

Hazardous Waste Section – Compliance Branch
Soil Cleanup Goals (SCG) for Hazardous Waste
Generators

October 2024

(based on May 2024 USEPA Regional Screening Tables)

***These notes must be used with the Hazardous Waste Section – Compliance Branch
SCG Table***

General Notes:

1. For clean closure, hazardous waste generators must remove ALL wastes and meet the LOWEST VALUE of: (1) the level most protective of human health through direct contact (residential level; unrestricted-use level) and (2) the soil screening level protective of groundwater. Before soil that contains hazardous waste is excavated for purposes of treatment, management, and/or subsequent disposal, Branch staff should be contacted for assistance to address Land Disposal Restrictions found in 40 CFR Part 268.
2. SCGs are obtained using the USEPA Regional Screening Level (RSL) table. If a contaminant does not have a SCG listed, then one or more of the contaminant-specific parameters are not available from the RSL table to calculate a soil SCG (these contaminants are only included in the SCG table to account for all entries on the RSL table). If a 02L Standard or IMAC is available for a contaminant with no SCG, protection of groundwater evaluation is still necessary. Please contact Branch staff for further guidance if a protection of groundwater evaluation is required. If no SCG is available and a protection of groundwater evaluation is not possible for a contaminant of concern, the Method Detection Limit (MDL) or a calculated Practical Quantitation Limit (PQL) will be used as the unrestricted-use-cleanup level.
3. Bold contaminants have SCGs below applicable LDRs found in 40 CFR Part 268. Please contact Branch staff prior to managing these contaminants of concern for further guidance before establishing unrestricted-use-cleanup levels.
4. The **health-based SCGs** (Residential; unrestricted-use) are based upon human health risk and do not address potential ecological risk. The SCGs listed are the lower of:
 - a. the carcinogenic target risk of 1.0E-06 (C), or
 - b. the non-carcinogenic target hazard quotient of 0.2 (N).

Residential health-based SCGs are unrestricted-use goals and must be met by hazardous waste generators for clean-closure.

5. The **protection of groundwater SCGs** are provided as a conservative indicator of soil leachability and are developed using a USEPA soil leaching model with conservative assumptions and default values appropriate for North Carolina (see **Equation 1** on page 4). The target groundwater concentration used in the equation is either:
 - a. **2L or IMAC** – a North Carolina 15A NCAC 02L Groundwater Quality Standard (2L) or an Interim Maximum Allowable Concentration (IMAC),

- b. **Calculated** – For contaminants with no 02L Standard or IMAC, a target groundwater concentration is calculated using the procedures and references in 15A NCAC 02L.0202 (see **Equation 2** on page 5). The calculated groundwater concentration is then used in **Equation 1**. Check with the appropriate remedial program guidance to ensure applicability of a calculated groundwater concentration prior to use. *Use of these calculated soil-to-groundwater SCGs in no way exempts compliance from the 15A NCAC 02L Standards or IMACs, including the use of the Practical Quantitation Limit (PQL) where no numeric standard currently exists. Contact Branch staff for further assistance if no standard exists.*
6. If more than five contaminants with non-carcinogenic effects are detected at a site, including those where a lab reporting limit is greater than a SCG, contact Branch staff so the SCG can be adjusted.

Compound-Specific Notes:

- Cadmium:** Compare soil results with the Cadmium (Diet) SCG.
- Chromium:** USEPA does not provide an RSL for Total Chromium (Cr). Non-speciated Total Cr results must be compared to the Cr (VI) SCG. Speciated results for Cr (III) and Cr (VI) should be compared to the corresponding SCG. Cr levels that have been demonstrated to be the result of natural background concentrations do not require screening.
- Cyanide:** Analyze for total cyanide and compare the results to the lowest SCG for a cyanide species on the table (unless analyses for both total cyanide and specific species are available to demonstrate all of the cyanide is of one particular species).
- Dioxins and Furans:** Multiply the individual dioxin and furan isomer concentrations by their associated toxicity equivalency factor (TEF) shown in the table below. Sum the adjusted concentrations of all dioxins and furans and compare the result to the 2,3,7,8-Tetrachlorodibenzo-p-dioxin (2,3,7,8-TCDD) SCG.

| CASRN | Dioxins and Furans | TEF |
|--------------------------------------|---------------------|--------|
| Chlorinated dibenzo-p-dioxins | | |
| 1746-01-6 | 2,3,7,8-TCDD | 1 |
| 40321-76-4 | 1,2,3,7,8-PeCDD | 1 |
| 39227-28-6 | 1,2,3,4,7,8-HxCDD | 0.1 |
| 57653-85-7 | 1,2,3,6,7,8-HxCDD | 0.1 |
| 57653-85-7 | 1,2,3,7,8,9-HxCDD | 0.1 |
| 35822-46-9 | 1,2,3,4,6,7,8-HpCDD | 0.01 |
| 3268-87-9 | OCDD | 0.0003 |
| CASRN | Dioxins and Furans | TEF |
| Chlorinated dibenzofurans | | |
| 51207-31-9 | 2,3,7,8-TCDF | 0.1 |
| 57117-41-6 | 1,2,3,7,8-PeCDF | 0.03 |
| 57117-31-4 | 2,3,4,7,8-PeCDF | 0.3 |
| 70648-26-9 | 1,2,3,4,7,8-HxCDF | 0.1 |
| 57117-44-9 | 1,2,3,6,7,8-HxCDF | 0.1 |
| 72918-21-9 | 1,2,3,7,8,9-HxCDF | 0.1 |

| | | |
|------------|---------------------|--------|
| 60851-34-5 | 2,3,4,6,7,8-HxCDF | 0.1 |
| 35822-46-9 | 1,2,3,4,6,7,8-HpCDF | 0.01 |
| 55673-89-7 | 1,2,3,4,7,8,9-HpCDF | 0.01 |
| 39001-02-0 | OCDF | 0.0003 |

5. Lead Compounds:

- a. Residential health-based: The SCG is the EPA RSL of 200 mg/kg (updated 1/17/24). However, where additional sources of lead are present (e.g., lead water service lines, lead-based paint and/or non-attainment areas where the air lead concentrations exceed the national ambient air quality standards set by the EPA), the SCG is 100 mg/kg.
- b. Industrial/commercial health-based: The SCG is the EPA RSL of 800 mg/kg.

6. **Manganese:** USEPA provides RSLs for both diet and non-diet. Use the non-diet SCG for comparison, which excludes the dietary contribution from the normal US diet.

7. **Mercury:** Unless it is known what species of mercury is present at the site, use the ‘Mercuric Chloride (and other Mercury salts)’ SCG. If methyl mercury formation or mercury vapor inhalation is likely, contact DEQ for further guidance. .

8. **Nickel:** Unless the release occurred in the last six months and it is clearly known which species of this metal was released, the SCG for "soluble salt" should be used.

9. **PCBs:** SCGs are provided for the 12 dioxin-like PCB congeners. Total the remaining 197 congener concentrations and compare the sum to the “PCBs (high risk)” SCG. Aroclor SCGs should only be used for gross screening and with approval from the appropriate remediation program within DEQ.

