAL PROPERTIES: . - Ref. Molecular weight: 92.13 3 -95.00 3 Melting point (C): Boiling point (C): 111.00 3 Vapor pressure (torr, 25C): 28.7000 3 Henry's law!const (atm m**3/mol, 25C): 0.59300E-02 3 535.00 3 - - - - Ref. Biodegradation rate in water (1/month): Significant 3

NONHALOGENATED SOLVENTS August 1987 CAS #: 95-47-6 o-Xylene

Other Names: o-Dimethylbenzenes

Haz Waste #: U239 Ind/EPA Gen #: F003,K093,K094

Formula: CH3C6H4CH3
Description: Colorless liquid

- - - Ref.

Molecular weight: 106.20 3 Melting point (C): -25.20 3 144.00 3 Boiling point (C): Vapor pressure (torr, 25C): 10.0000 3

Henry's law const (atm m**3/mol, 25C): 0.61200E-02 3 7.9 Reactive (Non-Cyanide)

August 1987 CAS #: 131-74-8

REACTIVE (NON-CYANIDE) Ammonium picrate

Other Names: 2,4,6-Trinitrophenol ammonium salt; Picric acid ammonium salt; ammonium picronitrate; ammonium carbazoate Haz Waste #: P009 Ind/EPA Gen #: C6H2 (NO2) 30NH4 Formula: Description: Explodes easily from heat or shock (as solid); bright yellow bitter scales or orthorhombic crystals (Ref. 2) Molecular weight: 246.14 2 Density (g/cc, 20C): 1.7200 2

August 1987

REACTIVE (NON-CYANIDE)

1 q/100 mL (20 C) 2

CAS #: 60-34-4 Methylhydrazine

Other Names: Monomethylhydrazine; MMH Haz Waste #: P068 Ind/EPA Gen #:

Formula: CH3NHNH2; CH6N2

Solubility in water (mg/L, 25C):

Description: Reducing agent; clear liquid; ignites spontaneously on

contact with strong oxidizing agents (Ref. 2)

- - - - Ref. Molecular weight: 46.07 2 Melting point (C): -52.40 1 Boiling point (C): 87.50 1 (g/mL, 25 C) 0.2900 1 Density (g/cc, 20C): Vapor pressure (torr, 25C): (kPa) 6.6200 1 Solubility in water (mg/L, 25C): Miscible 2

REACTIVE (NON-CYANIDE) Nitroglycerine

```
Other Names: 1,2,3-Propanetriol; 1,2,3-Propanetriol trinitrate; Glyceryl
             trinitrate; Glycerol nitric acid triester; Nitroclycerol;
             Glonoin; Trinitrin; Blasting gelatin; Blasting oil; S.N.G.;
             Angibid; Angiolingual; Anginine; Angorin; Cardamist;
             Corditrine; Gilucor "nitro"; Klavikordal; Lenitral;
             Myoglycerin; Nitrobid; Nitrodisc; Nitro-Dur; Nitroglin;
             Nitro Mack; Nitromel; Nitronal; NitroPRN;
             Nitrorectal; Nitroretard; Nitrostat; Nitrozell retard;
             Nysconitrine; Percutol; Perlinganit; Perglottal;
             Transderm-Nitro; Suscard; Sustac; Tridil; Trinalgon; others
Haz Waste #: PO81
                           Ind/EPA Gen #:
             CH2(0N02)CH(0N02)CH20N02; C3H5N309
Formula:
Description: Pale yellow, oily liquid; sweet, burning taste; produces
             headache on tasting; explodes on rapid heating or on
             concussion; crystallizes in two forms: labile and stable;
             appreciably volatile at 100 C; evolves nitrous yellow vapors
             at 135 C; acute poisonous; used for manufacture of
             dynamite (75% nitroglycerin) and as a vasodilator; begins to
             deocmpose at 100 C (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - -
                                                                - - - Ref.
Molecular weight:
                                                                227.09
Melting point (C):
                                 2.8 (labile form); 13.5 (stable form)
Boiling point (C):
                                                     (explodes) 218.00
                                                                        2
Vapor pressure (torr, 25C):
                                                (mm, 20 C) 0.26000E-03 2
Solubility in water (mg/L, 25C):
                                                            1 g/800 mL
August 1987
                                   REACTIVE (NON-CYANIDE)
CAS #:
        509-14-8
                                   Tetranitromethane
Haz Waste #: P112
                           Ind/EPA Gen #:
Formula:
          C (NO2) 4
Description: Liquid
                                                                  - - Ref.
PHYSICAL PROPERTIES: - -
Molecular weight:
                                                                196.04 2
Melting point (C):
                                                                13.80 2
Boiling point (C):
                                                                126.00 2
Density (g/cc, 20C):
                                                         (25 C) 1.6229 2
```

August 1987 REACTIVE (NON-CYANIDE) CAS #: 75-36-5 Acetyl chloride Ind/EPA Gen #: Haz Waste #: U006 Formula: CH3COC1; C2H3C10 Description: Flammable liquid; pungent odor; extremely irritating to the eyes; decomposed violently by water or alcohol; irritant; corrosive; causes severe burns (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. 78.50 2 Molecular weight: -112.00 2 Melting point (C): Boiling point (C): 52.00 2 Hydrolysis rate, water (1/month, 25C): Decomposes violently 2 REACTIVE (NON-CYANIDE) August 1987 CAS #: 98-09-9 Benzenesulfonyl chloride Other Names: Benzenesulfonic (acid) chloride; Benzene sulfonechloride Haz Waste #: U020 Ind/EPA Gen #: C6H5S02C1 Formula: Description: Colorless, oily liquid PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - - Ref. 176.62 2 Molecular weight: Melting point (C): 14.50 2 (decomposes) 251.00 to 252.00 2 Boiling point (C): Hydrolysis rate, water (1/month, 25C): Hydrolysis rate, .00294 sec**-1 10 REACTIVE (NON-CYANIDE) August 1987 CAS #: 98-07-7 Benzotrichloride Other Names: (Trichloromethyl)benzene; Phenylchloroform; alpha, alpha, alpha-Trichlorotoluene; Benzenyl trichloride; omega, omega, omega-Trichlorotoluene; Toluene trichloride Haz Waste #: U023 Ind/EPA Gen #: C7H5C13 Formula: Description: Liquid; fumes in air; vapors highly irritant to skin, mucous membranes (Ref. 2); suspected carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref. Molecular weight: 195.48 2 Melting point (C): -5.00 2 Boiling point (C): 220.80 2 Density (g/cc, 20C): 1.3756 2

RETENTION PROPERTIÉS: --------

Hydrolysis rate, water (1/month, 25C):

----- Ref.

Hydrolyzes 2

August 1987
CAS #: 1615-80-1

Other Names: 1,2-Diethylhydrazine; Hydrazoethane
Haz Waste #: U086

REACTIVE (NON-CYANIDE)
Diethylhydrazine
Hydrazoethane
Hydrazoethane

CH3CH2NHNHCH3CH2; C4H12N2

Formula:

Molecular weight:

Boiling point (C):

August 1987 REACTIVE (NON-CYANIDE)
CAS #: 80-15-9 alpha,alpha-Dimethylbenzyl hydroperoxide

Other Names: 1-Methyl-1-phenylethyl-hydroperoxide
Haz Waste #: U096 Ind/EPA Gen #:
Formula: C9H1202
PHYSICAL PROPERTIES: - - - - - - - Ref.
Molecular weight: 152.19 17

- - - - Ref.

88.00 13

86.00 13

85.00 to

August 1987 REACTIVE (NON-CYANIDE)
CAS #: 57-14-7 1,1-Dimethylhydrazine

Other Names: unsym-Dimethylhydrazine; asym-Dimethylhydrazine;

N,N-Dimethylhydrazine; UDMH; Dimazine

Haz Waste #: U098 Ind/EPA Gen #:

Formula: (CH3) 2NNH2; C2H8N2

Description: Flammable, hygroscopic, mobile liquid; fumes in air and

gradually turns yellow; highly corrosive and irritating to skin, eyes, mucous membranes; convulsant poison; base in

rocket fuel formulations (Ref. 2)

 PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.

 Molecular weight:
 60.10 2

 Melting point (C):
 -58.00 2

 Boiling point (C):
 63.90 2

 Density (g/cc, 20C):
 (22 C) 0.7910 2

 Solubility in water (mg/L, 25C):
 Miscible 2

August 1987 REACTIVE (NON-CYANIDE) CAS #: 540-73-8 1,2-Dimethyl hydrazine Other Names: N.N-dimethylhydrazine Haz Waste #: U099 Ind/EPA Gen #: Formula: CH3NHNHCH3 PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - - Ref. Molecular weight: 60.10 2 (753 torr) 81.00 2 Boiling point (C): 0.8274 2 Density (g/cc, 20C): Solubility in water (mg/L, 25C): Miscible 2 August 1987 REACTIVE (NON-CYANIDE) CAS #: 77-78-1 Dimethyl sulfate Other Names: Sulfuric acid mimethyl ester; Dimethyl monosulfate; Methyl sulfate; DMS Ind/EPA Gen #: Haz Waste #: U103 (CH3)2S04; C2H604S Formula: Description: Colorless, oily liquid; extremely hazardous; no warning characteristics; delayed appearance of symptoms; liquid produces severe blistering and necrosis of skin; listed as a carcinogen by the EPA (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - ----- Ref. Molecular weight: 126.13 2 Melting point (C): -27.00 2 (decomposes) 188.00 2 Boiling point (C): Density (g/cc, 20C): 1.3322 2

Solubility in water (mg/L, 25C): 2.8 g/100 mL (18 C) RETENTION PROPERTIES: -----Ref. Hydrolysis rate, water (1/month, 25C): Rapid hydrolysis 2

August 1987 REACTIVE (NON-CYANIDE) CAS #: 122-66-7 1,2-Dipehnylhydrazine

Other Names: Hydrazobenzene; Hydrazodibenzene; 1,1'-Hydrazobisbenzene;

(sym)-Dipehnylhydrazine; N,N'-Diphenylhydrazine;

NCI-C01854

Haz Waste #: U109 Ind/EPA Gen #: Formula: C12H12N2

Description: A carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - - - - -

Molecular weight: 184.24 10 Solubility in water (mg/L, 25C): 0.22100E-03 10 - - - Ref. RETENTION PROPERTIES: - - - - - -

log (octanol/water) partition (25C): 2.94 10

Partition coefficient, soil/water: 75.6000 17

- - - - Ref.

REACTIVE (NON-CYANIDE) August 1987 CAS #: 302-01-2 Hydrazine Other Names: Diamine Haz Waste #: U133 Ind/EPA Gen #: H2NNH2 Formula: Description: Liquid; reducing agent PHYSICAL PROPERTIES: - - - - - -- - Ref. Molecular weight: 32.05 2 Boiling point (C): 113.50 2 Density (g/cc, 20C): (25 C) 1.0036 2 Solubility in water (mg/L, 25C): Miscible 2 REACTIVE (NON-CYANIDE) August 1987 CAS #: 1338-23-4 Methyl ethyl ketone peroxide Other Names: 2-Butanone peroxide; NCI-C55447; Methy lethy | ketonhydroperoxide Haz Waste #: U160 Ind/EPA Gen #: Formula: C8H1604 Description: Skin and eye irritant PHYSICAL PROPERTIES: - - - - - -Molecular weight: 176.24 14 REACTIVE (NON-CYANIDE) August 1987 CAS #: 99-35-4 sym-Trinitrobenzene Other Names: 1,3,5-Trinitrobenzene; Benzite Haz Waste #: U234 Ind/EPA Gen #: Formula: C6H3(NO2)3; C6H3N3O6 Description: Orthorhombic bipyramidal plates; can be sublimed by careful heating; explodes when headed rapidly; less sensitive to impact than TNT, but more powerful and brisant (Ref. 2) Molecular weight: 213.11 2

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

1.7600 2

0.035 q/100 q 2

7.10 Nonhalogenated Pesticides

NONHALOG. PESTICIDES August 1987 Hydroxycoumarin salts CAS #: 81-81-2 Other Names: Warfarin; 4-Hydroxy-3-(3-oxo-1-phenylbutyl)-2H-1-benzopyran-2-one; 3-(alpha-acetonylbenzyl)-4-hydroxy-coumarin; 1-(4'hydroxy-3'-coumarinyl)-1-phenyl-3-butanone; 3-alpha-phenylbeta-acetylethyl-4-hydroxycoumarin: Compound 42: Coumadine: WARF compound 42; Athrombine-K; Sodium salt--Marevan; Prothromadin; Tintorane; Warfilone; Waran Haz Waste #: P001,U248 Ind/EPA Gen #: C19H1604; Sodium salt, C19H15Na04 Formula: Description: Crystals; Sodiums salt--Slightly bitter, crystalline powder Depresses formation of prothrombin and increases capillary fragility, leading to hemorrhages (Ref. 2) Molecular weight: 308.32 2 Melting point (C): 161.; 117.-118. for acidic enol which forms salts 2 NONHALOG. PESTICIDES August 1987 CAS #: 357-57-3 Brucine Other Names: 2,3-Dimethoxystrychnidin-10-one; 10,11-dimethoxystrychnine Haz Waste #: P018 Ind/EPA Gen #: C23H26N2O4 Formula: Description: Highly toxic alkaloid resembling strychnine; possible carcinogen PHYSICAL PROPERTIES: - - - - - - - - - - -- - - Ref. Molecular weight: 394.47 2 178.00 2 Melting point (C): Solubility in water (mg/L, 25C): 1 q/1320 mL 2 NONHALOG. PESTICIDES August 1987 CAS #: 88-85-7 Dinoseb

Other Names: 2,4-Dinitro-6-(1-methylpropyl)-phenol; DNBP; ENT 1122; 2-(1-Methylpropyl)-4,6-dinitrophenol; WSX 8365; Chemox PE; 2-sec-Butyl-4,6-dinitrophenol; Dow General; Premerge; Subitex; Caldon; Basanite Haz Waste #: P020 Ind/EPA Gen #:

Formula: C10H12N205 Description: Orange-brown viscous liquid

PHYSICAL PROPERTIES: - - - - - - -- - - - Ref. Molecular weight: 240.22 2 Melting point (C): 38.00 to 42.00 2

August 1987

NONHALOG. PESTICIDES Disulfoton

CAS #: 298-04-4

Solubility in water (mg/L, 25C):

Other Names: 0,0-Diethyl-S-[2-(ethylthio)ethyl]phosphoro dithioate Haz Waste #: P039 Ind/EPA Gen #: K036,K037 Formula: C8H1902PS3 Description: Oil PHYSICAL PROPERTIES: - - - - - - - -Molecular weight: 274.40 3 Melting point (C): > -25. 3 (1.5 torr) 132.00 to 133.00 Boiling point (C): 3 Density (g/cc, 20C): 1.1440 2 Vapor pressure (torr, 25C): (20 C) 0.18000E-03 3 Henry's law const (atm m**3/mol, 25C): 0.25000E-05 3 (23 C) 25.0000 3 Solubility in water (mg/L, 25C): August 1987 NONHALOG. PESTICIDES CAS #: 297-97-2 0,0-Diethyl P Other Names: 0,0-Diethyl-0-(2-pyrazinyl)phosphorthioate; Thionazin; Phosphorothioic acid 0,0-Diethyl 0-pyrazinyl ester; ethyl pyrazinyl phosphorothioate; EN-18133; ENT 25580; American Cyanamid 18133; Cynem; Nemafos; Zinophos Haz Waste #: PO40 Ind/EPA Gen #: C8H13N2O3PS Formula: Description: Amber liquid; cholinesterase inhibitor (Ref. 2) Molecular weight: 248.26 2 Melting point (C): -1.702Boiling point (C): 80.00 2 Vapor pressure (torr, 25C): (mm Hg at 30 C) 0.30000E-02 2 Solubility in water (mg/L, 25C): Slightly soluble 2 August 1987 NONHALOG. PESTICIDES CAS #: 311-45-5 Diethyl p-nitrophenyl phosphate Other Names: Phosphoric acid diethyl 4-nitrophenyl ester; Paraoxon; Phosphacol; E 600; Ester 25; Eticol; Fosfakol; Mintacol; Iostisal A; Soluglaucit Haz Waste #: PO41 Ind/EPA Gen #: C10H14N06P Formula: Description: Oily liquid; slight odor; poisonous; cholinesterase inhibitor. (Ref. 2) PHYSICAL PROPERTIES: - - - - -- - - Ref. Molecular weight: 275.21 2 Boiling point (C): 169.00 to 170.00 2 Density (g/cc, 20C): (25 C) 1.2683 2

2400. ug/mL 2

```
NONHALOG. PESTICIDES
August 1987
                                  Diisopropyl fluorophosphate
          55-91-4
CAS #:
Other Names: Isoflurophate; Phosphorofluoridic acid bis(1-methylethyl)
             ester: Isopropyl fluophosphate; Diisopropyl
             fluorophosphonate; Diisopropylphosphorofluoridate;
            Fluostigmine; Isofluorphate; DFP; Diflupyl; Dyflos;
             Floropryl; Fluropryl
Haz Waste #: PO43
                          Ind/EPA Gen #:
Formula:
            C6H14F03P
Description: Liquid; traces of vapor cause myosis; highly toxic;
             cholinesterase inactivator (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - -
Molecular weight:
                                                              184.15
Melting point (C):
                                                              -82.00
                                                                     2
                                                                     2
Boiling point (C):
                                            (by extrapolation) 183.00
                                                                     2
Density (g/cc, 20C):
                                                              1.0550
Vapor pressure (torr, 25C):
                                               (mm Hg at 20 C) 0.5790 2
Solubility in water (mg/L, 25C):
                       1.54% w/w (decomposes; pH about 2.5; forms HF) 2
                                  NONHALOG. PESTICIDES
August 1987
CAS #:
          60-51-5
                                  Dimethoate
Other Names: Phosphorodithioic acid 0,0-dimethyl S-[2-(methylamino)-2-
             oxoethyll ester; phosphorodithioic acid 0,0-dimethyl ester;
             ester with 2-mercapto-N-methylacetamide; 0,0-dimethyl S-
            methylcarbamoylmethyl phosphorodithioate; American
            Cyanamid 12880; Cygon; Fostion MM; Perfekthion; Rogor;
            Roxion
Haz Waste #: PO44
                          Ind/EPA Gen #:
Formula:
             C5H12N03PS2
Description: Crystals; cholinesterase inhibitor (Ref. 2)
Molecular weight:
                                                              229.28 2
```

Hydrolysis rate, water (1/month, 25C): Hydrolyzes in alkaline solution 2 Biodegradation rate in water (1/month): Very slightly soluble 2

Melting point (C):

RETENTION PROPERTIES: ------

log (octanol/water) partition (25C):

52.00 to

52.50 2

(65 C) 1.28 2

August 1987

Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

NONHALOG. PESTICIDES

CAS #: 39196-18-4 3,3-Dimethyl Other Names: 3,3-Dimethyl-1-(methylthio)-2-butanone-0-[(methylamino)carbonyl]oxime; Dacamox; Thiofanox; 3,3-Dimethyl-1-(methylthio)butanone 0-methylcarbamoyloxime Ind/EPA Gen #: Haz Waste #: PO45 C9H18N202S Formula: Description: Solid PHYSICAL PROPERTIES: - - - - - -- - - - Ref. Molecular weight: 218.00 13 Melting point (C): 56.50 to 57.50 13 NONHALOG. PESTICIDES August 1987 CAS #: 534-52-1 4,6-Dinitro-o-cresol Other Names: DNOC; Dinitrocresol; 2-Methyl-4,6-dinitrophenol; 3,5-Dinitro-2-hydroxytoluene; 3,5-Dinitro-o-cresol; DN; DNC; Antinonnin; Detal; Dinitrol; Elgetol; K III; K IV; Ditrosol; Prokarbol; Effusan; Lipan; Selinon; Sinox; Dekrysil Ind/EPA Gen #: Haz Waste #: PO47 Formula: C6H2(N02)2(OH)(CH3) Description: Yellow prisms; skin contact may lead to local necrosis and dangerous systemic effects; a cumulative poison (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. Molecular weight: 198.13 2 Melting point (C): 85.80 3 Henry's law const (atm m**3/mol, 25C): 0.14000E-05 3 Solubility in water (mg/L, 25C): (? C) 250.00 3 RETENTION PROPERTIES: - - - Ref. log (octanol/water) partition (25C): 2.85 3 7.4800 17 Partition coefficient, soil/water: Biodegradation rate in water (1/month): Not significant 3 NONHALOG. PESTICIDES August 1987 CAS #: 541-53-7 2.4-Dithiobiuret Other Names: Thioimidodicarbonic diamide Haz Waste #: PO49 Ind/EPA Gen #: Formula: NH2CSNHCSNH2 Description: Monoclinic or triclinic crystals; poisonous when ingested PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 135.20 2 Melting point (C): decomposes 181.00 2

(30 C) 1.5220 2

0.27 g/100 mL (27 C)

August 1987

NONHALOG. PESTICIDES

CAS #: 16752-77-5

Methomyl

Other Names: N-[[(Methylamino)carbonyl]oxy]ethan-imidothioic acid methyl ester; N-[(methylcarbamoyl)oxy]thio-acetimidic acid methyl ester; S-methyl N-[(methylcarbamoyl)oxy]thioacetimidate; methyl 0-(methylcarbamoyl)thiol-acetohydroxamate;

Insecticide 1179; Lannate; Nudrin

Haz Waste #: PO66 Ind/EPA Gen #:

Formula: C5H10N202S Description: Crystals

PHYSICAL PROPERTIES: - - - - - - -Molecular weight: 162.20 2 79.00 2 Melting point (C): 78.00 to Density (g/cc, 20C): (24 C) 1.2946 2

Solubility in water (mg/L, 25C):

5.8% 2

August 1987

NONHALOG. PESTICIDES

CAS #: 116-06-3 Aldicarb

Other Names: 2-Methyl-2-(methylthio)propanal O-[(methylamino)carbonyl]

oxime; 2-methyl-2-(methyltio)propionaldehyde 0-

(methylcarbamoyl)oxime; UC 21149; Temik

Haz Waste #: P070 Ind/EPA Gen #: CH3SC (CH3) 2CH=N0C0NHCH3 Formula:

Description: Crystals.

PHYSICAL PROPERTIES: - - - - - - - - - - -Molecular weight: 190.25 2 100.00 2 99.00 to Melting point (C): Solubility in water (mg/L, 25C): 0.6%w 2

August 1987 CAS #: 298-00-0 NONHALOG. PESTICIDES Methyl parathion

Other Names: 0,0-Dimethyl O-p-nitrophenyl phosphorothioate;

Phosphorothioic acid 0,0-dimethyl 0-(4-nitrophenyl) ester; 0,0-dimethyl 0-p-nitrophenyl thiophosphate; Dimethyl parathion; Parathion-methyl; Metaphos; E 601; ENT-17292;

Dalf; Folidol-M; Metacide; Metron; Nitrox 80

Haz Waste #: PO71 Ind/EPA Gen #:

(CH30) 2PS0-C6H4N02 Formula:

Description: Crystals; cholinesterase inhibitor (Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - -Molecular weight: 263.23 2 Melting point (C): 36.00 3 35.00 to Boiling point (C): Unstable 3 Density (g/cc, 20C): 1.3580 2 Vapor pressure (torr, 25C): (20 C) 0.97000E-05 Henry's lawwconst (atm m**3/mol, 25C): 0.54000E-07 3 Solubility in water (mg/L, 25C): 55.0000 to 60.0000 3

NONHALOG. PESTICIDES August 1987 CAS #: 86-88-4 1-Naphthy I-2-thiourea Other Names: BANTU; 1-Naphthalenylthiourea; ANTU; Krysid; Chemical 109; 1-(1-Naphthyl)-2-thiourea; alpha-Naphthylthiourea; Anturat; N-1-Naphthylthiourea; alpha-Naphthylthiocarbamide; Rattrack Haz Waste #: P072 Ind/EPA Gen #: Formula: C11H10N2S Description: Prisms; toxic; bitter taste PHYSICAL PROPERTIES: -- - Ref. Molecular weight: 202.27 2 198.00 2 Melting point (C): Solubility in water (mg/L, 25C): 0.06 g/100 mL 2 NONHALOG. PESTICIDES August 1987 CAS #: 54-11-5 Nicotine Other Names: 3-(1-Methyl-2-pyrrolidinyl)pyridine; 1-Methyl-2-(3-pyridyl)pyrrolidine; beta-Pyridyl-alpha-N-methyl-pyrrolidine Haz Waste #: P075 Ind/EPA Gen #: Formula: C10H14N2 Description: Colorless to pale yellow, oily liquid; very hygroscopic; turns brown on exposure to air or light; acrid, burning taste; develops odor of pyridine; highly toxic; base is absorbed through mucous membranes and intact skin, but the salts (hydrochloride, dihydrochloride, sulfate, tartarate, zinc chloride, salicylate) are not (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - -- - Ref. 162.23 2 Molecular weight: Boiling point (C): (745 torr), partial decomposition 247.00 2 1.0097 2 Density (g/cc, 20C): Miscible 2 Solubility in water (mg/L, 25C): NONHALOG. PESTICIDES August 1987

CAS #: 152-16-9 Octamethy I pyrophosphorami de

Other Names: Schradan; Octamethyl pyrophosphoramide; Pestox III; bis[bisdimethylaminophosphonous] anhydride; OMPA; Sytam; bis-N,N,N',N'-tetramethylphosphorodiamidic anhydride

Ind/EPA Gen #: Haz Waste #: PO85

Formula: C8H24N4O3P2 Description: Viscous liquid

PHYSICAL PROPERTIES: - - - - - - - - - - -Molecular weight: 286.26 2 2 Melting point (C): 14.00 to 20.00 Boiling point (C): 120.-125. (0.5 atm); 154. (2.0 atm) Density (g/cc, 20C): (25 C) 1.0900 2 Vapor pressure (torr, 25C): (mm Hg) 0.10000E-02 2 Solubility in water (mg/L, 25C): Miscible 2

NONHALOG. PESTICIDES August 1987 CAS #: 145-73-3 7-0xabicyclo Other Names: ENDOTHALL; 7-0xabicyclo[2.2.1]heptane-2,3-dicarboxylic acid Ind/EPA Gen #: Haz Waste #: P088 Formula: C8H1005 PHYSICAL PROPERTIES: - -- - - Ref. 186.16 2 Molecular weight: 10 q/100 mL (20 C) 2 Solubility in water (mg/L, 25C): NONHALOG. PESTICIDES August 1987 CAS #: 56-38-2 Parathion Other Names: 0.0-Diethyl O-p-nitrophenyl phosphorothioate; DNTP; S.N.P.; Phosphorothioic acid 0,0-diethyl 0-(4-nitrophenyl) ester; Diethyl-p-nitrophenyl monothiophosphate; E 605; AC 3422; ENT 15108; Alkron; Alleron; Aphamite; Etilon; Folidol; Fosferno; Niran; Paraphos; Rhodiatox; Thiophos Ind/EPA Gen #: Haz Waste #: P089 (C2H50)2PS0C6H4N02; C10H14N05PS Formula: Description: Pale, yellow liquid; highly toxic; cholinesterase inhibitor effects are cumulative; special precautions necessary to prevent inhalation and skin contamination (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - -- Ref. 291.30 3 Molecular weight: Melting point (C): 3 6.10 Boiling point (C): 375.00 Density (g/cc, 20C): (25 C) 1.2600 2 Vapor pressure (torr, 25C): (20 C) 0.37800E-04 3 Henry's law'const (atm m**3/mol, 25C): 0.61000E-06 24.0000 3 Solubility in water (mg/L, 25C): August 1987 NONHALOG. PESTICIDES CAS #: 103-85-5 N-Phenylthiourea

Other Names: Phenylthiourea; Phenylthiocarbamide Haz Waste #: P093 Ind/EPA Gen #: C6H5NHCSNH2: C7H8N2S

Description: Bitter or tasteless needles PHYSICAL PROPERTIES: - - - - - - - - -

Solubility in water (mg/L, 25C):

Formula:

Molecular weight:

Melting point (C):

Density (g/cc, 20C):

- - - Ref.

2

152.22

154.00

1 part/400 (? C) 2

1.3000 2

August 1987
CAS #: 298-02-2
Other Names: Phosphorodithioic acid 0,0-diethyl S-[(ethylthio)methyl]

ester Haz Waste #: P094 Ind/EPA Gen #: K038,K039,K040 (C2H50) 2PS-SCH2SC2H5 Formula: Description: Liquid - - - Ref. PHYSICAL PROPERTIES: - -Molecular weight: 260.40 2 (25°C) 1.1560 2 Density (g/cc, 20C): Vapor pressure (torr, 25C): (mm Hg) 0.84000E-03 2 (ppm) 50.0000 2 Solubility in water (mg/L, 25C): RETENTION PROPERTIES: - - - Ref. Hydrolysis rate, water (1/month, 25C): Hydrolyzes 2

August 1987 CAS #: 52-85-7 NONHALOG. PESTICIDES Famphur

Other Names: Phosphorothioic acid 0-[4-[(dimethylamino)sulfonyl]phenyl]
0,0-dimethyl ester; Phosphorothioic acid 0,0-dimethyl ester;

0-Ester with p-hydroxy-N, N-dimethy | benzenesul fonamide;

0,0-Dimethyl-0,p-(dimethylsulfamoyl)phenyl phosphorothioate; p-Hydroxy-N,N-dimethylbenzenesulfonamide ester with

phosphorothioic acid 0,0-dimethyl ester; Famophos; American Cyanamid 38023; ENT 25644; Warbex

Haz Waste #: P097 Ind/EPA Gen #:

Formula: (CH30) 2PS0C6H4S02N(CH3) 2

Description: Crystals; cholinesterase inhibitor (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - Ref.

Molecular weight: 325.36 2 Melting point (C): 52.50 to 53.50 2

```
August 1987
                                  NONHALOG. PESTICIDES
CAS #:
         57-24-9
                                  Strychnine and salts
Other Names: Strychnidin-10-one
                          Ind/EPA Gen #:
Haz Waste #: P108
Formula:
            C21H22N202
Description: Very bitter, orthorhombic, spehnoidal prisms; salts are
            acetate (crystals); arsenate dihydrate (efflorescent
            crystals); dichromate (orange-yellow needles); formate
             (crystals); gluconate pentahydrate (crystals);
            glycerophosphate hexahydrate; hydrobromide monohydrate
             (efflorescent crystans); hydrochloride dihydrate
             (efflorescent, trimetric prisms); nitrate; phosphate;
            salicylate (leaflets); valerate (crystals); extremely
            poisonous (Ref. 2)
- - - Ref.
Molecular weight:
                                                              334.40
                                                                     3
Melting point (C):
                                                268.00 to
                                                              290.00
                                                                     3
                                                     (5 torr) 270.00 3
Boiling point (C):
Density (g/cc, 20C):
                                                              1.3600 2
                                                        (? C) 156.00 3
Solubility in water (mg/L, 25C):
August 1987
                                  NONHALOG. PESTICIDES
CAS #: 3689-24-5
                                  Tetraethyldithiopyrophosphate
Other Names: Dithiopyrophosphoric acid tetraethyl ester;
            Tetraethyl thiodiphosphate; Tetraethyl dithiopyrophosphate;
            Sulfotep
                          Ind/EPA Gen #:
Haz Waste #: P109
Formula:
            C8H2O05P2S2
Description: Oil; corrodes iron; highly toxic (Ref. 13)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -
Molecular weight:
                                                              322.00 13
Boiling point (C):
                                                110.00 to
                                                             113.00 13
Solubility in water (mg/L, 25C):
                                         Slightly soluble (670 ppm) 13
                                  NONHALOG. PESTICIDES
August 1987
CAS #: 107-49-3
                                  Tetraethy | pyrophosphate
Other Names: Diphosphoric acid tetraethyl ester
Haz Waste #: P111
                         Ind/EPA Gen #:
             (C2H50) 2P000P (OC2H5) 2
Formula:
Description: Liquid
PHYSICAL PROPERTIES: - - - - - - - - - -
                                                                 – – Ref.
Molecular weight:
                                                              290.20 2
                                         decomposes 170.00 to 213.00 2
Melting point (C):
```

Density (g/cc, 20C):

RETENTION PROPERTIES:

Vapor pressure (torr, 25C):

Solubility in water (mg/L, 25C):

Hydrolysis rate, water (1/month, 25C):