10. **Per- and Polyfluoroalkyl Substances (PFAS):** The SCG table contains several species of individual PFAS compounds. The SCg values are typically the same for the salt, acid, and anionic species of each compound. Use the SCGs provided for the acid species (the one with an acronym in parentheses). Although environmental contamination is typically associated with anion species, the acid species is used because the acid species has had the most reference value information. An example of the species associated with a single PFAS compound is shown below with the preferred SCG form highlighted in blue. Protection of groundwater SCGs are not provided for PFAS. If PFAS are detected in soil, co-located groundwater should be tested for PFAS to determine if the protection of groundwater criterium is met or if nearby water supplies could be at risk

Example showing the PFAS species for PFBA

| | CAS No. | Analyte | Chemical Form |
|------|------------|---------------------------------|----------------|
| PFBA | 10495-86-0 | ~Ammonium perfluorobutanoate | Ammonium Salt |
| | 2966-54-3 | ~Potassium heptafluorobutanoate | Potassium Salt |
| | 2218-54-4 | ~Sodium perfluorobutanoate | Sodium Salt |
| | 45048-62-2 | ~Perfluorobutanoate | Anion |
| | 375-22-4 | ~Perfluorobutanoic acid (PFBA) | Acid |

11. **Phosphates:** Unless it is known which species is present, use the “Potassium Salts of Inorganic Phosphates” SCG. This includes laboratory results that report Phosphorus with a CAS of 7723-14-0.
12. **Thallium:** Unless it is clearly known which species of this metal was released, the SCG for "soluble salt" should be used.
13. **Total Petroleum Hydrocarbons (TPH):** TPH is a term intended to refer to the total mass of hydrocarbons present without identifying individual compounds. Because TPH is not a consistent entity, the assessment of health effects and development of toxicity values for mixtures of hydrocarbons are problematic. Therefore, the individual chemical constituents should be analyzed for risk assessment rather than rely on TPH data. In fact, most of the carcinogens in the TPH carbon range are individually listed on the SCG table. Combining TPH and individual constituent cancer risks would be overly protective.
- To better understand the TPH screening levels in the SCG Table, the associated carbon ranges and 2L Standards are provided below. More information on USEPA’s RSLs for TPH can be found here: [Regional Screening Levels \(RSLs\) - Frequent Questions | US EPA](#)
- Protection of groundwater SCGGs are only provided for the two TPH carbon ranges that have both an EPA RSL and a 2L Standard, Aliphatic Low and Aliphatic Medium. See table below for a comparison of the carbon ranges.

| EPA Designation for TPH | Associated Carbon Range (per EPA RSL FAQs) | Associated Carbon Range with NC 2L Groundwater Standard Calculation | NC 2L Groundwater Standard (mg/L) |
|-------------------------|--|---|-----------------------------------|
| Aliphatic Low | C5-C8 | C5-C8 | 0.4 |
| Aliphatic Medium | C9-C18 | C9-C18 | 0.7 |
| Aliphatic High | C19-C32 | C19-C36 | 10 |
| Aromatic Low | C6-C8 | NA | NA |
| Aromatic Medium | C9-C10 | C9-C22 | 0.2 |
| Aromatic High | C10-C32 | NA | NA |

NA - not available

Please contact Branch staff if you have a contaminant not listed on the SCG Table.

Equations Used to Calculate the Protection of Groundwater SCG

Equation 1 calculates a target soil concentration that serves as a leachability screening level. (from USEPA soil to groundwater RSL calculation):

$$C_{soil} = C_{gw} \left[k_s + \frac{(\theta_w + \theta_a H')}{P_b} \right] df$$

| | Parameters | Default Values | Units |
|------------|---|--|---------------------------------------|
| C_{soil} | Calculated Source Concentration for soil | not applicable | mg/kg - soil |
| C_{gw} | Applicable Groundwater Target Concentration: 15A NCAC 02L Standard | 02L standard or IMACs | mg/L - water |
| df | Dilution factor | 20 (0.5 acre source area) ¹ | unitless |
| k_s | Soil-water partition coefficient for organic constituents $k_s = k_{oc} \times f_{oc}$ for inorganic constituents $k_s = k_d$ | Calculated from k_{oc} and f_{oc} , or K_d | L/kg |
| k_{oc} | Soil organic carbon-water partition coefficient | contaminant-specific ² | L/kg |
| f_{oc} | Fraction of organic carbon in subsurface vadose soils | 0.002 (0.2%) ³ | kg/kg |
| k_d | Soil-water partition coefficient for inorganics | contaminant-specific ² (pH=5.5) | L/kg |
| θ_w | Water-filled soil porosity-vadose soils | 0.3 ³ | L _{water} /L _{soil} |
| θ_a | Air-filled soil porosity-vadose soils | 0.13 ³ | L _{air} /L _{soil} |
| P_b | Dry bulk density | 1.5 ³ | kg/L |
| H' | Henry's Law constant-dimensionless where: $H' = \text{Henry's Law constant (atm- m}^3/\text{mole)} \times \text{conversion factor of 41}$ | contaminant-specific ² | unitless |

1 - USEPA default value from USEPA 1996 Soil Screening Guidance

2 - USEPA value from the RS contaminant-specific parameter supporting table:

<https://semspub.epa.gov/work/HQ/199954.pdf>

3 - DEQ default value appropriate for North Carolina

Equation 2 calculates a target groundwater concentration that can be used in Equation 1 when no 2L standard or IMAC is available (based on 15A NCAC 02L .0202 methodology). C_{gw} in Equation 1 becomes the lower of the following calculated groundwater concentrations:

$$C_{gw}(C) = (TCR \times BW) / (CSFo \times IR)$$

$$C_{gw}(N) = (THQ \times RfDo \times BW \times RSC) / (IR)$$

| Symbol | Parameter | Default Values | Units |
|-------------|---|--|---------------------------|
| $C_{gw}(C)$ | Calculated groundwater concentration based on carcinogenic risk | Calculated | mg/L |
| $C_{gw}(N)$ | Calculated groundwater concentration based on non-carcinogenic risk | Calculated | mg/L |
| TCR | Target cancer risk | 1.0E-06 | unitless |
| THQ | Target hazard quotient | 1.0 | unitless |
| RfDo | Oral reference dose | Contaminant-specific | mg/kg-day |
| CSFo | Oral cancer Slope Factor | Contaminant-specific | (mg/kg-day) ⁻¹ |
| BW | Body weight | 70 | kg |
| IR | Ingestion Rate | 2 | L/day |
| RSC | Relative source contribution | 0.1 for inorganics and 0.2 for organics | unitless |

North Carolina Department of Environmental Quality

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(based on May 2024 USEPA Regional Screening Tables)

This table must be used with the SCG Notes/ See Compound-Specific Notes for chemicals in *blue*.

| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|---|--|--------------|---|--------------|--|
| 30560-19-1 | Acephate | 3.8E+00 | N | 9.2E-03 | Calculated | Yes |
| 75-07-0 | Acetaldehyde | 1.2E+01 | C | | | Yes |
| 34256-82-1 | Acetochlor | 2.5E+02 | N | 1.6E+00 | 2L or IMAC | Yes |
| 67-64-1 | Acetone | 1.4E+04 | N | 2.5E+01 | 2L or IMAC | Yes |
| 75-86-5 | Acetone Cyanohydrin | 2.5E+07 | N | | | Yes |
| 75-05-8 | Acetonitrile | 1.7E+02 | N | | | Yes |
| 98-86-2 | Acetophenone | 1.6E+03 | N | 4.3E+00 | 2L or IMAC | Yes |
| 53-96-3 | Acetylaminofluorene, 2- | 1.4E-01 | C | 8.5E-04 | Calculated | |
| 107-02-8 | Acrolein | 3.1E-02 | N | 1.6E-02 | 2L or IMAC | Yes |
| 79-06-1 | Acrylamide | 2.4E-01 | C | 3.4E-05 | 2L or IMAC | Yes |
| 79-10-7 | Acrylic Acid | 4.2E+00 | N | 1.4E+01 | Calculated | Yes |
| 107-13-1 | Acrylonitrile | 2.7E-01 | C | 2.8E-04 | Calculated | Yes |
| 111-69-3 | Adiponitrile | 7.4E+07 | N | | | Yes |
| 15972-60-8 | Alachlor | 9.7E+00 | C | 3.3E-02 | 2L or IMAC | Yes |
| 116-06-3 | Aldicarb | 1.3E+01 | N | 3.5E-02 | Calculated | Yes |
| 1646-88-4 | Aldicarb Sulfone | 1.3E+01 | N | 3.1E-02 | Calculated | Yes |
| 1646-87-3 | Aldicarb sulfoxide | | | | | |
| 309-00-2 | Aldrin | 3.9E-02 | C | 6.6E-03 | 2L or IMAC | Yes |
| 107-18-6 | Allyl Alcohol | 7.5E-01 | N | 1.1E-01 | Calculated | Yes |
| 107-05-1 | Allyl Chloride | 3.5E-01 | N | 1.1E-02 | Calculated | Yes |
| 7429-90-5 | Aluminum | 1.6E+04 | N | 1.1E+05 | Calculated | Yes |
| 20859-73-8 | Aluminum Phosphide | 6.3E+00 | N | | | Yes |
| 834-12-8 | Ametryn | 1.1E+02 | N | 1.3E+00 | Calculated | Yes |
| 92-67-1 | Aminobiphenyl, 4- | 2.6E-02 | C | 1.7E-04 | Calculated | |
| 591-27-5 | Aminophenol, m- | 1.0E+03 | N | 4.3E+00 | Calculated | Yes |
| 95-55-6 | Aminophenol, o- | 5.1E+01 | N | 2.2E-01 | Calculated | Yes |
| 123-30-8 | Aminophenol, p- | 2.5E+02 | N | 1.1E+00 | Calculated | Yes |
| 33089-61-1 | Amitraz | 3.2E+01 | N | 1.8E+02 | Calculated | Yes |
| 7664-41-7 | Ammonia | | | | | |
| 131-74-8 | Ammonium Picrate | 2.5E+01 | N | 1.3E+00 | Calculated | Yes |
| 7773-06-0 | Ammonium Sulfamate | 3.1E+03 | N | | | Yes |
| 75-85-4 | Amyl Alcohol, tert- | 1.7E+01 | N | | | Yes |
| 62-53-3 | Aniline | 8.8E+01 | N | 4.2E-02 | Calculated | Yes |
| 84-65-1 | Anthraquinone, 9,10- | 1.4E+01 | C | 1.8E-01 | Calculated | Yes |
| 7440-36-0 | Antimony (metallic) | 6.3E+00 | N | 9.0E-01 | 2L or IMAC | Yes |
| 1314-60-9 | Antimony Pentoxide | 7.8E+00 | N | | | Yes |
| 1332-81-6 | Antimony Tetroxide | 6.3E+00 | N | | | Yes |
| 1309-64-4 | Antimony Trioxide | 2.5E+06 | N | | | Yes |
| 7440-38-2 | Arsenic, Inorganic | 6.8E-01 | C | 5.8E+00 | 2L or IMAC | Yes |
| 7784-42-1 | Arsine | 5.5E-02 | N | | | Yes |
| 1332-21-4 | Asbestos (units in fibers) | | | | | |
| 3337-71-1 | Asulam | 4.6E+03 | N | 1.3E+01 | Calculated | Yes |
| 1912-24-9 | Atrazine | 2.4E+00 | C | 3.9E-02 | 2L or IMAC | Yes |
| 492-80-8 | Auramine | 6.2E-01 | C | 7.2E-03 | Calculated | |
| 65195-55-3 | Avermectin B1 | 5.1E+00 | N | 9.8E+01 | Calculated | Yes |
| 86-50-0 | Azinphos-methyl | 3.8E+01 | N | 1.3E-01 | Calculated | Yes |

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This table must be used with the SCG Notes/ See Compound-Specific Notes for chemicals in *blue*.

| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|---|--|--------------|---|--------------|--|
| 103-33-3 | Azobenzene | 5.6E+00 | C | 4.9E-02 | Calculated | |
| 123-77-3 | Azodicarbonamide | 1.1E+04 | N | 4.7E+01 | Calculated | Yes |
| 7440-39-3 | Barium | 3.1E+03 | N | 5.8E+02 | 2L or IMAC | Yes |
| 1861-40-1 | Benfluralin | 7.8E+01 | N | 2.3E+01 | Calculated | Yes |
| 17804-35-2 | Benomyl | 6.3E+02 | N | 6.1E+00 | Calculated | Yes |
| 83055-99-6 | Bensulfuron-methyl | 2.5E+03 | N | 7.2E+00 | Calculated | Yes |
| 25057-89-0 | Bentazon | 3.8E+02 | N | 9.2E-01 | Calculated | Yes |
| 100-52-7 | Benzaldehyde | 1.7E+02 | C | 3.1E+00 | 2L or IMAC | Yes |
| 71-43-2 | Benzene | 1.2E+00 | C | 1.0E-02 | 2L or IMAC | Yes |
| 25551-13-7 | Benzene, Trimethyl | 1.1E+01 | N | | | Yes |
| 6369-59-1 | Benzenediamine-2-methyl sulfate, 1,4- | 3.8E+00 | N | 1.9E-03 | Calculated | Yes |
| 108-98-5 | Benzenethiol | 1.6E+01 | N | 9.4E-02 | Calculated | Yes |
| 92-87-5 | Benzidine | 5.3E-04 | C | 7.9E-06 | Calculated | Yes |
| 65-85-0 | Benzoic Acid | 5.1E+04 | N | 1.2E+02 | 2L or IMAC | Yes |
| 98-07-7 | Benzotrichloride | 5.3E-02 | C | 1.2E-04 | Calculated | |
| 100-51-6 | Benzyl Alcohol | 1.3E+03 | N | 3.4E+00 | 2L or IMAC | Yes |
| 100-44-7 | Benzyl Chloride | 1.1E+00 | C | 4.5E-03 | Calculated | Yes |
| 7440-41-7 | Beryllium and compounds | 3.1E+01 | N | 6.3E+01 | 2L or IMAC | Yes |
| 42576-02-3 | Bifenox | 1.1E+02 | N | 9.5E+00 | Calculated | Yes |
| 82657-04-3 | Biphenrin | 1.9E+02 | N | 9.5E+03 | Calculated | Yes |
| 92-52-4 | Biphenyl, 1,1'- | 1.0E+01 | N | 8.4E+01 | 2L or IMAC | Yes |
| 108-60-1 | Bis(2-chloro-1-methylethyl) ether | 6.3E+02 | N | 2.1E+00 | Calculated | Yes |
| 111-91-1 | Bis(2-chloroethoxy)methane | 3.8E+01 | N | 9.6E-02 | Calculated | Yes |
| 111-44-4 | Bis(2-chloroethyl)ether | 2.4E-01 | C | 1.6E-04 | 2L or IMAC | |
| 542-88-1 | Bis(chloromethyl)ether | 8.8E-05 | C | 7.5E-07 | Calculated | |
| 80-05-7 | Bisphenol A | 6.3E+02 | N | 5.3E+02 | Calculated | Yes |
| 7440-42-8 | Boron And Borates Only | 3.1E+03 | N | 4.5E+01 | 2L or IMAC | Yes |
| 10294-34-5 | Boron Trichloride | 3.1E+04 | N | | | Yes |
| 7637-07-2 | Boron Trifluoride | 6.3E+02 | N | | | Yes |
| 15541-45-4 | Bromate | 9.9E-01 | C | 7.7E-03 | Calculated | Yes |
| 107-04-0 | Bromo-2-chloroethane, 1- | 7.5E-02 | N | 4.0E-03 | Calculated | Yes |
| 1073-06-9 | Bromo-3-fluorobenzene, 1- | 4.7E+00 | N | 4.0E-02 | Calculated | Yes |
| 460-00-4 | Bromo-4-fluorobenzene, 1- | 4.7E+00 | N | 4.0E-02 | Calculated | Yes |
| 79-08-3 | Bromoacetic acid | | | | | |
| 108-86-1 | Bromobenzene | 5.9E+01 | N | 7.6E-01 | Calculated | Yes |
| 74-97-5 | Bromochloromethane | 3.2E+01 | N | | | Yes |
| 75-27-4 | Bromodichloromethane | 3.1E-01 | C | 3.3E-03 | 2L or IMAC | Yes |
| 75-25-2 | Bromoform | 2.0E+01 | C | 2.1E-02 | 2L or IMAC | Yes |
| 74-83-9 | Bromomethane | 1.4E+00 | N | 5.0E-02 | 2L or IMAC | Yes |
| 2104-96-3 | Bromophos | 7.8E+01 | N | 3.0E+00 | Calculated | Yes |
| 106-94-5 | Bromopropane, 1- | 1.7E+00 | C | | | Yes |
| 1689-84-5 | Bromoxynil | 5.3E+00 | C | 5.8E-03 | Calculated | Yes |
| 1689-99-2 | Bromoxynil Octanoate | 6.7E+00 | C | 5.9E-02 | Calculated | Yes |
| 106-99-0 | Butadiene, 1,3- | 8.0E-02 | C | 6.3E-04 | Calculated | Yes |