Hydrolyzes quickly, t1/2 about 7 hrs (25 C) 2

1.1850

Miscible 2

0.00047 mm Hg (30 C)

2

2

August 1987 CAS #: 137-26-8

NONHALOG. PESTICIDES Thiram

Other Names: Tetramethylthioperoxydicarbonic diamide; TMTD; ENT 987; Bis(dimethylthiocarbamoyl) disulfide; SQ 1489; NSC 1771; Bis(dimethylthiocarbamyl) disulfide; Thiurad; Thiosan; Tetramethylthiuram disulfide; Thylate; Tiuramyl; Thiuramyl; Puralin; Fernasan; Nomersan; Rezifilm; Pomarsol; Tersan; Tuads: Tulisan: Arasan Haz Waste #: P117,U244 Ind/EPA Gen #: C6H12N2S4 Formula: Description: Crystals; irritant of mucous membranes and skin (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - -- - - - Ref. Molecular weight: 240.44 Melting point (C): (commercial grades, 146.) 155.00 to 156.00 2 Density (g/cc, 20C): 1.2900 2 Solubility in water (mg/L, 25C): Insoluble 2

August 1987 CAS #: 50-07-7 NONHALOG. PESTICIDES

Mitomycin C

Other Names: [1aR]-6-amino-8-[[(aminocarbonyl)oxy]methyl]-1,1a,2,8,8a,8b-

hexahydro-8a-methoxy-5-methylazirino[2',3':3,4]pyrrolo[1, 2-a]indole-4,7-dione, Ametycine; MMC; Mitocin-C; Mutamycin

Haz Waste #: U010 Ind/EPA Gen #:

Formula: C15H1805N4

Description: Blue-violet crystals; used in cancer chemotherapy (Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Melting point (C): > 360. 2

Solubility in water (mg/L, 25C): Soluble 2

```
NONHALOG. PESTICIDES
August 1987
          61-82-5
CAS #:
                                   Amitrole
Other Names: 3-Amino-1H-1,2,4-triazole; Aminotriazole; ATA; ENT 25445;
             Amizol; Cytrol; Weedazol; Amerol; 3-Amino-S-triazole;
             Amino triazole weedkiller 90; 3-Amino-1,2,4-triazole;
             5-Amino-1,2,4-triazole; Amitrol; Amitril; AT; ATA; Azaplant;
             Azole; Diurol; Domatol; Elmasil; Emisol; Fenamine;
             Herbizole; Kleer-lot; Orga-414; Ramizol; Simazol;
             USAF XR-22; Vorox; Weedazin; Weedoclor; (other commercial
             names)
Haz Waste #: U011
                           Ind/EPA Gen #:
             C2H4N4
Formula:
Description: Crystals; listed as a carcinogen by the EPA (Ref. 2);
             will persist in soil for several weeks and in water for
             more than 200 days after application; canceled for use on
             croplands July 1971 (Ref. 12)
                                                                --- - Ref.
PHYSICAL PROPERTIES: - - - - - - - -
Molecular weight:
                                                                  84.08 2
Melting point (C):
                                                                159.00 2
Solubility in water (mg/L, 25C):
                                                                Soluble 2
                                   NONHALOG. PESTICIDES
August 1987
CAS #:
         492-80-8
                                   Auramine
Other Names: 4,4'-(Imidocarbonyl)bis(N,N-dimethylaniline); Auramine base;
             4,4'-(Iminocarbonyl)bis(N,N-dimethylaniline); Auramine OAF;
             Apyonine auramine base; Auremine; Brilliant oil yellow;
             Bis (P-dimethylaminophenyl) methyleneimine; Glauramine;
             C.I. Solvent yelloww34; C.I. 41000B; Waxoline yellow 0;
             4,4'-Dimethylaminobenzophenonimide; Yellow pyoctanine;
             Tetramehtyldiaminodiphenylacetimine; Basic yellow 2
             4,4'-CarbonimidoyIbis(N,N-dimethyI)benzenamine
Haz Waste #: U014
                           Ind/EPA Gen #:
Formula:
             C17H21N3
```

Description: Yellow leaves or needles; used industrially as a dye or

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -

Molecular weight:

Solubility in water (mg/L, 25C):

dye intermediate; vapors harmful; a carcinogen (Ref. 12)

- - - Ref.

267.37 17

Insoluble 12

August 1987 NONHALOG. PESTICIDES Cyclophosphamide

Other Names: N,N-Bis(2-chloroethyl) tetrahydro-2H-1,3,2-oxazaphosphorin-2-amine 2-oxide; 2-[Bis(2-chloroethyl)amino]tetrahydro-2H-1,3,2-oxazophosphorine 2-oxide; 1-Bis(2-chloroethyl)amino-1-oxo-2-aza-5-oxaphosphoridin; Cyclophosphane; B518; Bis(2-chloroethyl)phosphamide cyclic propanolamide ester; Bis(2-chloroethyl)phosphoramide cyclic propanolamide ester; N, N-Bis (beta-chloroethyl)-N', O-propylenephosphoric acid ester diamide; N.N-Bis(beta-chloroethyl)-N', 0trimethylenephosphoric acid ester diamide; Cytophosphane; Cytoxan; Endoxan; Procytox; Sendoxan; ASTA; (many other) Haz Waste #: U058 Ind/EPA Gen #: C7H15C12N202P Formula: Description: Monohydrate; listed as a carcinogen by the EPA (Ref. 2); a widely used synthetic druge used in cancer chemotherapy and chemically related to the nitrogen mustards (Ref. 12) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 261.10 2 41.00 to 45.00 2 Melting point (C): Solubility in water (mg/L, 25C): 0.40000E+05 2 NONHALOG. PESTICIDES August 1987 CAS #: 3288-58-2 0,0-Diethyl Other Names: Phosphorodithioic acid, 0.0-diethyl S-methyl ester Haz Waste #: U087 Ind/EPA Gen #: C5H1302PS2 Formula: PHYSICAL PROPERTIES: - - - ---- Ref. Molecular weight: 200.25 17 NONHALOG. PESTICIDES August 1987 CAS #: 111-54-6 Ethylene bis-dithiocarbamate Other Names: Ethylene bis(dithiocarbamic acid); Ethylenebisdithiocarbamic acid; 1,2-Ethanediylbiscarbamodithioic acid Haz Waste #: U114 Ind/EPA Gen #: Formula: C4H8N2S4 Description: Parent acid very unstable; decomposes to CS2 and ethylenethiourea (Ref. 13) PHYSICAL PROPERTIES: - - - - - - - -- - - Ref. Molecular weight: 212.00 13

August 1987 CAS #:

NONHALOG. PESTICIDES Cacodylic acid

75-60-5

Other Names: Hydroxydimethylarsine oxide; Dimethylarsinic acid; Phytar

Ind/EPA Gen #: KO31 Haz Waste #: U136

(CH3)2As(0)0H; C2H2As02 Formula:

Description: Crystals; hygroscopic; poisonous; herbicide

PHYSICAL PROPERTIES: - - - - - - - - - - -- - Ref.

Molecular weight: 137.99 2

Melting point (C): 195.00 to 196.00 2

Solubility in water (mg/L, 25C): Soluble in 0.5 part water 2 7.11 Organo-Nitrogens

August 1987 CAS #: 122-09-8

ORGANO-NITROGENS

P-alpha-Dimethylphenethylamine

Other Names: alpha, alpha-Dimethylphenethylamine; Phentermine; alpha, alpha-Diemthy Ibenzeneethanamine; Phenyl-tertbutylamine; alpha-benzylisopropylamine Haz Waste #: PO46 Ind/EPA Gen #: CH3C6H4CH2(NH2)CHCH3 Formula: Description: Oily liquid PHYSICAL PROPERTIES: - - - - - - - - - - - ----- Ref. 149.23 2 Molecular weight: Boiling point (C): (10) 103.00 to 103.00 2

August 1987 ORGANO-NITROGENS 62-75-9 N-Nitrosodimethylamine CAS #:

Other Names: Dimethylnitrosoamine; N-Methyl-N-nitrosomethanamine; DMN; DMNA; Dimethylnitrosamine; N,N-Dimethylnitrosamine; NDMA;

N-Nitroso-N, N-dimethylamine Haz Waste #: P082,U100 Ind/EPA Gen #:

(CH3) 2NN=0; C2H6N20 Formula:

Description: Yellow liquid; potent carcinogen; listed as a carcinogen by

the EPA (Ref. 2); oily liquid (Ref. 12)

PHYSICAL PROPERTIES: - - - - - - - - - - - - - -- - - Ref. 74.08 3 Molecular weight: 153.00 3 Boiling point (C): 151.00 to Density (g/cc, 20C): 1.0048 2 Solubility in water (mg/L, 25C): Miscible 3 RETENTION PROPERTIES: - - - Ref. log (octanol/water) partition (25C): 0.06 3 Partition coefficient, soil/water: 0.1900 17 Biodegradation rate in water (1/month): Significant 3

August 1987

ORGANO-NITROGENS

CAS #: 86-30-6 N-nitroso-diphenylamine

Other Names: DiphenyInitrosoamine

Haz Waste #: PO83 Ind/EPA Gen #:

Formula: (C6H5)2N=0

Description: Note (4/7/87): These values may be for p-nitroso

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 198.22 2 145.00 2 Melting point (C): 144.00 to Solubility in water (mg/L, 25C):

RETENTION PROPERTIES: -----Ref. log (octanol/water) partition (25C): 2.57 3

Biodegradation rate in water (1/month):

Significant 3

August 1987

log (octanol/water) partition (25C):

Biodegradation rate in water (1/month):

Partition coefficient, soil/water:

ORGANO-NITROGENS

CAS #: 4549-40-0 N-Nitroso-methylvinylamine Other Names: N-Methyl-N-nitrosoethenamine: MVNA: NMVA: N-Methyl-N-nitroso-ethenylamine Haz Waste #: P084 Ind/EPA Gen #: Formula: C3H6N20 Description: Pale yellow liquid; very volatine; photolabile; a carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - -- - - - Ref. 86.09 17 Molecular weight: Solubility in water (mg/L, 25C): Soluble 12 ORGANO-NITROGENS August 1987 CAS #: 124-40-3 Dimethylamine Other Names: N-Methylmethanamine Haz Waste #: U092 Ind/EPA Gen #: Formula: (CH3) 2NH PHYSICAL PROPERTIES: - -- - - Ref. Molecular weight: 45.08 2 Melting point (C): -96.00 2 Boiling point (C): 7.00 2 Solubility in water (mg/L, 25C): Very soluble 2 August 1987 ORGANO-NITROGENS CAS #: 121-14-2 2,4-Dinitrotoluene Other Names: DNT; 1-methyl-2,4-dinitrotoluene Haz Waste #: U105 Ind/EPA Gen #: Formula: C6H3 (NO2) 2CH3 PHYSICAL PROPERTIES: - - - - - -- - - Ref. Molecular weight: 182.10 3 Melting point (C): 70.00 3 Boiling point (C): 300.00 3 Vapor pressure (torr, 25C): (59 C) 0.13000E-02 3 Henry's law const (atm m**3/mol, 25C): 0.46800E-04 3 Solubility in water (mg/L, 25C): (22 C) 270.00 3 RETENTION PROPERTIES: -------

2.01 3

10.9000 17

Significant 3

ORGANO-NITROGENS August 1987 606-20-2 CAS #: 2,6-Dinitrotoluene Other Names: Dinitrotoluol Haz Waste #: U106 Ind/EPA Gen #: Formula: C6H3 (N02) 2CH3 PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - Ref. 182.10 3 Molecular weight: Melting point (C): 63.00 3 285.00 3 Boiling point (C): Vapor pressure (torr, 25C): (20 C) 0.35000E-03 3 Solubility in water (mg/L, 25C): RETENTION PROPERTIES: ------(20 C) 0.10000E-02 3 - - - Ref. log (octanol/water) partition (25C): 2.05 3 Significant 3 Biodegradation rate in water (1/month): ORGANO-NITROGENS August 1987 CAS #: 142-84-7 N-Dipropylamine Other Names: N-Propylpropanamine; Dipropylamine Haz Waste #: U110 Ind/EPA Gen #: Formula: (CH3CH2CH2) 2NH Description: Liquid PHYSICAL PROPERTIES: - - - - - - -– – – Ref. 101.19 2 Molecular weight: Melting point (C): -63.00 2 110.00 2 Boiling point (C): Density (g/cc, 20C): 0.7380 2 Solubility in water (mg/L, 25C): Freely soluble 2 August 1987 ORGANO-NITROGENS CAS #: 621-64-7 di-n-PropyInitrosamine Other Names: N-nitroso-di-n-propylamine; Dipropylnitrosamine; N, N-DipropyInitrosamine; DPN; DPNA; NDPA; Nitrosodipropylamine; N-Nitrosodipropylamine; N-Nitroso-N-propyl-1-propanamine Haz Waste #: U111 Ind/EPA Gen #: Formula: (C3H7)2N=0

Description: Yellow liquid; a carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - - - - - - - -

RETENTION PROPERTIES: --------

Solubility in water (mg/L, 25C):

Partition coefficient, soil/water:

log (octanol/water) partition (25C):

Biodegradation rate in water (1/month):

Molecular weight:

Boiling point (C):

---- Ref.

130.20 3 205.00 3

9900.00 3

- - - Ref.

1.31 3

2.5400 17

Not significant 3

ORGANO-NITROGENS August 1987 CAS #: 134-32-7 1-Naphthylamine Other Names: 1-Naphthalenamine Haz Waste #: U167 Ind/EPA Gen #: Formula: C10H8NH2 PHYSICAL PROPERTIES: - - - - - - -Molecular weight: 143.18 2 50.00 2 Melting point (C): 301.00 2 Boiling point (C): Solubility in water (mg/L, 25C): 1 part/590 (7 C) 2 ORGANO-NITROGENS August 1987 CAS #: 91-59-8 2-Naphthylamine Other Names: 2-Aminonaphthalene; C.I. 37270; Fast scarlet base B; NA; 2-Naphthalamine; 2-Naphthalenamine; beta-Naphthylamine; 6-Naphthylamine; 2-Naphthylamine mustard; USAF CB-22 Haz Waste #: U168 Ind/EPA Gen #: C10H7NH2; C10H9N Formula: Description: White to reddish crystals; volatile with steam; listed as a carcinogen by the EPA (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - -- - - Ref. Molecular weight: 143.18 2 111.00 to 113.00 Melting point (C): Boiling point (C): 306.00 2 Density (g/cc, 20C): (98 C) 1.0610 2 Soluble in hot water Solubility in water (mg/L, 25C): August 1987 ORGANO-NITROGENS CAS #: 98-95-3 Nitrobenzene Other Names: Oil of Mirbane; Nitrobenzol; Essence of mirbane Haz Waste #: U169 Ind/EPA Gen #: F004,K025,K104 Formula: C6H5N02 Description: Colorless to pale yellow, oily liquid; odor of volatile oil almond; poisonous; rapidly absorbed through skin; vapor hazardous (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - -Molecular weight: 123.11 2 Melting point (C): 5.60 3 Boiling point (C): 211.00 3 Density (g/cc, 20C): (15 C) 1.2050 2 Vapor pressure (torr, 25C): (20 C) 0.1500 3 Henry's law const (atm m**3/mol, 25C): 0.24000E-04 3 (20 C) 1900.00 3 -----Ref. log (octanol/water) partition (25C): 1.85 3 Partition coefficient, soil/water: 7.8200 17 Significant 3 Biodegradation rate in water (1/month):