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|---|--|--------------|---|--------------|--|
| 71-36-3 | Butanol, N- | 1.6E+03 | N | 2.4E+00 | 2L or IMAC | Yes |
| 75-65-0 | Butyl Alcohol, t- | 1.4E+03 | C | 4.1E-02 | 2L or IMAC | Yes |
| 78-92-2 | Butyl alcohol, sec- | 2.7E+04 | N | 4.1E+01 | 2L or IMAC | Yes |
| 2008-41-5 | Butylate | 7.8E+02 | N | 6.8E+00 | Calculated | Yes |
| 25013-16-5 | Butylated hydroxyanisole | 2.7E+03 | C | 6.6E+00 | Calculated | |
| 128-37-0 | Butylated hydroxytoluene | 1.5E+02 | C | 5.8E+00 | Calculated | Yes |
| 104-51-8 | Butylbenzene, n- | 7.8E+02 | N | 4.5E+00 | 2L or IMAC | Yes |
| 135-98-8 | Butylbenzene, sec- | 1.6E+03 | N | 4.1E+00 | 2L or IMAC | Yes |
| 98-06-6 | Butylbenzene, tert- | 1.6E+03 | N | 3.1E+00 | 2L or IMAC | Yes |
| 75-60-5 | Cacodylic Acid | 2.5E+02 | N | 8.1E-01 | Calculated | Yes |
| 7440-43-9 | Cadmium (Diet) | 1.4E+00 | N | 3.0E+00 | 2L or IMAC | Yes |
| 7440-43-9 | Cadmium (Water) | | | | | |
| 105-60-2 | Caprolactam | 6.3E+03 | N | 2.0E+01 | 2L or IMAC | Yes |
| 2425-06-1 | Captafol | 3.6E+00 | C | 8.2E-03 | Calculated | Yes |
| 133-06-2 | Captan | 2.4E+02 | C | 2.1E-01 | Calculated | Yes |
| 63-25-2 | Carbaryl | 1.3E+03 | N | 1.3E+01 | Calculated | Yes |
| 1563-66-2 | Carbofuran | 6.3E+01 | N | 3.1E-01 | 2L or IMAC | Yes |
| 75-15-0 | Carbon Disulfide | 1.6E+02 | N | 4.1E+00 | 2L or IMAC | Yes |
| 56-23-5 | Carbon Tetrachloride | 6.9E-01 | C | 2.3E-03 | 2L or IMAC | Yes |
| 463-58-1 | Carbonyl Sulfide | 1.4E+01 | N | | | Yes |
| 55285-14-8 | Carbosulfan | 1.3E+02 | N | 3.4E+01 | Calculated | Yes |
| 5234-68-4 | Carboxin | 1.3E+03 | N | 7.5E+00 | Calculated | Yes |
| 1306-38-3 | Ceric oxide | 1.1E+07 | N | | | Yes |
| 302-17-0 | Chloral Hydrate | 1.6E+03 | N | 2.8E+00 | Calculated | Yes |
| 133-90-4 | Chloramben | 1.9E+02 | N | 5.1E-01 | Calculated | Yes |
| E701235 | Chloramines, Organic | | | | | |
| 118-75-2 | Chloranil | 1.3E+00 | C | 1.4E-03 | Calculated | |
| 5103-71-9 | Chlordane (alpha) | 7.1E+00 | N | 9.5E+00 | Calculated | Yes |
| 5103-74-2 | Chlordane (gamma) | 7.1E+00 | N | 9.5E+00 | Calculated | Yes |
| 12789-03-6 | Chlordane (technical mixture) | 1.7E+00 | C | 2.7E-01 | 2L or IMAC | Yes |
| 143-50-0 | Chlordecone (Kepone) | 5.4E-02 | C | 2.5E-03 | Calculated | Yes |
| 470-90-6 | Chlorfenvinphos | 8.8E+00 | N | 2.7E-01 | Calculated | Yes |
| 90982-32-4 | Chlorimuron, Ethyl- | 1.1E+03 | N | 4.3E+00 | Calculated | Yes |
| 7782-50-5 | Chlorine | 3.9E-02 | N | 3.4E+00 | Calculated | Yes |
| 10049-04-4 | Chlorine Dioxide | 4.7E+02 | N | | | Yes |
| 7758-19-2 | Chlorite (Sodium Salt) | 4.7E+02 | N | | | Yes |
| 75-68-3 | Chloro-1,1-difluoroethane, 1- | 1.1E+04 | N | | | Yes |
| 126-99-8 | Chloro-1,3-butadiene, 2- (Chloroprene) | 1.1E-02 | C | 1.5E+00 | Calculated | Yes |
| 3165-93-3 | Chloro-2-methylaniline HCl, 4- | 1.2E+00 | C | 1.4E-03 | Calculated | |
| 95-69-2 | Chloro-2-methylaniline, 4- | 5.4E+00 | C | 4.0E-03 | Calculated | Yes |
| 107-20-0 | Chloroacetaldehyde, 2- | 2.6E+00 | C | 5.2E-04 | Calculated | |
| 79-11-8 | Chloroacetic Acid | | | | | |
| 532-27-4 | Chloroacetophenone, 2- | 3.7E+05 | N | | | Yes |
| 106-47-8 | Chloroaniline, p- | 2.7E+00 | C | 1.5E-03 | Calculated | Yes |
| 108-90-7 | Chlorobenzene | 5.8E+01 | N | 6.8E-01 | 2L or IMAC | Yes |

North Carolina Department of Environmental Quality

Soil Cleanup Goals (SCG)

October 2024

(based on May 2024 USEPA Regional Screening Tables)

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|-------------|---|---------------------------------------|-------|--|------------|---|
| 98-66-8 | Chlorobenzene sulfonic acid, p- | 1.3E+03 | N | 3.2E+00 | Calculated | Yes |
| 510-15-6 | Chlorobenzilate | 4.9E+00 | C | 2.1E-02 | Calculated | Yes |
| 74-11-3 | Chlorobenzoic Acid, p- | 3.8E+02 | N | 1.1E+00 | Calculated | Yes |
| 98-56-6 | Chlorobenzotrifluoride, 4- | 2.4E+00 | C | 1.5E+00 | Calculated | Yes |
| 109-69-3 | Chlorobutane, 1- | 6.3E+02 | N | 2.3E+00 | Calculated | Yes |
| 75-45-6 | Chlorodifluoromethane | 1.0E+04 | N | | | Yes |
| 107-07-3 | Chloroethanol, 2- | 3.1E+02 | N | 5.7E-01 | Calculated | Yes |
| 67-66-3 | Chloroform | 3.4E-01 | C | 3.9E-01 | 2L or IMAC | Yes |
| 74-87-3 | Chloromethane | 2.3E+01 | N | 1.5E-02 | 2L or IMAC | Yes |
| 107-30-2 | Chloromethyl Methyl Ether | 2.1E-02 | C | 6.2E-05 | Calculated | |
| 88-73-3 | Chloronitrobenzene, o- | 1.8E+00 | C | 2.2E-03 | Calculated | Yes |
| 100-00-5 | Chloronitrobenzene, p- | 8.8E+00 | N | 1.1E-02 | Calculated | Yes |
| 95-57-8 | Chlorophenol, 2- | 7.8E+01 | N | 7.8E-03 | 2L or IMAC | Yes |
| 76-06-2 | Chloropicrin | 4.2E-01 | N | | | Yes |
| 1897-45-6 | Chlorothalonil | 3.2E+01 | C | 9.4E-02 | Calculated | Yes |
| 95-49-8 | Chlorotoluene, o- | 3.1E+02 | N | 2.0E+00 | 2L or IMAC | Yes |
| 106-43-4 | Chlorotoluene, p- | 3.1E+02 | N | 4.6E-01 | 2L or IMAC | Yes |
| 54749-90-5 | Chlorozotocin | 2.3E-03 | C | 6.4E-07 | Calculated | |
| 101-21-3 | Chlorpropham | 6.3E+01 | N | 6.3E-01 | Calculated | Yes |
| 2921-88-2 | Chlorpyrifos | 1.3E+01 | N | 2.1E+00 | Calculated | Yes |
| 5598-13-0 | Chlorpyrifos Methyl | 1.3E+02 | N | 6.4E+00 | Calculated | Yes |
| 64902-72-3 | Chlorsulfuron | 6.3E+02 | N | 5.9E+00 | Calculated | Yes |
| 1861-32-1 | Chlorthal-dimethyl | 1.3E+02 | N | 1.7E+00 | Calculated | Yes |
| 60238-56-4 | Chlorthiophos | 1.0E+01 | N | 2.9E+00 | Calculated | Yes |
| 16065-83-1b | Chromium(III) (Soluble Compounds) | 7.4E+05 | N | 3.6E+05 | 2L or IMAC | Yes |
| 16065-83-1 | Chromium(III), Insoluble Salts | 2.3E+04 | N | 3.6E+05 | 2L or IMAC | Yes |
| 18540-29-9 | Chromium(VI) | 3.1E-01 | C | 3.8E+00 | 2L or IMAC | Yes |
| 7440-47-3 | Chromium, Total | | | | | |
| 74115-24-5 | Clofentezine | 1.6E+02 | N | 1.1E+02 | Calculated | Yes |
| 7440-48-4 | Cobalt | 4.7E+00 | N | 9.0E-01 | 2L or IMAC | Yes |
| E649830 | Coke Oven Emissions | | | | | |
| 7440-50-8 | Copper | 6.3E+02 | N | 7.0E+02 | 2L or IMAC | Yes |
| 108-39-4 | Cresol, m- | 6.3E+02 | N | 6.4E+00 | 2L or IMAC | Yes |
| 95-48-7 | Cresol, o- | 6.3E+02 | N | 6.5E+00 | 2L or IMAC | Yes |
| 106-44-5 | Cresol, p- | 2.5E+02 | N | 6.4E-01 | 2L or IMAC | Yes |
| 59-50-7 | Cresol, p-chloro-m- | 1.3E+03 | N | 1.7E+01 | Calculated | Yes |
| 1319-77-3 | Cresols | 1.3E+03 | N | 1.1E+01 | Calculated | Yes |
| 123-73-9 | Crotonaldehyde, trans- | 3.7E-01 | C | 7.5E-05 | Calculated | Yes |
| 98-82-8 | Cumene | 4.1E+02 | N | 2.3E+00 | 2L or IMAC | Yes |
| 135-20-6 | Cupferron | 2.5E+00 | C | 5.5E-03 | Calculated | |
| 21725-46-2 | Cyanazine | 6.5E-01 | C | 3.9E-04 | Calculated | Yes |
| | Cyanides | | | | | |
| 592-01-8 | ~Calcium Cyanide | 1.6E+01 | N | | | Yes |
| 544-92-3 | ~Copper Cyanide | 7.8E+01 | N | | | Yes |

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|------------|--|---------------------------------------|-------|--|------------|---|
| 57-12-5 | ~Cyanide (CN-) | 4.7E+00 | N | 1.4E+01 | 2L or IMAC | Yes |
| 460-19-5 | ~Cyanogen | 1.6E+01 | N | | | Yes |
| 506-68-3 | ~Cyanogen Bromide | 1.4E+03 | N | | | Yes |
| 506-77-4 | ~Cyanogen Chloride | 7.8E+02 | N | | | Yes |
| 74-90-8 | ~Hydrogen Cyanide | 4.7E+00 | N | 4.2E-01 | Calculated | Yes |
| 151-50-8 | ~Potassium Cyanide | 3.1E+01 | N | | | Yes |
| 506-61-6 | ~Potassium Silver Cyanide | 7.8E+01 | N | | | Yes |
| 506-64-9 | ~Silver Cyanide | 1.6E+03 | N | | | Yes |
| 143-33-9 | ~Sodium Cyanide | 1.6E+01 | N | | | Yes |
| E1790665 | ~Thiocyanates | 3.1E+00 | N | | | Yes |
| 463-56-9 | ~Thiocyanic Acid | 3.1E+00 | N | | | Yes |
| 557-21-1 | ~Zinc Cyanide | 7.8E+02 | N | | | Yes |
| 110-82-7 | Cyclohexane | 1.4E+03 | N | | | Yes |
| 87-84-3 | Cyclohexane, 1,2,3,4,5-pentabromo-6-chloro- | 2.7E+01 | C | 2.0E-01 | Calculated | Yes |
| 108-94-1 | Cyclohexanone | 6.0E+03 | N | 1.6E+02 | Calculated | Yes |
| 110-83-8 | Cyclohexene | 6.3E+01 | N | 4.6E-01 | Calculated | Yes |
| 108-91-8 | Cyclohexylamine | 3.1E+03 | N | 7.4E+00 | Calculated | Yes |
| 68359-37-5 | Cyfluthrin | 3.2E+02 | N | 9.1E+02 | Calculated | Yes |
| 66215-27-8 | Cyromazine | 6.3E+03 | N | 1.8E+01 | Calculated | Yes |
| 75-99-0 | Dalapon | 3.8E+02 | N | 8.3E-01 | 2L or IMAC | Yes |
| 1596-84-5 | Daminozide | 3.0E+01 | C | 8.6E-03 | Calculated | Yes |
| 1163-19-5 | Decabromodiphenyl ether, 2,2',3,3',4,4',5,5',6,6'- (BDE-209) | 8.8E+01 | N | 5.4E+02 | Calculated | Yes |
| 8065-48-3 | Demeton | 5.1E-01 | N | | | Yes |
| 103-23-1 | Di(2-ethylhexyl)adipate | 4.5E+02 | C | 4.2E+01 | Calculated | Yes |
| 2303-16-4 | Diallate | 8.9E+00 | C | 1.7E-02 | Calculated | |
| 333-41-5 | Diazinon | 8.8E+00 | N | 6.1E-01 | Calculated | Yes |
| 96-12-8 | Dibromo-3-chloropropane, 1,2- | 5.6E-03 | C | 3.5E-04 | 2L or IMAC | Yes |
| 631-64-1 | Dibromoacetic acid | | | | | |
| 108-36-1 | Dibromobenzene, 1,3- | 6.3E+00 | N | 5.3E-02 | Calculated | Yes |
| 106-37-6 | Dibromobenzene, 1,4- | 1.6E+02 | N | 1.3E+00 | 2L or IMAC | Yes |
| 124-48-1 | Dibromochloromethane | 8.3E+00 | C | 2.1E-03 | 2L or IMAC | Yes |
| 106-93-4 | Dibromoethane, 1,2- | 3.8E-02 | C | 1.1E-04 | 2L or IMAC | Yes |
| 74-95-3 | Dibromomethane (Methylene Bromide) | 5.0E+00 | N | 3.4E-01 | 2L or IMAC | Yes |
| E1790661 | Dibutyltin Compounds | 3.8E+00 | N | | | Yes |
| 1918-00-9 | Dicamba | 3.8E+02 | N | 1.1E+00 | Calculated | Yes |
| 3400-09-7 | Dichloramine | | | | | |
| 764-41-0 | Dichloro-2-butene, 1,4- | 2.3E-03 | C | | | |
| 1476-11-5 | Dichloro-2-butene, cis-1,4- | 7.9E-03 | C | | | |
| 110-57-6 | Dichloro-2-butene, trans-1,4- | 7.9E-03 | C | | | |
| 79-43-6 | Dichloroacetic Acid | 1.1E+01 | C | 2.9E-03 | 2L or IMAC | Yes |
| 95-50-1 | Dichlorobenzene, 1,2- | 3.8E+02 | N | 3.9E-01 | 2L or IMAC | Yes |
| 106-46-7 | Dichlorobenzene, 1,4- | 2.8E+00 | C | 1.2E-01 | 2L or IMAC | Yes |