August 1987 ORGANO-NITROGENS CAS #: 924-16-3 N-Nitroso-di-n-butylamine Other Names: Azo-stat; Azodyne; N-Butyl-N-nitroso-1-butamine; DBN; DBNA; DibutyInitrosamine; Di-n-butyInitrosamine; N, N-Di-n-butyInitrosamine; N, N-DibutyInitrosoamine; NDBA; Nitrosodibutylamine; N-Nitrosodibutylamine; N-Butyl-N-nitroso-1-butanamine Ind/EPA Gen #: Haz Waste #: U172 Formula: C8H18N20 Description: Yellow oil; a carcinogen (Ref. 12) PHYSICAL PROPERTIES: - -158.24 17 Molecular weight: Soluble 12 Solubility in water (mg/L, 25C): ORGANO-NITROGENS August 1987 CAS #: 1116-54-7 N-Nitrosoethanolamine Other Names: 2,2'-(Nitrosimino)bisethanol; Diethanolnitrosamine; Bis(beta-hydroxyethyl)nitrosamine; N,N-Diethanolnitrosamine; 2,2'-Dihydroxy-N-nitroso-diethylamine; NCI-C5583; NDEA; NDELA; N-Nitrosoaminodiethanol; Nitrosoimino diethanol; N-Nitrosobis (2-hydroxyethyl) amine; Nitrosodiethanolamine; 2,2'-Nitrosiminodiethanol; Di-(2-ydroxyethyl)nitrosamine Haz Waste #: U173 Ind/EPA Gen #: (HOCH2CH2) 2NNO; C4H1ON2O3 Formula: Description: Light yellow, viscous oil; listed as a carcinogen by the EPA (Ref. 2) PHYSICAL PROPERTIES: - - -- - - Ref. Molecular weight: 134.13 2 (.01 mm Hg) 125.00 2 Boiling point (C): Solubility in water (mg/L, 25C): Miscible 12 ORGANO-NITROGENS August 1987 CAS #: 55-18-5 N-Nitrosodiethylamine Other Names: N-Ethyl-N-nitrosoethanamine; DANA; DEN; DENA; NDEA; Diethylnitrosamide; Diethylnitrosamine; N, N-DIethylnitrosamine; Diethylnitrosoamine; Nitrosodiethylamine; N-Nitroso-N, N-diethylamine Ind/EPA Gen #: Haz Waste #: U174 (C2H5) 2NNO; C4H10N2O Formula: Description: Slightly yellow liquid; listed as a carcinogen by the EPA

(Ref. 2); decomposes in the presence of sunlight; volatile

102.14

177.00

0.9422 2

Soluble 2

2

175.00 to

(Ref. 12)

Solubility in water (mg/L, 25C):

PHYSICAL PROPERTIES: - -

Molecular weight:

Boiling point (C):

Density (g/cc, 20C):

ORGANO-NITROGENS August 1987 CAS #: 759-73-9 N-Nitroso-N-ethylurea Other Names: N-Ethyl-N-nitroso-carbamide; ENU; N-Ethyl-N-nitroso-urea; N-Ethyl-N-nitrosourea; 1-Ethyl-1-Nitrosourea; NEU; Ethylnitrosourea; N-Ethylnitrosourea; Nitrosoethylurea; NSC 45403 Haz Waste #: U176 Ind/EPA Gen #: C3H7N302 Formula: Description: Crystal; a carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - - - - - - -- - - Ref. 117.11 17 Molecular weight: Solubility in water (mg/L, 25C): Soluble 12 ORGANO-NITROGENS August 1987 CAS #: 684-93-5 N-Nitroso-N-methylurea Other Names: N-Methyl-N-nitrosocarbamide; Methylnitrosourea; N-Methyl-N-nitrosourea; 1-Methyl-1-nitrosourea; MNU; Nitrosomethylurea; 1-Nitroso-1-methylurea; NMH; NMU; NSC 23909 Haz Waste #: U177 Ind/EPA Gen #: Formula: C2H5N302 Description: A carcinogen; considered for cancer chemotherapy (Ref. 12) - - - - Ref. Molecular weight: 103.08 17 Solubility in water (mg/L, 25C): Soluble 12

August 1987

CAS #: 615-53-2

ORGANO-NITROGENS
N-Nitroso-N-methylurethane

Other Names: Methylnitroso carbamic acid, ethyl ester; N-Methyl-N-nitrosoethylcarbamate; Methylnitrosourethan;

Ethyl ester of methylnitroso-carbamic acid;

N-Methyl-N-nitroso-urethane; Nitrosomethylurethan

Haz Waste #: U178 Ind/EPA Gen #: Formula: C4H8N2O3

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Molecular weight: 132.14 14

August 1987 CAS #: 100-75-4 ORGANO-NITROGENS
N-Nitrosopiperidine

Other Names: 1-Nitrosopiperidine; NO-PIP; NPIP

Haz Waste #: U179 Ind/EPA Gen #:

Formula: C5H10N20

Description: Rapidly decomposed by sunlight; a carcinogen (Ref. 12)

PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.

Molecular weight:

114.15 17

August 1987

ORGANO-NITROGENS

CAS #: 930-55-2

N-Nitrosopyrrolidine

Other Names: 1-Nitrosopyrrolidine; NO-PYR; NPYR;

N-Nitroso-1-pyrrolidinamine

Haz Waste #: U180

Ind/EPA Gen #:

Formula: C4H8NN0

Description: Yellow liquid; listed as a carcinogen by the EPA (Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Molecular weight:

100.11 2

Boiling point (C):

(2) 104.00 to 106.00 2

Solubility in water (mg/L, 25C):

Soluble 2

August 1987

ORGANO-NITROGENS

CAS #: 99-55-8

5-Nitrotoluidine

Other Names: 2-Methyl-5-nitrobenzenamine; 2-Methyl-5-nitroaniline;

5-Nitro-o-toluidine

Haz Waste #: U181 Ind/EPA Gen #:

Formula:

C7H8N202

Description: Yellow prisms

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Molecular weight:

152.00 13 107.00 13

256.13

2

Melting point (C):

August 1987

7 ORGANO-NITROGENS

CAS #: 23950-58-5

Pronamide

Other Names: 3,5-Dichloro-N-(1,1-dimethyl-2-propynyl)benzamide; Propamid;

RH 315; Kerb

Haz Waste #: U192 Ind/EPA Gen #:

Formula: C12H11Cl2N0

Description: White solid; herbicide

PHYSICAL PROPERTIES: - - - - - - - - - - - - Ref.

Molecular weight:

Melting point (C): 155.00 to 156.00 2

Solubility in water (mg/L, 25C): 15 ppm 2

August 1987 CAS #: 107-10-8 ORGAND-NITROGENS n-Propylamine

Other Names: 1-Propanamine; Propylamine; 1-Aminopropane

Haz Waste #: U194 Ind/EPA Gen #:

Formula: CH3CH2CH2NH2; C3H9N

Description: Colorless, alkaline liquid; strong ammonia odor; strong

irritant; possible skin sensitizer (Ref. 2)

Density (g/cc, 20C): 0.7190 2 Solubility in water (mg/L, 25C): Miscible 2

147

7.12 Organo-Sulfurs

ORGANO-SULFURS August 1987 CAS #: 591-08-2 1-Acetyl-2-thiourea Other Names: N-(Aminothioxomethyl)acetamide Haz Waste #: P002 Ind/EPA Gen #: Formula: C3H6N20S Description: Needles; severe poison PHYSICAL PROPERTIES: - - - - - - - - - - - -Melting point (C): 165.00 13 ORGANO-SULFURS August 1987 CAS #: 108-98-5 Thiophenol Other Names: Benzenethiol; Phenylmercaptan Haz Waste #: P014 Ind/EPA Gen #: Formula: C6-H5SH Description: Liquid; repulsive, penetrating, garlic-like odor; oxidizes in air (Ref. 2) PHYSICAL PROPERTIES: -------110.17 2 Molecular weight: Boiling point (C): 168.30 2 Density (g/cc, 20C): (25 C) 1.0728 2 Solubility in water (mg/L, 25C): Insoluble 2 ORGANO-SULFURS August 1987 CAS #: 96-45-7 Ethylene thioruea Other Names: 2-Imidazolidinethione Haz Waste #: U116 Ind/EPA Gen #: Formula: C3H6N2S Description: Needles, prisms PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref. Molecular weight: 102.17 2 Melting point (C): 203.00 to 204.00 2 2 g/100 mL at 30 C 2 Solubility in water (mg/L, 25C): ORGANO-SULFURS August 1987 CAS #: 62-50-0 Ethyl methanesulfonate Other Names: Methanesulfonic acid ethyl ester; Ethyl methanesulfonic acid; ethyl mesylate; EMS; NSC 26805 Haz Waste #: U119 Ind/EPA Gen #: CH3S020CH2CH3 Description: Liquid; used experimentally as a mutagen PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. Molecular weight: 124.15 2

Boiling point (C): Density (g/cc, 20C): 213.00 to 213.50 2

(22 C) 1.1452 2

August 1987 ORGANO-SULFURS CAS #: 74-93-1 Methyl mercaptan Other Names: Methanethiol; Mercaptomethane; Thiomethyl alcohol; Methyl sulfhydrate Haz Waste #: U153 Ind/EPA Gen #: CH3SH: CH4S Formula: Description: Flammable gas; odor of rotten cabbage (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - -– – - Ref. Molecular weight: 48.11 2 Melting point (C): -123.00 3 Boiling point (C): 5.95 0.8665 Density (g/cc, 20C): 2 Vapor pressure (torr, 25C): 1500.00 3 Henry's law!const (atm m**3/mol, 25C): 0.38500E-02 3 Solubility in water (mg/L, 25C): (20 C) 0.23330E+05 August 1987 ORGANO-SULFURS CAS #: 1120-71-4 1,3-Propane sulfone Other Names: 2,2-Dioxide-1,2-oxathiolane Haz Waste #: U193 Ind/EPA Gen #: Formula: C3H603S PHYSICAL PROPERTIES: - - - - - - - - -- - - Ref. Molecular weight: 122.14 17 August 1987 ORGANO-SULFURS CAS #: 62-55-5 Thioacetamide

August 1987 ORGANO-SULFURS
CAS #: 62-55-5 Thioacetamide

Other Names: Ethanethioamide
Haz Waste #: U218 Ind/EPA Gen #:
Formula: CH3CSNH2; C2H5NS

Description: Crystals; slight odor of mercaptans
PHYSICAL PROPERTIES: - - - - - - - - Ref.
Molecular weight: 75.14 2
Melting point (C): 113.00 to 114.00 2

Solubility in water (mg/L, 25C):

16.3 g/100 mL 2

August 1987

ORGANO-SULFURS

CAS #: 62-56-6 Thiourea

Other Names: Thiocarbamide; Thiocarbonic diamide Haz Waste #: U219 Ind/EPA Gen #:

Formula: H2NCSNH2; CH4N2S

Description: Rhombohedron or needles; experimental carcinogen (Ref. 13)

PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - Ref.

Molecular weight: 76.00 13

Melting point (C): 180.00 13

Solubility in water (mg/L, 25C): Soluble 13 7.13 Halogenated N.O.S.

August 1987 CAS #: 598-31-2 HALOGENATED N.O.S.

Bromoacetone

Other Names: Bromo-2-propanone

Ind/EPA Gen #: Haz Waste #: P017

Formula: CH3COCH2Br

Description: Liquid; violent lacrimator

PHYSICAL PROPERTIES: - - - - - - - -136.99 2 Molecular weight: -36.50 2 Melting point (C): 137.00 2 Boiling point (C): (23 C) 1.6340 2 Density (g/cc, 20C): Vapor pressure (torr, 25C): 0.00000E+00 2 0.00000E+00 2 Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): Sparingly soluble 2

August 1987

HALOGENATED N.O.S.

CAS #: 107-20-0 Chloroacetaldehyde

Other Names: 2-Chloro-1-ethanal; monochloroacetaldehyde

Ind/EPA Gen #: Haz Waste #: P023

Formula: CICH2CH0

Description: Liquid; acrid, penetrating odor; very corrosive to mucous

membranes; intensely irritating to eyes, skin.

- - - Ref.

Molecular weight: 78.50 2 Boiling point (C): (760 mm Hg) 85.00 to 6.00 2

August 1987

HALOGENATED N.O.S.

CAS #: 106-47-8 p-Chloroaniline

Other Names: 4-Chlorobenzenamine

Haz Waste #: P024 Ind/EPA Gen #:

C6H6CIN Formula:

Description: Orthorhombic crystals

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.

127.57 2 Molecular weight:

Melting point (C): 72.50 2

232.00 2 Boiling point (C):

Density (g/cc, 20C): (liquid, 77 C) 1.1690 2

Solubility in water (mg/L, 25C): Soluble in hot water 2

HALOGENATED N.O.S. August 1987 CAS #: 53-86-1 Indomethacin Other Names: 1-(4-Chlorobenzoyl)-5-methoxy-2-methyl-1H-indole-3-acetic acid; 1-(p-Chlorobenzoyl)-5-methoxy-2-methyl-3-indolylacetic acid; Amuno; Artacin; Artinovo; Artrivia; Confortid; Idomethine; Imbrilon; Inacid; Indacin; Indocid; Indocin; Indomed; Indomee; Indoptic; Indo-Tablinen; Inflazon; Infrocin; Inteban SP; Lausit; Mezolin; Mikametan; Mobilan; Tannex Haz Waste #: P025 Ind/EPA Gen #: C19H16CIN04 Formula: Description: Crystals exhibiting polymorphism PHYSICAL PROPERTIES: - - - - - - - - - - -Molecular weight: 357.81 2 Melting point (C): One crystalline form melts about 155., another about 162. Solubility in water (mg/L, 25C): Practically insoluble 2 August 1987 HALOGENATED N.O.S. CAS #: 5344-82-1 1-(o-Chlorophenyl)thiourea Other Names: 2-Chlorophenyl thiourea Haz Waste #: P026 Ind/EPA Gen #: C7H7CIN2S Formula: PHYSICAL PROPERTIES: - - - - - - - - - - - - - - -– – – Ref. Molecular weight: 186.67 14 August 1987 HALOGENATED N.O.S. CAS #: 542-76-7 3-Chloropropionitrile Other Names: beta-Chloropropionitrile; 3-Chloropropanenitrile; 3-Chloropropanonitrile Haz Waste #: P027 Ind/EPA Gen #: Formula: C1CH2CH2CN Description: Liquid; acrid, characterisitic odor; poisonous; readily penetrates skin to produce systemic cyanide poisoning, death (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.

Molecular weight:

Melting point (C):

Boiling point (C):

Density (g/cc, 20C):

Vapor pressure (torr, 25C):

Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): 89.53

2

-51.00

(decomposes) 176.00 2

·(25 C) 1.1363 2

0.00000E+00 2 0.00000E+00 2

4.5 g/100 mL 2

August 1987 CAS #: 100-44-7

Hydrolysis rate, water (1/month, 25C):

HALOGENATED N.O.S. Benzyl chloride

Other Names: (Chloromethyl)benzene; alpha-Chlorotoluene Haz Waste #: PO28 Ind/EPA Gen #: KO15 Formula: C6H5CH2C1 Description: Very refractive liquid; rather unpleasant, irritating odor; intensely irritating to skin, eyes, mucous membranes; large doses can cause CNS depression (Ref. 2); suspected carcinogen (Ref. 12) PHYSICAL PROPERTIES: - - - - -- - - - Ref. Molecular weight: 126.58 2 Melting point (C): -43.00 to -48.00 3 Boiling point (C): 179.00 3 Density (g/cc, 20C): 1.1000 2 (22 C) 1.0000 3 Vapor pressure (torr, 25C): Henry's lawwconst (atm m**3/mol, 25C): 0.52200E-03 3 RETENTION PROPERTIES: -------- - - - - - Ref. Hydrolysis rate, water (1/month, 25C): 0.02 15 August 1987 HALOGENATED N.O.S. CAS #: 75-44-5 Phosgene Other Names: Carbonyl chloride; Carbonic dichloride; Chloroformyl chloride Haz Waste #: P095 Ind/EPA Gen #: Formula: C12C=0Description: Colorless, highly toxic gas; suffocating odor; insidious poison as it is not irritating immediately, even when fatal concentrations are inhaled (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - ----- Ref. Molecular weight: 98.92 3 Melting point (C): -118.00 3 Boiling point (C): 3 8.10 Density (g/cc, 20C): (0 C; clear, colorless, fuming liquid) 1.4320 Vapor pressure (torr, 25C): (20 C) 1220.00 3 Solubility in water (mg/L, 25C): Decomposes 3 RETENTION PROPERTIES: - - - - ----- Ref.

Rapid 3

August 1987 CAS #: 594-42-3

HALOGENATED N.O.S. Trichloromethanethiol

CN3 #. 394-42-3	TI TCITO Olle Glane Gitto
Thiocarbonyl tetrachl Haz Waste #: P118 Ind/EPA Formula: Cl3CSCl; CCl4S Description: Pale yellow oil; high	
200 C (Ref. 13)	D A
PHYSICAL PROPERTIES:	186.00 13 149.00 13 1.6900 13
August 1987 CAS #: 98-87-3	HALOGENATED N.O.S. Benzal chloride
Haz Waste #: U017 Ind/EPA Formula: C6H5CHC12; C7H6C12	oha-Dichlorotoluene; Benzylene chloride A Gen #:
	d; fumes in air; vapors irritate the
eyes; pungent odor (R PHYSICAL PROPERTIES:	
Molecular weight:	. 161.03 2
Melting point (C): Boiling point (C):	-17.00 2 205.00 2
Density (g/cc, 20C):	1.2600 2
Solubility in water (mg/L, 25C):	Insoluble 2
August 1987 CAS #: 111-91-1	HALOGENATED N.O.S. bis-(2-chloroethyoxy)methane
Formula: (CICH2CH2O)2CH2	A Gen #:
PHYSICAL PROPERTIES:	Ref. 173.10 3
Melting point (C):	-32.80 3
Boiling point (C): Vapor pressure (torr, 25C):	218.00 3 <0.1 (20 C) 3
Henry's law const (atm m**3/mol, 2	25C): 0.37800E-06 3
Solubility in water (mg/L, 25C): RETENTION PROPERTIES:	(? C) 0.81000E+05 3
log (octanol/water) partition (250	2): 1.26 3
Partition coefficient, soil/water:	2.2900 17
Hydrolysis rate, water (1/month, 2 Biodegradation rate in water (1/mo	

August 1987 CAS #: 111-44-4

HALOGENATED N.O.S. bis(2-chloroethyl)ether

Other Names: Dichloroethyl ether; bis(beta-chloroethyl)ether Ind/EPA Gen #: Haz Waste #: U025 (CICH2CH2)20 Formula: PHYSICAL PROPERTIES: - - ----- Ref. 143.00 3 Molecular weight: -46.80 3 Melting point (C): Boiling point (C): 178.00 (20 C) 0.7100 3 Vapor pressure (torr, 25C): 0.21600E-04 3 Henry's law const (atm m**3/mol, 25C): Solubility in water (mg/L, 25C): 0.10200E+05 3 RETENTION PROPERTIES: - - - - - -- - - Ref. log (octanol/water) partition (25C): 1.58 3 4.4600 17 Partition coefficient, soil/water: Hydrolysis rate, water (1/month, 25C): Slow hydrolysis 3 Biodegradation rate in water (1/month): Significant 3 HALOGENATED N.O.S. August 1987 CAS #: 494-03-1 N, N-Bis (2-chloroethyl) -2-naphthylamine Other Names: Chlornaphazine; Dichloroethyl-beta-naphthylamine; CB 1048 beta-Naphthyldi(2-chloroethyl)amine; R48; Clornaftina; beta-Naphthylbis(beta-chloroethyl)amine; Erysan; Di(2-chloroethyl)-b-naphthylamine Haz Waste #: U026 Ind/EPA Gen #: C14H15C12N Formula: Description: Platelets; can produce bladder tumors; listed as a carcinogen by the EPA (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - -Molecular weight: 268.20 2 54.00 to Melting point (C): 56.00 2 Boiling point (C): (5 mm Hg) 210.00 2 Solubility in water (mg/L, 25C): Very sparingly soluble 2

HALOGENATED N.O.S.

bis-(2-chloroisopropyl)ether

```
Other Names: 2,2'-Dichloroisopropyl ether
                          Ind/EPA Gen #:
Haz Waste #: U027
             (C1CH2CH(CH3))20
Formula:
PHYSICAL PROPERTIES: - - - - - - - - - - - -
                                                              - - - Ref.
                                                              171.00
Molecular weight:
                                                              -97.00
Melting point (C):
                                                                     3
Boiling point (C):
                                                              189.00 3
                                                       (20 C) 0.8500 3
Vapor pressure (torr, 250):
Henry's law const (atm m**3/mol, 25C):
                                                         0.15300E-03 3
Solubility in water (mg/L, 25C):
                                                     (? C) 1700.00 3
                                                       RETENTION PROPERTIES: - - - - - -
log (octanol/water) partition (250):
                                                               2.58 3
                                                            35.7500 17
Partition coefficient, soil/water:
Hydrolysis rate, water (1/month, 25C):
                                                     Slow hydrolysis 3
Biodegradation rate in water (1/month):
                                                              Rapid 3
August 1987
                                  HALOGENATED N.O.S.
CAS #:
         74-83-9
                                  Bromomethane
Other Names: Methyl bromide; Monobromomethane; Embafume
Haz Waste #: UO29
                          Ind/EPA Gen #:
Formula:
             CH3Br
Description: Colorless gas; usually odorless; burning taste;
             non-flammable in air; forms a hydrate below 4 C;
             inhalation causes dizziness, hadache, nausea, etc. leading
             to death; chronich exposure can cause CNS depression or
             kidney injury (Ref. 2)
PHYSICAL PROPERTIES: - - - - - -
                                                              - - - Ref.
Molecular weight:
                                                              94.94 3
Melting point (C):
                                                              -93.66 2
Boiling point (C):
                                                                4.60 3
Density (g/cc, 20C):
                                                   (g/L, gas) 3.9740 2
                                                      (20 C) 1420.00 3
Vapor pressure (torr, 250):
                                                              0.2200 3
Henry's law const (atm m**3/mol, 25C):
                                                       (20 C) 900.00 3
Solubility in water (mg/L, 25C):
RETENTION PROPERTIES: - - - - -
                                                            ---- Ref.
log (octanol/water) partition (25C):
                                                               1.10 3
Partition coefficient, soil/water:
                                                              1.6400 17
Hydrolysis rate, water (1/month, 25C):
                                                             8232.00 15
```

August 1987

CAS #: 101-55-3

HALOGENATED N.O.S.

4-Bromophenal phenyl ether

Other Names: 1-Bromo-4-phenoxy benzene; 4-Bromophenyl ether Haz Waste #: U030 Ind/EPA Gen #: Formula: C6H50C6H4Br PHYSICAL PROPERTIES: - - - -- - - - Ref. Molecular weight: 249.10 3 Melting point (C): 18.70 3 310.00 3 Boiling point (C): (20 C) 0.15000E-02 3 Vapor pressure (torr, 250): - - - Ref. RETENTION PROPERTIES: - - - - - - - log (octanol/water) partition (25C): 4.28 3 Partition coefficient, soil/water: 1235.00 17 Hydrolysis rate, water (1/month, 25C): Not significant 3 Biodegradation rate in water (1/month): Not significant 3

August 1987 HALOGENATED N.O.S CAS #: 353-50-4 Carbonyl fluoride

Other Names: Carbon oxyfluoride; C. Haz Waste #: UO33 Ind/EPA Formula: COF2; CF20		
Description: Pungent, very hygrosc	opic gas; strong irritant to skin,	
eyes, mucous membrane	s, respiratory tract (Ref. 2)	
PHYSICAL PROPERTIES:		
Molecular weight:	66.01 2	
Melting point (C):	-114.00 2	
Boiling point (C):	-83.10 2	
Density (g/cc, 20C):	(liquid, -114 C) 1.1390 2	
RETENTION PROPERTIES:	Ref	:
Hydrolysis rate, water (1/month, 2		·

HALOGENATED N.O.S. August 1987 CAS #: 75-87-6 Trichloroacetaldehyde Other Names: Chloral; Anhydrous chloral Haz Waste #: U034 Ind/EPA Gen #: C13CCH0; C2HC130 Formula:

addiction PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref. 147.40 2 Molecular weight: -57.50 2 Melting point (C):

Description: Oily liquid; pungent, irritating odor; abuse may lead to

Boiling point (C): 97.80 2 1.5100 2 Density (g/cc, 20C):

Solubility in water (mg/L, 25C):

Freely soluble forming chloral hydrate 2

HALOGENATED N.O.S. August 1987 Chlorambucil CAS #: 305-03-3

Other Names: Ambochlorin; 4-[Bis(2-chloroethyl)amino]benzenebutanoic

acid; 4-[p-[Bis(2-chloroethyl)amino]phenyl]butyric acid; gamma-[p-di(2-Chloroethyl)aminophenyl]butyric acid;

N, N-di-2-Chloroethyl-gamma-p-aminophenylbutyric acid; Chloraminophene; CB 1348; Amboclorin; Leukeran; Amboclorin;

Chlorbutin; Bhlorbutine; Ecloril; Elcoril; Linfolizin;

Linfolysin; NCI-CO3485; NSC-3088; Chloraminophen;

Phenylbutyric acid nitrogen mustard; Leukersan; Leukoran

Haz Waste #: U035 Ind/EPA Gen #:

C14H19C|2N02 Formula:

Description: Flattened needles; listed as a carcinogen by the EPA; used

in cancer chemotherapy

- - - - Ref. Molecular weight: 304.23 2 Melting point (C): 66.00 2 64.00 to

HALOGENATED N.O.S. Chlorodibromomethane

Other Names: Dibromochloromethane Haz Waste #: U040,U065		
August 1987 HALOGENATED N CAS #: 106-89-8 1-Chloro-2,3-		
Other Names: Epichlorohydrin; Chloromethyloxirane; dl-alpha-Epichlorohydrin; gamma-Chloropropylene oxide Haz Waste #: UO41		

HALOGENATED N.O.S.

2-Chloroethyl vinyl ether

```
Other Names: Vinyl 2-chloroethyl ether
Haz Waste #: UO42 Ind/EPA Gen #:
            C1CH2CH2OCH=CH2
Formula:
PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.
Molecular weight:
                                                           106.60 3
                                                           -69.70 3
Melting point (C):
Boiling point (C):
                                                           108.00 3
                                                    (20 C) 26.7500 3
Vapor pressure (torr, 25C):
log (octanol/water) partition (25C):
                                                            1.28 3
Partition coefficient, soil/water:
                                                           2.4300 17
Hydrolysis rate, water (1/month, 25C): Minor hydrolysis, if adsorbed 3
Biodegradation rate in water (1/month):
                                                      Significant 3
August 1987
                                HALOGENATED N.O.S.
CAS #:
        75-01-4
                                Vinyl chloride
Other Names: Chloroethylene; Airex; Armodour; Astralon; Bakelite; Benvic;
            Boltaron; Bonloid; Breon; Caliplast; Carina; Chlorethene;
            Chlorethylene; Chloroethene; Chlorostop; Conoco 5385;
            Conoco 5425; Conoco 7200; Contizell; Dacovin; Daran; Darvic;
            Daycell; Cecatone; Cecelith H; Diamond Shamrock 40; Diamond
            Shamrock 71; 1,1-Dichloroethene polymer with chlorethene;
            Dorlyl; Dow latex 874; Dow 874; Duraform; Durvyl; Dynadur;
            E-PVC; Esteproz; Flocor; Flovic; Genotherm; Geon; Kaylite;
            Klegecell; Koroseal; Laplen; Latex SVKH; Marvinal; Marvinol;
            Norvinyl; Opalon; PVC; Vinyon; Vinoflex; (many more)
                         Ind/EPA Gen #: KO20
Haz Waste #: U043
Formula:
            CH2=CHCI; C2H3CI
Description: Colorless gas, flammable, polymerizes in light or in
            presence of catalyst (Ref. 1); causes "vinyl chloride
            disease"; listed as a carcinogen by the EPA (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref. ,
                                                            62.50 2
Molecular weight:
Melting point (C):
                                                          -153.80 2
Boiling point (C):
                                                           -13.37 2
Vapor pressure (torr, 25C):
                                                          2660.00 3
Henry's law const (atm m**3/mol, 25C):
                                                           0.0360 3
Solubility in water (mg/L, 25C): 1.1000 3 RETENTION PROPERTIES: ----- Ref.
log (octanol/water) partition (25C):
                                                             0.60 3
Partition coefficient, soil/water:
                                                           0.5900 17
```

HALOGENATED N.O.S. Chloromethane

Other Names: Methyl chloride	
Haz Waste #: UO45 Ind/EPA	Gen #:
Formula: CH3C1	
Description: Colorless gas; compre	
ethereal odor and swe	et taste; poisonous; injury to liver,
	r; used as a refrigerant and local
anesthetic (Ref. 2)	D. C
PHYSICAL PROPERTIES:	Ref.
Molecular weight:	50.49 2 -97.70 3
Melting point (C): Boiling point (C):	-97.70 3 -24.20 3
Vapor pressure (torr, 25C):	(20 C) 3765.00 3
Henry's lawwconst (atm m**3/mol, 2	
Solubility in water (mg/L, 25C):	6450.00 to 7250.00 3
RETENTION PROPERTIES:	Ref.
log (octanol/water) partition (250): 0.91 3
Partition coefficient, soil/water:	1.1300 17
Hydrolysis rate, water (1/month, 2	
Biodegradation rate in water (1/mo	nth): Not important 3
August 1987	HALOGENATED N.O.S.
CAS #: 107-30-2	Chloromethyl methyl ether
	· · · · · · · · · · · · · · · · · · ·
Other Names: Chloromethoxymethane;	Methyl chloromethyl ether;
	yl ether; Chlorodimethyl ether; CMME;
	imethyl ether; Methoxychloromethane;
	e; Monochlorodimethyl ether
Haz Waste #: UO46 Ind/EPA	Gen #:
Formula: CH30CH2C1; C2H5C10	
	hnical grade listed as a carcinogen by
	hnical grade contains 1-8% BCME
PHYSICAL PROPERTIES:	er; CAS#542-88-1) (Ref. 12) Ref.
Molecular weight:	80.51 2
Boiling point (C):	59.00 2
Density (g/cc, 20C):	1.0605 2
Solubility in water (mg/L, 25C):	Soluble 12
	22.02.0

August 1987 CAS #: 91-58-7 HALOGENATED N.O.S. 2-Chloronaphthalene

Other Names: Beta-chloronaphthalene Haz Waste #: UO47 Ind/EPA Gen #: C10H7C1 Formula: Description: Monoclinic plates, leaflets PHYSICAL PROPERTIES: - - - - - - - - - - ---- Ref. Molecular weight: 162.61 2 Melting point (C): 61.00 3 Boiling point (C): 256.00 3 Vapor pressure (torr, 25C): (20 C) 0.0170 3 Henry's law const (atm m**3/mol, 25C): 0.61200E-03 3 Solubility in water (mg/L, 25C): 6.7400 3 RETENTION PROPERTIES: --------- - - Ref. log (octanol/water) partition (25C): 4.12 3 Partition coefficient, soil/water: 882.00 17 Biodegradation rate in water (1/month): Significant 3 August 1987 HALOGENATED N.O.S. CAS #: 3165-93-3 4-Chloro-o-toluidine hydrochloride Other Names: 4-Chloro-2-methyl benzenamine hydrochloride Haz Waste #: U049 Ind/EPA Gen #: Formula: C7H8C1N.C1H - - - - Ref. Molecular weight: 178.06 17 August 1987 HALOGENATED N.O.S. CAS #: 106-93-4 1,2-Dibromoethane Other Names: Ethylene dibromide; Aadibroom; Bromofume; Celmide; DBE; Dibromoethane; alpha, beta-Dibromoethane; sym-Dibromoethane; Dowfume EDB; Dowfume W 85; E-D-Bee; EDB; EDB-85; ENT 15349; Ethylene bromide; Fumo-gas; Glycol bromide; Iscobrome D; Glycol dibromide; Kopfume; NCI-CO0522; Nefis; Nephis: Pestmaster; Sanhyuum; Soilbrom; Soilbrome-85; Soilfume; Unifume; Soilbrom-90EC; Soilbrom-40; Dowfume 40 Haz Waste #: U067 Ind/EPA Gen #: Formula: BrCH2CH2Br; C2H4Br2 Description: Heavy liquid; chloroform odor; listed as a carcinogen by the EPA (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - - -- - - Ref. Molecular weight: 187.90 3 Melting point (C): 9.97 3 Boiling point (C): 132.00 3 Density (g/cc, 20C): (25 C) 2.1720 2 Vapor pressure (torr, 25C): (20 C) 11.0000 3 Henry's law const (atm m**3/mol, 25C): 0.88200E-03 3 Solubility in water (mg/L, 25C): (30 C) 4310.00 3