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|--------------|---|--|--------------|---|--------------|--|
| 91-94-1 | Dichlorobenzidine, 3,3'- | 1.2E+00 | C | 1.0E-02 | Calculated | |
| 90-98-2 | Dichlorobenzophenone, 4,4'- | 1.1E+02 | N | 7.6E+00 | Calculated | Yes |
| 75-71-8 | Dichlorodifluoromethane | 1.8E+01 | N | 3.0E+01 | 2L or IMAC | Yes |
| 72-54-8 | Dichlorodiphenyldichloroethane, p,p'- (DDD) | 2.3E+00 | C | 4.7E-01 | 2L or IMAC | Yes |
| 72-55-9 | Dichlorodiphenyldichloroethylene, p,p'- (DDE) | 2.0E+00 | C | 4.7E-01 | 2L or IMAC | Yes |
| 50-29-3 | Dichlorodiphenyltrichloroethane, p,p'- (DDT) | 1.9E+00 | C | 6.7E-01 | 2L or IMAC | Yes |
| 75-34-3 | Dichloroethane, 1,1- | 3.8E+00 | C | 3.4E-02 | 2L or IMAC | Yes |
| 107-06-2 | Dichloroethane, 1,2- | 4.9E-01 | C | 2.3E-03 | 2L or IMAC | Yes |
| 75-35-4 | Dichloroethylene, 1,1- | 4.8E+01 | N | 2.5E+00 | 2L or IMAC | Yes |
| 156-59-2 | Dichloroethylene, cis-1,2- | 1.3E+01 | N | 4.1E-01 | 2L or IMAC | Yes |
| 156-60-5 | Dichloroethylene, trans-1,2- | 1.5E+01 | N | 6.2E-01 | 2L or IMAC | Yes |
| 120-83-2 | Dichlorophenol, 2,4- | 3.8E+01 | N | 9.7E-03 | 2L or IMAC | Yes |
| 94-75-7 | Dichlorophenoxy Acetic Acid, 2,4- | 1.4E+02 | N | 3.6E-01 | 2L or IMAC | Yes |
| 78-87-5 | Dichloropropane, 1,2- | 2.6E+00 | C | 4.0E-03 | 2L or IMAC | Yes |
| 142-28-9 | Dichloropropane, 1,3- | 3.1E+02 | N | 9.7E-01 | Calculated | Yes |
| 616-23-9 | Dichloropropanol, 2,3- | 3.8E+01 | N | 8.9E-02 | Calculated | Yes |
| 542-75-6 | Dichloropropene, 1,3- | 1.9E+00 | C | 2.9E-03 | 2L or IMAC | Yes |
| 62-73-7 | Dichlorvos | 1.9E+00 | C | 7.4E-04 | Calculated | Yes |
| 141-66-2 | Dicrotophos | 3.8E-01 | N | 9.8E-04 | Calculated | Yes |
| 77-73-6 | Dicyclopentadiene | 2.7E-01 | N | 3.9E+01 | Calculated | Yes |
| 60-57-1 | Dieldrin | 3.4E-02 | C | 1.6E-03 | 2L or IMAC | Yes |
| E17136615 | Diesel Engine Exhaust | | | | | |
| 111-42-2 | Diethanolamine | 2.5E+01 | N | 5.7E-02 | Calculated | Yes |
| 112-34-5 | Diethylene Glycol Monobutyl Ether | 3.8E+02 | N | 9.2E-01 | Calculated | Yes |
| 111-90-0 | Diethylene Glycol Monoethyl Ether | 7.6E+02 | N | 1.7E+00 | Calculated | Yes |
| 617-84-5 | Diethylformamide | 1.6E+01 | N | 2.9E-02 | Calculated | Yes |
| 56-53-1 | Diethylstilbestrol | 1.6E-03 | C | 1.1E-03 | Calculated | |
| 43222-48-6 | Difenzoquat | 1.0E+03 | N | 1.8E+03 | Calculated | Yes |
| 35367-38-5 | Diflubenzuron | 2.5E+02 | N | 3.2E+00 | Calculated | Yes |
| 75-37-6 | Difluoroethane, 1,1- | 1.0E+04 | N | | | Yes |
| 420-45-1 | Difluoropropane, 2,2- | 5.0E+03 | N | | | Yes |
| 94-58-6 | Dihydrosafrole | 1.0E+01 | C | 9.8E-03 | Calculated | |
| 108-20-3 | Diisopropyl Ether | 4.8E+02 | N | 3.6E-01 | 2L or IMAC | Yes |
| 1445-75-6 | Diisopropyl Methylphosphonate | 1.3E+03 | N | 3.2E+00 | Calculated | Yes |
| 55290-64-7 | Dimethipin | 2.8E+02 | N | 6.7E-01 | Calculated | Yes |
| 60-51-5 | Dimethoate | 2.8E+01 | N | 6.9E-02 | Calculated | Yes |
| 119-90-4 | Dimethoxybenzidine, 3,3'- | 3.4E-01 | C | 5.3E-04 | Calculated | |
| 756-79-6 | Dimethyl methylphosphonate | 3.2E+02 | C | 8.7E-02 | Calculated | Yes |