HALOGENATED N.O.S. Dibromomethane

Other Names: Methylene bromide Haz Waste #: U068 Ind/EPA Formula: CH2Br2 Description: Liquid PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Boiling point (C): Density (g/cc, 20C): Solubility in water (mg/L, 25C): RETENTION PROPERTIES: Hydrolysis rate, water (1/month, 25)	1	Ref. 173.86 2 -52.70 2 97.00 2 2.4956 2 1.93 g/kg, 30 C 2 Ref. 2196.00 15
August 1987 CAS #: 91-94-1	HALOGENATED N.O.S. 3,3'-Dichlorobenzidi	ne
Other Names: C.I. 23060; Curithane 4,4'-Diamino-3,3'-dick 3,3'-Dichloro-4,4'-Dia Haz Waste #: U073 Ind/EPA Formula: H2N(CI)C6H3-C6H3NH2(C Description: Needles; listed as a constant and appropriate.	hlorobiphenyl; Dichlo aminobiphenyl; Dichlo Gen #: I); C12H1OCI2N2	probenzidine base; probenzidine; DCB
PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Solubility in water (mg/L, 25C): RETENTION PROPERTIES: Hydrolysis rate, water (1/month, 25) Biodegradation rate in water (1/month)		

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August 1987
                                   HALOGENATED N.O.S.
          75-09-2
CAS #:
                                   Methylene chloride
Other Names: Dichloromethane; Methylene dichloride;
             Methylene bichloride
                           Ind/EPA Gen #: F001,F002
Haz Waste #: U080
             CH2C12
Formula:
Description: Colorless liquid; vapor not flammable (Ref. 1)
                                                                 - - - Ref.
PHYSICAL PROPERTIES: - - - - - -
Molecular weight:
                                                                 84.94
                                                                        2
                                                                -96.70
Melting point (C):
                                                                        1
Boiling point (C):
                                                                 39.75
                                                                        2
Density (g/cc, 20C):
                                               specific gravity 1.3200
                                                                        1
Vapor pressure (torr, 25C):
                                                                362.00
Henry's law const (atm m**3/mol, 25C):
                                                           0.31900E-02
                                                                        3
Solubility in water (mg/L, 25C):
                                                           0.16700E+05
                                                                       3
RETENTION PROPERTIES:
                                                               --- Ref.
log (octanol/water) partition (25C):
                                                                  1.25 3
Partition coefficient, soil/water:
                                                                2.2400 17
Hydrolysis rate, water (1/month, 25C):
                                                               8448.00 15
Biodegradation rate in water (1/month):
                                                           Significant 3
                                   HALOGENATED N.O.S.
August 1987
CAS #:
         148-82-3
                                   Melphalan
Other Names: 4-[Bis(2-chloroethyl)amino]-L-phenylalanine; L-PAM;
             p-Di(2-chloroethyl)amino-L-phenylalanine; Melfalan;
             L-phenylalanine mustard; Alanine nitrogen mustard;
             L-Sarcolysine; NSC-8806; CB 3025; Alkeran; Sarcoclorin;
             AT-290; L-3-(p-(bis(2-chloroethyl)amino)phenyl)alanine;
             p-N-Bis(2-chloroethyl)amino-L-phenylalanine; 3025 C.B.;
             3-(p-(Bis(2-chloroethyl)amino)phenyl)-L-alanine; L-PAM;
             Levofalan; Mel; Phenylalanine mustard; p-L-Sarcolysin;
             L-Sarcolysin; Sarcolysine; Sarkolysin; SK-15673;
```

Phenylalanine nitrogen mustard; L-Sarkolysin

depression may occur (Ref. 2); used in cancer chemotherapy

- - - Ref.

305.20 2

2

(decomposes) 182.00 to 183.00

Practically insoluble 2

Ind/EPA Gen #:

Description: Needles; listed as a carcinogen by the EPA; bone marrow

Haz Waste #: U150

Molecular weight:

Melting point (C):

C13H18C12N202

Solubility in water (mg/L, 25C):

Formula:

HALOGENATED N.O.S. Methyl Chlorocarbonate

Haz Waste #: U156 Ind/EPA Formula: CICO2CH3; C2H3ClO2	I, methyl ester; Methyl chloroformate Gen #: strongly irritating to eyes (Ref. 2)
August 1987 CAS #: 101-14-4	HALOGENATED N.O.S. 4,4'-Methylenebis[2-chloroaniline]
chlorophenyl)methane; Haz Waste #: U158 Ind/EPA Formula: C13H12Cl2N2 Description: Flakes	chloro-diphenylmethane; di-4-(Amino-3- ; Methylenebis(o-chloroaniline)
August 1987 CAS #: 608-93-5	HALOGENATED N.O.S. Pentachlorobenzene
Haz Waste #: U183 Ind/EPA Formula: 1,2,3,4,5-C6HCI5 PHYSICAL PROPERTIES: Molecular weight: Melting point (C): Boiling point (C): Vapor pressure (torr, 25C): Henry's law const (atm m**3/mol, 2 Solubility in water (mg/L, 25C):	250.30 3 85.00 1 276.00 1 (kPa) 0.88900E-02 3

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HALOGENATED N.O.S.
August 1987
         76-01-7
                                 Pentachloroethane
CAS #:
Other Names: Pentalin
                        Ind/EPA Gen #:
Haz Waste #: U184
            CCI3CHCI2; C2HCI5
Formula:
Description: Liquid; chloroform-like odor; irritant; narcotic (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                            202.31 2
Melting point (C):
                                                            -29.00
                                                                   2
                                               161.00 to
                                                            162.00 2
Boiling point (C):
                                                      (25 C) 1.6712 2
Density (q/cc, 20C):
                                                         Insoluble 2
Solubility in water (mg/L, 25C):
                                 HALOGENATED N.O.S.
August 1987
CAS #: 95-94-3
                                 1,2,4,5-Tetrachlorobenzene
                 Ind/EPA Gen #: K085,K042
Haz Waste #: U207
Formula: 1,2,4,5-C6H2C14
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                            215.90
Melting point (C):
                                                            140.00
                                                                   3
Boiling point (C):
                                                            243.00
                                              (kg/L, liquid) 1.8330 1
Density (g/cc, 20C):
Vapor pressure (torr, 25C):
                                                       (kPa) 0.0101 3
Henry's law const (atm m**3/mol, 25C):
                                              (kPa m**3/mol) 0.2610
Solubility in water (mg/L, 25C):
                                                    (g/m**3) 0.5950 3
August 1987
                                 HALOGENATED N.O.S.
CAS #:
        66-75-1
                                 Uracil mustard
Other Names: 5-[Bis(2-chloroethyl)amino]-2,4-(1H,3H)-pyrimidinedione;
            5-[Bis(2-chloroethyl)amino]uracil; Uramustine;
            2,6-Dihydroxy-5-bis[2-chloroethyl]aminopyramidine;
            5-[di-(beta-chloroethyl)amino]uracil; Demethyldopan;
```

Desmethyldopan; NSC-34462; U-8344

PHYSICAL PROPERTIES: - - - - - - - - - - - - - -

Ind/EPA Gen #:

Haz Waste #: U237

Molecular weight:

Melting point (C):

Description: Crystals

Formula:

C8H11C12N302

Solubility in water (mg/L, 25C):

252.10 2

(decomposes) 206.00 2

Sparingly soluble 2

August 1987 CAS #: 120-82-1 HALOGENATED N.O.S. 1,2,4-Trichlorobenzene

Other Names: unsym-Trichlorobenzene Haz Waste #: X015 Ind/EPA Gen #: K085		
Formula: 1,2,4-C6H3Cl3		
PHYSICAL PROPERTIES:		Ref.
Molecular weight:	181.45	1
Melting point (C):	17.15	1
Boiling point (C):	213.80	- 1
Density (g/cc, 20C):	(liquid) 1.4483	1
Vapor pressure (torr, 25C):	0.4200	3
Henry's law const (atm m**3/mol, 25C):	0.14200E-02	3
Solubility in water (mg/L, 25C):	30.0000	3
RETENTION PROPERTIES:		Ref.
log (octanol/water) partition (25C):	4.26	3
Partition coefficient, soil/water:	1180.00	17
Biodegradation rate in water (1/month):	Bioaccumulates	3

7.14 Pharmaceuticals

PHARMACEUTICALS . August 1987 5-(Aminomethyl)-3-isoxazolol CAS #: 2763-96-4 Other Names: Muscimol; 5-(Aminomethyl)-3(2H)-isoxazolone; 5-Aminomethyl-3-hydroxy-isoxazole; 3-Hydroxy-5-aminomethylisoxazole; Agarin; Pantherine Ind/EPA Gen #: Haz Waste #: P007 C4H6N202 Formula: Description: Crystals; potent CNS depressant and GABA antagonist (Ref. 2) PHYSICAL PROPERTIES: - - - - - - - - - - -- - - - Ref. 114.10 2 Molecular weight: decomposes 175.00 2 Melting point (C): **PHARMACEUTICALS** August 1987 504-24-5 CAS #: 4-Aminopyridine Other Names: 4-Pyridinamine; gamma-Pyridylamine Haz Waste #: P008 Ind/EPA Gen #: C5H6N2 Description: Crystals; highly toxic (Ref. 13) PHYSICAL PROPERTIES: - - - - - - - -- - - Ref. Molecular weight: 94.00 13 Melting point (C): 158.00 13 Boiling point (C): (12 mm Hg) 180.00 13 August 1987 **PHARMACEUTICALS** CAS #: 51-43-4 Epinephrine Other Names: 4-[1-Hydroxy-2-(methylamino)ethyl]-1,2-benzenediol; 3,4-Dihydroxy-alpha-[(methylamino)methyl]benzyl alcohol; 1-(3,4-dihydroxyphenyl)-2-(methylamino)ethanol; 3-4-dihydroxy-1-[1-hydroxy-2-(methylamino)-ethyl]benzene; Methylaminoethanolcatechol; Adrenalin I-Form--Adnephrine; Adrena; Adrenamine; Adrenine; Adrin; Epinephran; Epirenan; Hemostasin; Renaleptine; Renalina; Suprarenaline; Surrenine; Vasoconstrictine; etc. Haz Waste #: PO42 Ind/EPA Gen #: Formula: C9H13N03

Description: 1-Form occurs naturally (described here); minute crystals;

undergo oxidation (Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - -

Solubility in water (mg/L, 25C):

Molecular weight:

Melting point (C):

combines with acids forming water soluble salts; solutions

211.-212.; approx. 215. (decomposes) when heated rapidly

Sparingly soluble 2

183.20 2

August 1987 PHARMACEUTICALS
CAS #: 115-02-6 Azaserine

Other Names: L-Serine diazoacetate(ester); O-Diazo-acetyl-L-serine;
C1 337; CN 15757; P 165
Haz Waste #: U015 Ind/EPA Gen #:
Formula: N2CHC00CH2CH(NH2)C00H; C5H7N304
Description: Orthorhombic, pale yellow to green crystals

August 1987 PHARMACEUTICALS
CAS #: 20830-81-3 Daunomycin

Other Names: 8-Acetyl-10-[(3-amino-2,3,6-trideoxy-alpha-L-lyxo-

hexopyranosyl)oxy]-7,8,9,10-tetrahydro-6,8,11-trihydroxy-1-methoxy-5,12-naphtacenedione; Daunorubicin; Rubidomycin;

- - Ref.

173.13 2

Very soluble 2

(decomposes) 146.00 to 162.00 2

Leukaemomycin C; Cerubidin

Haz Waste #: U059 Ind/EPA Gen #:

Formula: C27H29N010

Molecular weight: Melting point (C):

Solubility in water (mg/L, 25C):

PHYSICAL PROPERTIES: - - - - - - - - - - - - - - Ref.
Molecular weight: 527.53 17

August 1987 PHARMACEUTICALS
CAS #: 56-53-1 Diethylstilbestrol

Other Names: 4,4'-(1,2-Diethyl-2,3-ethenediyl)bisphenol; alpha,alpha'Diethylstilbenediol; Stilbestrol; 3,4-bis(p-hydroxyphenyl)3-hexene; 4,4'-dihydroxy-alpha,beta-diethylstilbene; DES;
Antigestil; Bio-des; Bufon; Cyren A; Distilbene; Domestrol;

Estrobene; Estrosyn; Fonatol; Grafestrol; Hi-Bestrol;

Makarol; Micrest; Milestrol; Neo-Destranol I; Destrogenine;

Oestromenin; Oestromensyl; Oestromon; Palestrol; Serral; Sexocretin; Sibol; Stilbetin; Stilboefral;

Stilboestroform; Stilkap; Synestrin (tablets); Synthoestrin; Vagestrol; Acnestrol; Agostilben; DEB; DESMA; (many more)

Haz Waste #: U089 Ind/EPA Gen #:

Formula: C18H2OO2

Description: Small plates; banned by the FDA in 1979 as a growth

promotant in livestock; listed as a carcinogen by the EPA;

used to prevent miscarriage in humans; an estrogen (Ref. 2)

PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.

Molecular weight: 268.34 2 Melting point (C): 169.00 to 172.00 2

Solubility in water (mg/L, 25C): Almost insoluble 2

```
PHARMACEUTICALS
August 1987
CAS #:
          94-58-6
                                   Dihydrosafrole
Other Names: 1,2-(Methylenedioxy)-4-propylbenzene;
             5-Propyl-1,3-benzodioxole;
             4-Propyl-1,2-methylenedioxybenzene
                          Ind/EPA Gen #:
Haz Waste #: U090
             C10H1202
Formula:
Description: Oily liquid
PHYSICAL PROPERTIES: - -
                                                                - - - Ref
Molecular weight:
                                                                164.22 14
                                                                228.00 14
Boiling point (C):
Density (g/cc, 20C):
                                                                1.0695 14
                                   PHARMACEUTICALS
August 1987
CAS #:
         120-58-1
                                   Isosafrole
Other Names: 5-(1-Propenyl)-1,3-benzodioxole; 1,2-(Methylenedioxy)-4-
             propenylbenzene
Haz Waste #: U141
                           Ind/EPA Gen #:
             C10H1002
Formula:
Description: Liquid; trans-form--odor of anise (Ref. 2)
                                                                 - - - Ref.
PHYSICAL PROPERTIES: - - - - - - - - - - -
Molecular weight:
                                                                162.18
                                               trans--8.2; cis-- -21.5
Melting point (C):
Boiling point (C):
                            trans--253.; cis--77. to 79. at 3.5 mm Hg
Density (g/cc, 20C):
                                            trans--1.1206; cis--1.1182 2
                                   PHARMACEUTICALS
August 1987
CAS #:
         303-34-4
                                   Lasiocarpine
Other Names: 2-Methyl-2-butenoic acid 7-[[2,3-dihydroxy-2-(1-
             methoxyethyl)-3-methyl-1-oxobutoxy]methyl]-2,3,5,7a-
             tetrahydro-1H-pyrrolizin-1-yl ester
                           Ind/EPA Gen #:
Haz Waste #: U143
             C21H33N07
Formula:
Description: Colorless leaflets
PHYSICAL PROPERTIES: - - - - - - - -
                                                                     – Ref.
Molecular weight:
                                                                411.40
                                                   94.00 to
                                                                 95.50 2
Melting point (C):
Solubility in water (mg/L, 25C):
                                                   Difficultly soluble 2
```

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PHARMACEUTICALS
August 1987
         70-25-7
CAS #:
                                  N-Methyl-N'-nitro-N-nitrosoguanidine
Other Names: N-Nitroso-N-methyl-N'-nitro-guanidine; N-Methyl-N-nitroso-
            N'-nitroguanidine; MNNG
Haz Waste #: U163
                          Ind/EPA Gen #:
            C2H5N503
Formula:
Description: Yellow crystals; carcinogen and mutagen (Ref. 2)
- - - Ref.
Molecular weight:
                                                              147.10 2
Melting point (C):
                                                  (decomposes) 118.00 2
                                  PHARMACEUTICALS
August 1987
CAS #:
         56-04-2
                                  Methylthiouracil
Other Names: 2,3-Dihydro-6-methyl-2-thioxo-4(1H)-pyrimidinone;
            6-methyl-2-thiouracil; 4-methyl-2-thiouracil; MTU; Alkiron;
            Antibason; Basecil; Basethyrin; Methiacil; Methicil;
            Methiocil; Muracil; Prostrumyl; Strumacil; Thimecil;
            Thyreostat I
Haz Waste #: U164
                          Ind/EPA Gen #:
            C5H6N20S
Formula:
Description: Crystals; bitter taste; thyroid inhibitor (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - -
Molecular weight:
                                                              142.18 2
Melting point (C):
                      (decomposes, sublimes readily) 326.00 to 331.00 2
Solubility in water (mg/L, 25C):
                        Very slightly soluble; 1 part in 150 at 100 C 2
August 1987
                                  PHARMACEUTICALS
CAS #:
         62-44-2
                                  Phenacetin
Other Names: N-(4-Ethoxyphenyl)acetamide; p-Acetophenetidide;
            p-Ethoxyacetanilide; Acetophenetidin; para-Acetphenetidin;
            Acet-p-phenetidin; p-Acetophenetide; Acet-p-phenalide;
            1-Acetamido-4-ethoxybenzene; Aceto-para-phenalide; Anapac;
            APC; Bromo seltzer; Buff-a-comp; Citra-fort; Contradol;
            Dasin; Empiral; Empirin compound; Malex; Gelonida; Pamprin;
            Norgesic; Sinutab; Sinutab II; Soma; Phenin; Phenodyne;
            Tetracydin; Viden; Wigraine; Xaril; Zactirin compound;
            Terracydin; Super anahist; Supralgin; Synalogos; Reformin;
            Phenazetin; Phenacon; Pehnaphen; Phenacet; (many more)
                          Ind/EPA Gen #:
Haz Waste #: U187
            Slightly bitter cyrstalline scales or powder; listed as a
Formula:
Description: carcinogen by the EPA (Ref. 2); widely used analgesic and
            antipyretic (Ref. 12)
PHYSICAL PROPERTIES: - - - - -
                                                                - - - Ref.
Molecular weight:
                                                              179.21
                                                                     2
                                                 134.00 to
                                                              135.00 2
Melting point (C):
Solubility in water (mg/L, 25C):
                                                         1 g/1310 mL 2
```

PHARMACEUTICALS Reserpine

```
Other Names: 11,17-Dimethoxy-18-[(3,4,5-trimethoxybenzoyl)oxy]yohimban-
             16-carboxylic acid methyl ester; 3,4,5-trimethoxybenzoyl
             methyl reserpate; Alserin; Austrapine; Crystoserpine;
             Eskaserp; Hiserpia; Orticalm; Quiescin; Rau-sed; Reserpex;
             Reserpoid; Rivasin; Roxinoid; Sandril; Sedaraupin; Serfin; Serolfia; Serpanray; Serpasil; Serpasol; Serpate; Serpen;
             Serpine; Serpiloid; Abesta; Abicol; Adelfan; ENT-50146;
             Enipresser; Gammaserpine; Anquil; R-E-S; Rausan; Roxel;
             Tenserlix; USAF CB-27; Triserpin; Unilord; Terbolan;
             Vio-serpine; V-Serp; Tensional; Tensionorme; (many more)
Haz Waste #: U200
                           Ind/EPA Gen #:
             C33H40N209
Formula:
Description: Long prisms; listed as a carcinogen by the EPA (Ref. 2);
             marketed under at least 150 trade names; used for treating
             hypertension and as a tranquilizer and sedative (Ref. 12)
Molecular weight:
                                                                 608.70 2
                                          (decomposes) 264.00 to 265.00
Boiling point (C):
Solubility in water (mg/L, 25C):
                                                Very sparingly soluble 2
                                    PHARMACEUTICALS
August 1987
CAS #:
          81-07-2
                                    Saccharin
Other Names: 1,2-Benzisothiazol-3(2H)-one 1,1-dioxide; Benzosulfimide;
             2,3-Dihydro-3-oxobenzisosulfonazole; Saccharin insoluble;
             1,2-Dihydro-2-ketobenzisosulfonazole; o-Sulfobenzimide;
             Benzoic sulfimide; o-Sulfobenzoic acid imide; Gluside;
             Glucid; Garantose; Saccharinol; Saccharinose; Saccharol;
             Saxin; Sykose; Hermesetas; Anhydro-o-sulfaminebenzoic acid;
             Natreen; Kandiset; Gluside; Benzosulfinide; Benzosulphimide;
             3-Hydroxybenzisothiazole-S,S-dioxide; O-Benzoyl sulfimide;
             Benzo-2-sulphimide; 3-Benzisothiazolinone 1.1-dioxide
                           Ind/EPA Gen #:
Haz Waste #: U202
Formula:
             C7H5N03S
Description: Monoclinic crystals; 500 time as sweet as sugar; listed as a
             carcinogen by the EPA (Ref. 2); white crystalline powder
             (Ref. 12)
PHYSICAL PROPERTIES: - -
                                                               - - - Ref.
Molecular weight:
                                                                 183.18
                                                                        2
                                                                 229.70 2
Melting point (C):
                                                  228.80 to
Density (g/cc, 20C):
                                                                0.8280 2
Solubility in water (mg/L, 25C):
                                                       1 g/290 ml (? C)
```

```
PHARMACEUTICALS
August 1987
         94-59-7
                                Safrole
CAS #:
Other Names: 5-(2-Propenyl)-1,3-benzodioxole; 4-Allyl-1,2-
            methylenedioxybenzene; Allylcatechol methylene ether;
            Allyldioxybenzene methylene ether; m-Allylpyrocatechin
            methylene ether
Haz Waste #: U203
                         Ind/EPA Gen #:
Formula:
            C10H1002
Description: Colorless or slightly yellow liquid; sassafras odor; listed
            as a carcinogen by the EPA (Ref. 2)
- - - Ref.
Molecular weight:
                                                            162.18
Melting point (C):
                                                             11.00
Boiling point (C):
                                                            234.00 2
                                              232.00 to
Density (g/cc, 20C):
                                                            1.0960 2
Solubility in water (mg/L, 25C):
                                                         Insoluble 2
```

PHARMACEUTICALS August 1987 CAS #: 18883-66-4 Streptozotocin Other Names: 2-Deoxy-2-[[(methylnitrosoamino)-carbonyl]amino]-Dglucopyranose; 2-Deoxy-2-(3-methyl-3-nitrosoureido)-Dglucopyranose; Streptozocin; NSC-85998; U-9889; Zanosar N-D-Glucosyl-(2)-N'-nitrosomethylurea; NCI-C03167; STR; **STRZ** Haz Waste #: U206 Ind/EPA Gen #: C8H15N307 Formula: Description: Pointed platelets or prisms; listed as a carcinogen by the EPA (Ref. 2); antibiotic with potential for use in cancer chemotherapy (Ref. 12) PHYSICAL PROPERTIES: - - - - - -Molecular weight: 265.22

Melting point (C):

Solubility in water (mg/L, 25C):

2

(decomposes) 115.00 2

Soluble 2

7.15 Dyes

August 1987 **DYES** CAS #: 131-89-5 2-Cyclohexyl-4,6-dinitrophenol Other Names: 2,4-Dinitro-6-cyclohexylphenol; Dinitro-o-cyclohexylphenol; DNOCHP: SN 46 Haz Waste #: P034 Ind/EPA Gen #: Formula: C12H14N205 Description: Cyrstals - Ref. PHYSICAL PROPERTIES: - - -Molecular weight: 266.25 107.50 2 Melting point (C): 106.50 to Solubility in water (mg/L, 25C): Very slightly soluble 2 **DYES** August 1987 CAS #: 100-01-6 p-Nitroaniline Other Names: 4-Nitrobenzenamine; p-Nitraniline Ind/EPA Gen #: Haz Waste #: P077 C6H6N202 Formula: Description: Bright yellowwpowder; forms water soluble salts with mineral acids; highly toxic; absorbed through skin (Ref. 2) 138.12 2 Molecular weight: 146.00 2 Melting point (C): Boiling point (C): (calculated) 332.00 2 Solubility in water (mg/L, 25C): 1 g/1250 mL, 45 mL boiling water August 1987 **DYES** CAS #: 79-19-6 Thiosemicarbazide Other Names: Hydrazinecarbothioamide Haz Waste #: P116 Ind/EPA Gen #: Formula: NH2CSNHNH2; CH5N3S Description: White, crystalline powder PHYSICAL PROPERTIES: - - - - - - - - -- Ref. Molecular weight: 91.14 2

Melting point (C):

Solubility in water (mg/L, 25C):

182.00 to

184.00 2

Soluble 2

```
DYES
August 1987
CAS #:
          92-87-5
                                   Benzidine
Other Names: [1,1'-Bipehnyl]-4,4'-diamine; p-Diaminodiphenyl;
             p,p-Bianiline; 4,4'-Bianiline; 4,4'-Diaminobiphenyl;
             Fast Corinth base B; p,p'-Dianiline; p,p'-Diaminobiphenyl
                           Ind/EPA Gen #:
Haz Waste #: U021
             H2NC6H4.C6H4NH2 (Dihydrochloride .2HCI); C12H12N2
Formula:
Description: White or slightly-reddish, crystalline powder; poisonous;
             (Dihydrochloride--crystals); solid and vapor rapidly
             absorbed through skin; listed as a carcinogen by the EPA
PHYSICAL PROPERTIES: -
Molecular weight:
                                                                 184.20
                                                                 129.00
Melting point (C):
                                                                         3
Boiling point (C):
                                                                 402.00 3
Solubility in water (mg/L, 25C):
                                                          (12<sub>c</sub>C) 400.00 3
RETENTION PROPERTIES:
                                                                 - - - Ref.
log (octanol/water) partition (25C):
                                                                   1.81 3
Partition coefficient, soil/water:
                                                                 7.2000 17
Hydrolysis rate, water (1/month, 25C):
                                                          Not important 3
Biodegradation rate in water (1/month):
                                                             Degradable 3
                                   DYES
August 1987
CAS #:
         119-90-4
                                   3,3'-Dimethyoxybenzidine
Other Names: Dianisidine; 3,3'-Dimethoxy-[1,1'-biphenyl]-4,4'-diamine;
             3,3'-Dimethoxy-4,4'-diaminobiphenyl
Haz Waste #: U091
                           Ind/EPA Gen #:
Formula:
             C14H16N202
Description: Crystals
PHYSICAL PROPERTIES: - - - - - - - -
Molecular weight:
                                                                 244.28
                                                                         2
                                                                 138.00
Melting point (C):
                                                  137.00 to
```

Solubility in water (mg/L, 25C):

Practically insoluble 2