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|------------|---|---------------------------------------|-------|--|------------|---|
| 60-11-7 | Dimethylamino azobenzene [p-] | 1.2E-01 | C | 6.5E-04 | Calculated | |
| 21436-96-4 | Dimethylaniline HCl, 2,4- | 9.4E-01 | C | 1.1E-03 | Calculated | |
| 95-68-1 | Dimethylaniline, 2,4- | 2.7E+00 | C | 2.0E-03 | Calculated | Yes |
| 121-69-7 | Dimethylaniline, N,N- | 2.6E+01 | C | 9.3E-03 | Calculated | Yes |
| 119-93-7 | Dimethylbenzidine, 3,3'- | 4.9E-02 | C | 4.2E-04 | Calculated | |
| 68-12-2 | Dimethylformamide | 5.4E+02 | N | 2.8E+00 | Calculated | Yes |
| 57-14-7 | Dimethylhydrazine, 1,1- | 1.2E-02 | N | 3.1E-03 | Calculated | Yes |
| 540-73-8 | Dimethylhydrazine, 1,2- | 9.0E-04 | C | 2.9E-07 | Calculated | |
| 105-67-9 | Dimethylphenol, 2,4- | 2.5E+02 | N | 2.4E+00 | 2L or IMAC | Yes |
| 576-26-1 | Dimethylphenol, 2,6- | 7.6E+00 | N | 1.0E-01 | Calculated | Yes |
| 95-65-8 | Dimethylphenol, 3,4- | 1.3E+01 | N | 1.7E-01 | Calculated | Yes |
| 513-37-1 | Dimethylvinylchloride | 1.2E+00 | C | 5.1E-03 | Calculated | |
| 534-52-1 | Dinitro-o-cresol, 4,6- | 1.0E+00 | N | 1.9E-02 | Calculated | Yes |
| 131-89-5 | Dinitro-o-cyclohexyl Phenol, 4,6- | 2.5E+01 | N | 9.3E+00 | Calculated | Yes |
| 618-87-1 | Dinitroaniline, 3,5- | 5.1E+00 | N | 3.0E-02 | Calculated | Yes |
| 528-29-0 | Dinitrobenzene, 1,2- | 1.3E+00 | N | 1.3E-02 | Calculated | Yes |
| 99-65-0 | Dinitrobenzene, 1,3- | 1.3E+00 | N | 1.3E-02 | Calculated | Yes |
| 100-25-4 | Dinitrobenzene, 1,4- | 1.3E+00 | N | 1.3E-02 | Calculated | Yes |
| 51-28-5 | Dinitrophenol, 2,4- | 2.5E+01 | N | 3.1E-01 | Calculated | Yes |
| E1615210 | Dinitrotoluene Mixture, 2,4/2,6- | 8.0E-01 | C | 1.4E-03 | Calculated | |
| 121-14-2 | Dinitrotoluene, 2,4- | 1.7E+00 | C | 1.4E-03 | 2L or IMAC | Yes |
| 606-20-2 | Dinitrotoluene, 2,6- | 3.6E-01 | C | 1.4E-03 | 2L or IMAC | Yes |
| 35572-78-2 | Dinitrotoluene, 2-Amino-4,6- | 1.5E+00 | N | 1.1E-02 | Calculated | Yes |
| 19406-51-0 | Dinitrotoluene, 4-Amino-2,6- | 1.5E+00 | N | 1.1E-02 | Calculated | Yes |
| 25321-14-6 | Dinitrotoluene, Technical grade | 1.2E+00 | C | 2.1E-03 | Calculated | Yes |
| 88-85-7 | Dinoseb | 1.3E+01 | N | 1.2E+00 | 2L or IMAC | Yes |
| 123-91-1 | Dioxane, 1,4- | 5.4E+00 | C | 1.2E-02 | 2L or IMAC | Yes |
| | Dioxins | | | | | |
| 34465-46-8 | ~Hexachlorodibenzo-p-dioxin, Mixture | 1.0E-04 | C | 1.6E-04 | Calculated | |
| 1746-01-6 | ~TCDD, 2,3,7,8- | 4.8E-06 | C | 2.0E-06 | 2L or IMAC | Yes |
| 957-51-7 | Diphenamid | 3.8E+02 | N | 4.1E+01 | Calculated | Yes |
| 101-84-8 | Diphenyl Ether | 7.2E+00 | N | 1.5E+01 | 2L or IMAC | Yes |
| 127-63-9 | Diphenyl Sulfone | 1.0E+01 | N | 2.7E-01 | Calculated | Yes |
| 122-39-4 | Diphenylamine | 1.3E+03 | N | 2.6E+01 | Calculated | Yes |
| 122-66-7 | Diphenylhydrazine, 1,2- | 6.8E-01 | C | 2.8E-03 | Calculated | |
| 2764-72-9 | Diquat | 2.8E+01 | N | 3.3E+00 | 2L or IMAC | Yes |
| 1937-37-7 | Direct Black 38 | 7.3E-02 | C | 4.6E+01 | Calculated | |
| 2602-46-2 | Direct Blue 6 | 7.3E-02 | C | 1.5E+02 | Calculated | |
| 16071-86-6 | Direct Brown 95 | 8.1E-02 | C | 1.5E+00 | Calculated | |
| 298-04-4 | Disulfoton | 5.1E-01 | N | 1.1E-02 | 2L or IMAC | Yes |
| 505-29-3 | Dithiane, 1,4- | 1.6E+02 | N | 6.9E-01 | Calculated | Yes |
| 330-54-1 | Diuron | 2.5E+01 | N | 1.2E-01 | Calculated | Yes |

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|--------------|---|--|--------------|---|--------------|--|
| 2439-10-3 | Dodine | 2.5E+02 | N | 1.4E+01 | Calculated | Yes |
| 759-94-4 | EPTC | 7.8E+02 | N | 3.7E+00 | Calculated | Yes |
| 115-29-7 | Endosulfan | 9.4E+01 | N | 1.1E+01 | 2L or IMAC | Yes |
| 1031-07-8 | Endosulfan Sulfate | 7.6E+01 | N | 1.6E+01 | 2L or IMAC | Yes |
| 145-73-3 | Endothall | 2.5E+02 | N | 4.8E-01 | 2L or IMAC | Yes |
| 72-20-8 | Endrin | 3.8E+00 | N | 1.6E+00 | 2L or IMAC | Yes |
| 106-89-8 | Epichlorohydrin | 4.0E+00 | N | 1.8E-02 | 2L or IMAC | Yes |
| 106-88-7 | Epoxybutane, 1,2- | 3.4E+01 | N | | | Yes |
| 111-77-3 | Ethanol, 2-(2-methoxyethoxy)- | 5.1E+02 | N | 1.1E+00 | Calculated | Yes |
| 16672-87-0 | Ethephon | 6.3E+01 | N | 1.5E-01 | Calculated | Yes |
| 563-12-2 | Ethion | 6.3E+00 | N | 1.4E-01 | Calculated | Yes |
| 111-15-9 | Ethoxyethanol Acetate, 2- | 5.4E+02 | N | 2.9E+00 | Calculated | Yes |
| 110-80-5 | Ethoxyethanol, 2- | 5.4E+02 | N | 2.5E+00 | Calculated | Yes |
| 141-78-6 | Ethyl Acetate | 1.3E+02 | N | 1.3E+01 | 2L or IMAC | Yes |
| 140-88-5 | Ethyl Acrylate | 9.8E+00 | N | 1.6E-01 | Calculated | Yes |
| 75-00-3 | Ethyl Chloride (Chloroethane) | 1.1E+03 | N | 1.7E+01 | 2L or IMAC | Yes |
| 60-29-7 | Ethyl Ether | 3.1E+03 | N | 6.3E+00 | Calculated | Yes |
| 97-63-2 | Ethyl Methacrylate | 3.8E+02 | N | | | Yes |
| 637-92-3 | Ethyl Tertiary Butyl Ether (ETBE) | 1.4E+02 | C | 2.3E-01 | 2L or IMAC | Yes |
| 2104-64-5 | Ethyl-p-nitrophenyl Phosphonate | 1.3E-01 | N | 4.4E-02 | Calculated | Yes |
| 100-41-4 | Ethylbenzene | 6.1E+00 | C | 1.3E+01 | 2L or IMAC | Yes |
| 109-78-4 | Ethylene Cyanohydrin | 8.8E+02 | N | 2.0E+00 | Calculated | Yes |
| 107-15-3 | Ethylene Diamine | 1.4E+03 | N | 2.9E+00 | Calculated | Yes |
| 107-21-1 | Ethylene Glycol | 1.0E+04 | N | 4.0E+01 | 2L or IMAC | Yes |
| 111-76-2 | Ethylene Glycol Monobutyl Ether | 1.3E+03 | N | 2.9E+00 | Calculated | Yes |
| 75-21-8 | Ethylene Oxide | 2.2E-03 | C | 4.7E-04 | Calculated | Yes |
| 96-45-7 | Ethylene Thiourea | 1.0E+00 | N | 2.5E-03 | Calculated | Yes |
| 151-56-4 | Ethyleneimine | 2.8E-03 | C | 2.3E-06 | Calculated | |
| 84-72-0 | Ethylphthalyl Ethyl Glycolate | 3.8E+04 | N | 9.4E+02 | Calculated | Yes |
| 22224-92-6 | Fenamiphos | 3.2E+00 | N | 3.5E-02 | Calculated | Yes |
| 39515-41-8 | Fenpropathrin | 3.2E+02 | N | 1.6E+02 | Calculated | Yes |
| 51630-58-1 | Fenvalerate | 3.2E+02 | N | 2.2E+03 | Calculated | Yes |
| 2164-17-2 | Fluometuron | 1.6E+02 | N | 1.4E+00 | Calculated | Yes |
| 16984-48-8 | Fluoride | 6.3E+02 | N | 6.0E+03 | 2L or IMAC | Yes |
| 7782-41-4 | Fluorine (Soluble Fluoride) | 9.4E+02 | N | 6.3E+02 | Calculated | Yes |
| 59756-60-4 | Fluridone | 1.0E+03 | N | 1.3E+03 | Calculated | Yes |
| 56425-91-3 | Flurprimidol | 5.1E+02 | N | 2.6E+01 | Calculated | Yes |
| 85509-19-9 | Flusilazole | 2.5E+01 | N | 4.5E+01 | Calculated | Yes |
| 66332-96-5 | Flutolanil | 6.3E+03 | N | 3.7E+02 | Calculated | Yes |
| 69409-94-5 | Fluvalinate | 1.3E+02 | N | 2.0E+03 | Calculated | Yes |
| 133-07-3 | Folpet | 1.1E+03 | N | 3.0E+00 | Calculated | Yes |
| 72178-02-0 | Fomesafen | 1.3E+02 | N | 4.6E+00 | Calculated | Yes |
| 944-22-9 | Fonofos | 2.5E+01 | N | 5.4E-01 | Calculated | Yes |