```
DYES
August 1987
CAS #:
         60-11-7
                                 p-Dimethylaminoazobenzene
Other Names: N.N-Dimethyl-4-(phenylazo)benzenamine; Butter yellow;
            Methyl yellow; C.I. Solvent Yellow 2; C.I. 11020;
            4-Dimethylaminoazobenzene; Atul fast yellow R; DAB; DMAB;
            Brilliant fast oil yellow; Brilliant fast yellow;
            Fast yellow; Fat yellow; Oil yellow; P.D.A.B. Sudan yellow;
            Toyo oil yellow G; USAF EK-338; Waxoline yellow AD;
            Methyl yellow; Dimethyl yellow; Oleal yellow 2G;
            Resinol yellow GR; Resoform yellow GGA; Fat yellow A;
            N,N-Dimethyl-4-aminoazobenzene; Cerasine yellow GG;
            Somalia yellow A; Stear yellow JB; (many others)
Haz Waste #: U093
                          Ind/EPA Gen #:
            C14H15N3
Formula:
Description: Yellow crystalline leaflets; listed as a carcinogen by the
            EPA (Ref. 2)
PHYSICAL PROPERTIES: - - - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             225.28 2
Melting point (C):
                                               114.00 to
                                                           117.00 2
                                                          Insoluble 2
Solubility in water (mg/L, 25C):
August 1987
                                 DYES
CAS #: 119-93-7
                                 3.3'-Dimethylbenzinidine
Other Names: o-Tolidine; 3,3'-Dimethyl-[1,1'-biphenyl]-4,4'-diamine;
            4,4'-Diamino-3,3'-dimethylbiphenyl
Haz Waste #: U095
                          Ind/EPA Gen #:
Formula:
            C14H16N2
Description: White to reddish crystals or crystalline powder
- - - - Ref.
Molecular weight:
                                                             212.28 2
Melting point (C):
                                               129.00 to
                                                            131.00 2
Solubility in water (mg/L, 25C):
                                                   Slightly soluble 2
August 1987
                                 DYES
CAS #:
        130-15-4
                                 1,4-Naphthoquinone
Other Names: 1,4-Naphthalenedione; alpha-Naphthoquinone; 1,4-Dihydro-
            1,4-diketonaphthalene
Haz Waste #: U166
                         Ind/EPA Gen #:
Formula:
            C10H602
Description: Yellow triclinic needles; odor like that of benzoquinone
PHYSICAL PROPERTIES: - - - - - - - - - - - - - Ref.
Molecular weight:
                                                             158.15
Melting point (C):
                             (begins to sublime below 100 C) 126.00
                                                                    2
Density (g/cc, 20C):
                                                                    2
                                                             1.4220
```

Solubility in water (mg/L, 25C):

Sparingly soluble 2

DYES August 1987 CAS #: 72-57-1 Trypan blue Other Names: 3,3'-[(3,3'-Dimethyl[1,1'-biphenyl]-4,4'-diyl)bis(azo)] bis[5-amino-4-hydroxy-2,7-naphthalenedisulfonic acid] tetrasodium salt; C.I. Direct Blue 14; C.I. 23850; 3,3'-[(3,3'-dimethyl-4,4'-biphenylene)bis(azo)]bis-(5-amino-4-hydroxy-2,7-naphthalenedisulfonic acid) tetrasodium salt; Tetrasodium 3,3'-[(3,3'-dimethyl-4,4'-biphenylene)bis(azo)] bis(5-amino-4-hydroxy-2,7-naphthalenedisulfonate); Sodium ditolyl-diazobis-8-amino-1-naphthol-3,6-disulfonate; Benzamine Blue; Diamine Blue; Benzo Blue; Congo Blue; Dianil Blue; Naphthylamine Blue; Niagara Blue Haz Waste #: U236 Ind/EPA Gen #: Formula: C34H24N6Na4014S4 Description: Bluish-gray powder, forming a deep blue solution with violet tinge in water; a biological stain (Ref. 2) PHYSICAL PROPERTIES: - - - -- - Ref. Molecular weight: 960.83 2

Soluble 2

Solubility in water (mg/L, 25C):

INDEX TO ORGANICS

(including primary name, P/U number and category)

1,1,1,2-Tetrachloroethane	U208	HALOGENATED SOLVENTS
1,1,1-Trichloroethane	U226	HALOGENATED SOLVENTS
1,1,2,2-Tetrachloroethane		HALOGENATED SOLVENTS
1,1,2-Trichloroethane		HALOGENATED SOLVENTS
1,1-Dichloroethane		HALOGENATED SOLVENTS
1,1-Dichloroethylene		HALOGENATED SOLVENTS
1,1-Dimethylhydrazine		REACTIVE (NON-CYANIDE)
1,2,4,5-Tetrachlorobenzene		HALOGENATED N.O.S.
1,2,4-Trichlorobenzene		HALOGENATED N.O.S.
1,2-Dibromo-3-chloropropane pesticides		HALOGENATED PESTICIDES
1,2-Dibromoethane		HALOGENATED N.O.S.
1,2-Dichlorobenzene		HALOGENATED SOLVENTS
1,2-Dichloroethane		HALOGENATED SOLVENTS
1,2-Dichloropropane		HALOGENATED SOLVENTS
1,2-Dimethyl hydrazine		REACTIVE (NON-CYANIDE)
1,2-Dipehnylhydrazine		REACTIVE (NON-CYANIDE)
1,3-Dichlorobenzene		HALOGENATED SOLVENTS
1,3-Dichloropropene		HALOGENATED SOLVENTS
1,3-Pentadiene	U186	POLYMERIZABLE
1,3-Propane sulfone		ORGAND-SULFURS
1,4-Dichloro-2-butene		HALOGENATED SOLVENTS
1,4-Dichlorobenzene		HALOGENATED PESTICIDES
1,4-Dioxane		NONHALOGENATED SOLVENTS
1,4-Naphthoquinone	U166	DYES
1-(o-Chlorophenyl)thiourea		HALOGENATED N.O.S.
1-Acetyl-2-thiourea	P002	ORGANO-SULFURS
1-Chloro-2,3-epoxypropane	U041	HALOGENATED N.O.S.
1-Naphthyl-2-thiourea	P072	NONHALOG. PESTICIDES
1-Naphthylamine	U167	ORGANO-NITROGENS
2,3,4,6-Tetrachlorophenol pesticides	U212	HALOGENATED PESTICIDES
2,4,5-Trichlorophenol pesticides	U230	HALOGENATED PESTICIDES
2,4,5-Trichlorophenoxy acetic acid pest	U232	HALOGENATED PESTICIDES
2,4,6-Trichlorophenol Pesticides	U231	HALOGENATED PESTICIDES
2,4-D Pesticides	P035	HALOGENATED PESTICIDES
2,4-Dichlorophenol Pesticides		HALOGENATED PESTICIDES
2,4-Dimethylphenol		PHENOLICS
2,4-Dinitrophenol	P048	PHENOLICS
2,4-Dinitrotoluene		ORGANO-NITROGENS
2,4-Dithiobiuret		NONHALOG. PESTICIDES
2,6-Dichlorophenol pesticides		HALOGENATED PESTICIDES
2,6-Dinitrotoluene		ORGANO-NITROGENS
2-Acetylaminofluorene		POLYNUCLEAR AROMATICS
2-Chloroethyl vinyl ether		HALOGENATED N.O.S.
2-Chloronaphthalene		HALOGENATED N.O.S.
2-Chlorophenol Pesticides		HALOGENATED PESTICIDES
2-Cyclohexyl-4,6-dinitrophenol	P034	
2-Methylaziridine		POLYMERIZABLE
2-Methyllactonitrile		POLYMERIZABLE
2-Naphthylamine		ORGANO-NITROGENS
2-Nitropropane		NONHALOGENATED SOLVENTS
2-Picoline		NONHALOGENATED SOLVENTS
3,3'-Dichlorobenzidine		HALOGENATED N.O.S.
3,3'-Dimethylbenzinidine	U095	
3,3'-Dimethyoxybenzidine	U091	
3,3-Dimethyl		NONHALOG. PESTICIDES
3-Chloropropionitrile		HALOGENATED N.O.S.
3-Methylcholanthrene		POLYNUCLEAR AROMATICS
4,4'-DDD Pesticides 4,4'-DDT Pesticides		HALOGENATED PESTICIDES
		HALOGENATED ALOS
4,4'-Methylenebis[2-chloroaniline]	0128	HALOGENATED N.O.S.

PO47 NONHALOG. PESTICIDES 4,6-Dinitro-o-cresol 4-Aminopyridine POOR PHARMACEUTICALS 4-Bromophenal phenyl ether U030 HALOGENATED N.O.S. 4-Chloro-o-toluidine hydrochloride U049 HALOGENATED N.O.S. 4-Nitrophenol U170 PHENOLICS 5-(Aminomethyl)-3-isoxazolol POO7 PHARMACEUTICALS 5-Nitrotoluidine U181 ORGANO-NITROGENS 7,12-Dimethylbenz[a]anthracene U094 POLYNUCLEAR AROMATICS PO88 NONHALOG. PESTICIDES 7-0xabicyclo Aceta I dehyde UOO1 DXYGENATED U002 NONHALOGENATED SOLVENTS Acetone UOO3 NONHALOGENATED SOLVENTS Acetonitrile UOO4 DXYGENATED Acetophenone Acetyl chloride UOO6 REACTIVE (NON-CYANIDE) Acrolein POO3 POLYMERIZABLE Acrylamide U007 POLYMERIZABLE Acrylic acid U008 POLYMERIZABLE Acrylonitrile U009 POLYMERIZABLE Aldicarb PO70 NONHALOG. PESTICIDES Aldrin Pesticides POO4 HALOGENATED PESTICIDES Allyl alcohol POO5 POLYMERIZABLE U011 NONHALOG. PESTICIDES Amitrole Ammonium picrate POO9 REACTIVE (NON-CYANIDE) Aniline U012 NONHALOGENATED SOLVENTS Auramine U014 NONHALOG. PESTICIDES **U015 PHARMACEUTICALS** Azaserine U016 POLYNUCLEAR AROMATICS Benz(c)acridine Benz[a]anthracene U018 POLYNUCLEAR AROMATICS Benzal chloride UO17 HALOGENATED N.O.S. Benzene UO19 NONHALOGENATED SOLVENTS Benzenesulfonyl chloride UO20 REACTIVE (NON-CYANIDE) Benzidine U021 DYES Benzo[a]pyrene UO22 POLYNUCLEAR AROMATICS Benzotrichloride UO23 REACTIVE (NON-CYANIDE) PO28 HALOGENATED N.O.S. Benzyl chloride Bromoacetone PO17 HALOGENATED N.O.S. U225 HALOGENATED SOLVENTS Bromoform Bromomethane U029 HALOGENATED N.O.S. Brucine PO18 NONHALOG. PESTICIDES Cacodylic acid U136 NONHALOG. PESTICIDES Carbon disulfide PO22 NONHALOGENATED SOLVENTS Carbon tetrachloride U211 HALOGENATED SOLVENTS Carbonyl fluoride U033 HALOGENATED N.O.S. U035 HALOGENATED N.O.S. Chlorambucil Chlordane Pesticides U036 HALOGENATED PESTICIDES Chloroacetaldehyde PO23 HALOGENATED N.O.S. Chlorobenzene U037 HALOGENATED SOLVENTS Chlorobenzilate pesticides U038 HALOGENATED PESTICIDES Chlorodibromomethane UO40 HALOGENATED N.O.S. Chloroform UO44 HALOGENATED SOLVENTS Chloromethane UO45 HALOGENATED N.O.S. Chloromethyl methyl ether UO46 HALOGENATED N.O.S. Chrysene U050 POLYNUCLEAR AROMATICS Creosote U051 PHENOLICS Creosote U051 PHENOLICS Crotonal dehyde U053 POLYMERIZABLE U055 OXYGENATED Cumene Cyanogen PO31 OXYGENATED PO33 OXYGENATED Cyanogen chloride Cyclohexane U056 NONHALOGENATED SOLVENTS

UO57 NONHALOGENATED SOLVENTS Cyclohexanone Cyclophosphamide UO58 NONHALOG. PESTICIDES Daunomycin U059 PHARMACEUTICALS Diallate pesticides U062 HALOGENATED PESTICIDES U063 POLYNUCLEAR AROMATICS Dibenz[a,h]anthracene U064 POLYNUCLEAR AROMATICS Dibenzo(a, i) pyrene UO68 HALOGENATED N.O.S. Dibromomethane U075 HALOGENATED SOLVENTS Dichlorodifluoromethane Dieldrin Pesticides PO37 HALDGENATED PESTICIDES Diepoxybutane UO85 OXYGENATED Diethyl p-nitrophenyl phosphate PO41 NONHALOG. PESTICIDES Diethyl phthalate **U088 PHTHALATES** Diethylhydrazine UO86 REACTIVE (NON-CYANIDE) UO89 PHARMACEUTICALS Diethylstilbestrol U090 PHARMACEUTICALS Dihydrosafrole Diisopropyl fluorophosphate PO43 NONHALOG. PESTICIDES Dimethoate PO44 NONHALOG. PESTICIDES Dimethyl phthalate U102 PHTHALATES U103 REACTIVE (NON-CYANIDE) Dimethyl sulfate U092 ORGANO-NITROGENS Dimethylamine **U097 HALOGENATED PESTICIDES** Dimethylcarbamoyl chloride pesticides Dinoseb PO20 NONHALOG. PESTICIDES PO39 NONHALOG. PESTICIDES Disulfoton Endrin Pesticides PO51 HALOGENATED PESTICIDES Epinephrine PO42 PHARMACEUTICALS Ethyl acetate U112 NONHALOGENATED SOLVENTS Ethyl acrylate U113 POLYMERIZABLE U117 NONHALOGENATED SOLVENTS Ethyl ether Ethyl methacrylate U118 POLYMERIZABLE U119 ORGANO-SULFURS Ethyl methanesulfonate Ethylene bis-dithiocarbamate U114 NONHALOG. PESTICIDES PO53 NONHALOGENATED SOLVENTS Ethylene diamine Ethylene oxide U115 POLYMERIZABLE Ethylene thioruea U116 ORGANO-SULFURS Ethylenimine PO54 POLYMERIZABLE Famphur PO97 NONHALOG. PESTICIDES Fluoranthene U120 POLYNUCLEAR AROMATICS Fluoroacetamide pesticides PO57 HALOGENATED PESTICIDES Fluorotrichloromethane U121 HALDGENATED SOLVENTS Formal dehyde U122 OXYGENATED Formic acid U123 OXYGENATED Furan U124 NONHALOGENATED SOLVENTS Furfural U125 NONHALOGENATED SOLVENTS Glycidaldehyde U126 OXYGENATED Heptachlor Pesticides PO59 HALOGENATED PESTICIDES Hexachlorobenzene U127 HALOGENATED PESTICIDES Hexachlorobutadiene Pesticides U128 HALOGENATED PESTICIDES Hexachlorocyclopentadiene Pesticides U130 HALOGENATED PESTICIDES Hexachlorophene pesticides U132 HALOGENATED PESTICIDES Hexachloropropene pesticides U243 HALOGENATED PESTICIDES Hydrazine U133 REACTIVE (NON-CYANIDE) Hydroxycoumarin salts POO1 NONHALOG. PESTICIDES Indeno (1,2,3-c,d) pyrene U137 POLYNUCLEAR AROMATICS Indomethacin PO25 HALOGENATED N.O.S. Isobutanol U140 NONHALOGENATED SOLVENTS Isocyanic acid methyl ester PO64 DXYGENATED PO60 HALOGENATED PESTICIDES Isodrin pesticides Isosafrole U141 PHARMACEUTICALS Kepone Pesticides U142 HALOGENATED PESTICIDES Lasiocarpine U143 PHARMACEUTICALS

Lindane pesticides U129 HALOGENATED PESTICIDES Maleic anhydride U147 POLYMERIZABLE Malononitrile U149 POLYMERIZABLE U150 HALOGENATED N.O.S. Melphalan U152 POLYMERIZABLE Methacrylonitrile U154 NONHALOGENATED SOLVENTS Methanol PO66 NONHALDG. PESTICIDES Methomy I U247 HALOGENATED PESTICIDES Methoxychlor Methyl Chlorocarbonate U156 HALOGENATED N.O.S. U159 NONHALOGENATED SOLVENTS Methyl ethyl ketone U160 REACTIVE (NON-CYANIDE) Methyl ethyl ketone peroxide Methyl isobutyl ketone U161 NONHALOGENATED SOLVENTS Methyl mercaptan U153 ORGAND-SULFURS Methyl methacrylate U162 POLYMERIZABLE Methyl parathion PO71 NONHALOG. PESTICIDES Methylene chloride UO80 HALDGENATED N.O.S. PO68 REACTIVE (NON-CYANIDE) Methylhydrazine Methylthiouracil **U164 PHARMACEUTICALS** Mitomycin C U010 NONHALOG. PESTICIDES N, N-Bis (2-chloroethyl) -2-naphthylamine UO26 HALOGENATED N.O.S. N-Dipropylamine U110 ORGAND-NITROGENS N-Methyl-N'-nitro-N-nitrosoguanidine U163 PHARMACEUTICALS U176 ORGANO-NITROGENS N-Nitroso-N-ethylurea U177 ORGANO-NITROGENS N-Nitroso-N-methylurea N-Nitroso-N-methylurethane U178 ORGANO-NITROGENS N-Nitroso-di-n-butylamine U172 ORGANO-NITROGENS N-Nitroso-methylvinylamine PO84 ORGANO-NITROGENS U174 ORGANO-NITROGENS N-Nitrosodiethylamine N-Nitrosodimethylamine PO82 ORGAND-NITROGENS U173 ORGANO-NITROGENS N-Nitrosoethanolamine U179 ORGANO-NITROGENS N-Nitrosopiperidine N-Nitrosopyrrolidine U180 ORGANO-NITROGENS N-Phenylthiourea PO93 NONHALOG. PESTICIDES N-nitroso-diphenylamine PO83 ORGANO-NITROGENS Na Fluoracetic acid pesticides PO58 HALOGENATED PESTICIDES Naphthalene U165 POLYNUCLEAR AROMATICS Nicotine PO75 NONHALOG. PESTICIDES U169 ORGANO-NITROGENS Nitrobenzene Nitroglycerine PO81 REACTIVE (NON-CYANIDE) U087 NONHALOG. PESTICIDES 0,0-Diethyl 0,0-Diethyl P PO40 NONHALOG. PESTICIDES Octamethylpyrophosphoramide PO85 NONHALOG. PESTICIDES PO46 ORGANO-NITROGENS P-alpha-Dimethylphenethylamine Paraldehyde U182 OXYGENATED Parathion PO89 NONHALOG. PESTICIDES Pentachlorobenzene U183 HALOGENATED N.O.S. Pentachloroethane U184 HALOGENATED N.O.S. U185 HALOGENATED PESTICIDES Pentachloronitrobenzene pesticides U242 HALOGENATED PESTICIDES Pentachlorophenol Pesticides Phenacetin U187 PHARMACEUTICALS Phenol **U188 PHENOLICS** Phorate PO94 NONHALOG. PESTICIDES Phosgene PO95 HALOGENATED N.O.S. Phthalic anhydride **U190 PHTHALATES** Pronamide U192 ORGANO-NITROGENS Propanenitrile P101 DXYGENATED P102 POLYMERIZABLE Propargyl alcohol Propylene glycol P100 NONHALOGENATED SOLVENTS Pyridine U196 NONHALOGENATED SOLVENTS Reserpine U200 PHARMACEUTICALS

Resorcinol U201 PHENOLICS Saccharin U202 PHARMACEUTICALS Safrole U203 PHARMACEUTICALS Silvex pesticides U233 HALOGENATED PESTICIDES Streptozotocin U206 PHARMACEUTICALS Strychnine and salts P108 NONHALOG. PESTICIDES Tetrachloroethylene U210 HALOGENATED SOLVENTS Tetraethy I dithiopyrophosphate P109 NONHALOG. PESTICIDES Tetraethy I pyrophosphate P111 NONHALOG, PESTICIDES Tetrahydrofuran U213 NONHALOGENATED SOLVENTS P112 REACTIVE (NON-CYANIDE) Tetranitromethane Thioacetamide U218 ORGANO-SULFURS P014 ORGANO-SULFURS Thiophenol Thiosemicarbazide P116 DYES Thiourea U219 ORGANO-SULFURS Thiram P117 NONHALOG. PESTICIDES Toluene U220 NONHALOGENATED SOLVENTS Toluene diamine U221 POLYMERIZABLE U223 POLYMERIZABLE Toluene diisocyanate Toxaphene Pesticides P123 HALOGENATED PESTICIDES Trichloroacetaldehyde U034 HALOGENATED N.O.S. Trichloroethylene U228 HALOGENATED SOLVENTS Trichloromethanethiol P118 HALOGENATED N.O.S. Trypan blue U236 DYES Uracil mustard U237 HALDGENATED N.O.S. Urethane U238 POLYMERIZABLE Vinvl chloride UO43 HALOGENATED N.O.S. alpha, alpha-Dimethylbenzyl hydroperoxid UO96 REACTIVE (NON-CYANIDE) alpha-Endosulfan Pesticides POSO HALOGENATED PESTICIDES bis(2-chloroethyl)ether UO25 HALOGENATED N.O.S. bis(Chloromethyl) ether PO16 HALOGENATED SOLVENTS bis-(2-chloroethyoxy) methane UO24 HALOGENATED N.O.S. bis-(2-chloroisopropyl)ether UO27 HALOGENATED N.O.S. bis-(2-ethylhexyl) phthalate U028 PHTHALATES cis-1,2-Dichloroethylene U079 HALOGENATED SOLVENTS d-n-Octyl phthalate U107 PHTHALATES di-n-Butyl phthalate U069 PHTHALATES di-n-PropyInitrosamine U111 ORGANO-NITROGENS n-Butyl alcohol U031 NONHALOGENATED SOLVENTS n-Propylamine U194 ORGANO-NITROGENS o-Toluidine hydrochloride U222 OXYGENATED o-Xylene U239 NONHALOGENATED SOLVENTS p-Chloro-m-cresol pesticides U039 HALOGENATED PESTICIDES p-Chloroaniline PO24 HALOGENATED N.O.S. p-Cresol U052 PHENOLICS p-Dimethylaminoazobenzene U093 DYES p-Nitroaniline P077 DYES U197 OXYGENATED p-Quinone sym-Trinitrobenzene U234 REACTIVE (NON-CYANIDE) trans-1,2-Dichloroethylene U079 HALDGENATED SOLVENTS

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