North Carolina Department of Environmental Quality

Soil Cleanup Goals (SCG)

October 2024

(based on May 2024 USEPA Regional Screening Tables)

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|------------|---|---------------------------------------|-------|--|------------|---|
| 50-00-0 | Formaldehyde | 1.2E+01 | C | 2.4E+00 | 2L or IMAC | Yes |
| 64-18-6 | Formic Acid | 6.1E+00 | N | 2.5E+01 | Calculated | Yes |
| 39148-24-8 | Fosetyl-AL | 3.2E+04 | N | 4.6E+03 | Calculated | Yes |
| | Furans | | | | | |
| 132-64-9 | ~Dibenzofuran | 1.6E+01 | N | 1.0E+01 | 2L or IMAC | Yes |
| 110-00-9 | ~Furan | 1.6E+01 | N | 5.3E-02 | Calculated | Yes |
| 109-99-9 | ~Tetrahydrofuran | 3.9E+03 | N | 8.9E+00 | 2L or IMAC | Yes |
| 67-45-8 | Furazolidone | 1.4E-01 | C | 3.5E-04 | Calculated | |
| 98-01-1 | Furfural | 4.3E+01 | N | 8.9E-02 | Calculated | Yes |
| 531-82-8 | Furium | 3.6E-01 | C | 6.3E-04 | Calculated | |
| 60568-05-0 | Furmecyclox | 1.8E+01 | C | 2.5E-02 | Calculated | |
| 77182-82-2 | Glufosinate, Ammonium | 7.6E+01 | N | 1.8E-01 | Calculated | Yes |
| 111-30-8 | Glutaraldehyde | 1.3E+03 | N | 2.8E+00 | Calculated | Yes |
| 765-34-4 | Glycidaldehyde | 4.7E+00 | N | 1.1E-02 | Calculated | Yes |
| 1071-83-6 | Glyphosate | 1.3E+03 | N | 6.2E+01 | Calculated | Yes |
| 113-00-8 | Guanidine | 1.6E+02 | N | 3.1E-01 | Calculated | Yes |
| 50-01-1 | Guanidine Chloride | 2.5E+02 | N | | | Yes |
| 506-93-4 | Guanidine Nitrate | 3.8E+02 | N | 1.0E+00 | Calculated | Yes |
| 69806-40-2 | Haloxypop, Methyl | 6.3E-01 | N | 7.8E-02 | Calculated | Yes |
| 76-44-8 | Heptachlor | 1.4E-01 | C | 1.3E-02 | 2L or IMAC | Yes |
| 1024-57-3 | Heptachlor Epoxide | 7.1E-02 | C | 1.6E-03 | 2L or IMAC | Yes |
| 111-71-7 | Heptanal, n- | 5.2E+00 | N | | | Yes |
| 142-82-5 | Heptane, N- | 4.4E+00 | N | 6.2E+01 | 2L or IMAC | Yes |
| 87-82-1 | Hexabromobenzene | 3.1E+01 | N | 1.6E+00 | Calculated | Yes |
| 68631-49-2 | Hexabromodiphenyl ether, 2,2',4,4',5,5'- (BDE-153) | 2.5E+00 | N | | | Yes |
| 118-74-1 | Hexachlorobenzene | 1.6E-01 | N | 5.0E-03 | 2L or IMAC | Yes |
| 87-68-3 | Hexachlorobutadiene | 1.3E+00 | C | 1.5E-02 | 2L or IMAC | Yes |
| 319-84-6 | Hexachlorocyclohexane, Alpha- | 8.6E-02 | C | 7.0E-04 | 2L or IMAC | Yes |
| 319-85-7 | Hexachlorocyclohexane, Beta- | 3.0E-01 | C | 2.3E-03 | 2L or IMAC | |
| 58-89-9 | Hexachlorocyclohexane, Gamma- (Lindane) | 5.7E-01 | C | 3.5E-03 | 2L or IMAC | Yes |
| 608-73-1 | Hexachlorocyclohexane, Technical | 3.0E-01 | C | 2.3E-03 | 2L or IMAC | |
| 77-47-4 | Hexachlorocyclopentadiene | 3.8E-01 | N | 2.6E+00 | Calculated | Yes |
| 67-72-1 | Hexachloroethane | 1.9E+00 | C | 1.1E-02 | Calculated | Yes |
| 70-30-4 | Hexachlorophene | 3.8E+00 | N | 5.6E+01 | Calculated | Yes |
| 121-82-4 | Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX) | 8.3E+00 | C | 3.3E-03 | Calculated | Yes |
| 822-06-0 | Hexamethylene Diisocyanate, 1,6- | 6.7E-01 | N | | | Yes |
| 4035-89-6 | Hexamethylene diisocyanate biuret | 4.9E+06 | N | | | Yes |
| 3779-63-3 | Hexamethylene diisocyanate isocyanurate | 4.9E+06 | N | | | Yes |
| 680-31-9 | Hexamethylphosphoramide | 5.1E+00 | N | 1.2E-02 | Calculated | Yes |

North Carolina Department of Environmental Quality

Soil Cleanup Goals (SCG)

October 2024

(based on May 2024 USEPA Regional Screening Tables)

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|-------------|--|---------------------------------------|-------|--|------------|---|
| E5241997 | Hexane, Commercial | 1.2E+01 | C | | | Yes |
| 110-54-3 | Hexane, N- | 1.3E+02 | N | 5.5E+01 | 2L or IMAC | Yes |
| 124-04-9 | Hexanedioic Acid | 2.5E+04 | N | 7.0E+01 | Calculated | Yes |
| 104-76-7 | Hexanol, 1-,2-ethyl- (2-Ethyl-1-hexanol) | 3.2E+00 | N | 2.0E-02 | Calculated | Yes |
| 591-78-6 | Hexanone, 2- | 4.2E+01 | N | 1.8E-01 | 2L or IMAC | Yes |
| 51235-04-2 | Hexazinone | 4.2E+02 | N | 2.1E+00 | Calculated | Yes |
| 78587-05-0 | Hexythiazox | 3.2E+02 | N | 1.6E+01 | Calculated | Yes |
| 67485-29-4 | Hydramethylnon | 2.1E+02 | N | 8.6E+05 | Calculated | Yes |
| 302-01-2 | Hydrazine | 3.4E-02 | C | 4.8E-05 | Calculated | Yes |
| 10034-93-2 | Hydrazine Sulfate | 2.3E-01 | C | | | |
| 7647-01-0 | Hydrogen Chloride | 2.5E+08 | N | | | Yes |
| 7664-39-3 | Hydrogen Fluoride | 6.3E+02 | N | | | Yes |
| 7783-06-4 | Hydrogen Sulfide | 2.5E+07 | N | | | Yes |
| 123-31-9 | Hydroquinone | 9.0E+00 | C | 7.9E-03 | Calculated | Yes |
| 35554-44-0 | Imazalil | 8.9E+00 | C | 2.0E-01 | Calculated | Yes |
| 81335-37-7 | Imazaquin | 3.2E+03 | N | 1.7E+02 | Calculated | Yes |
| 81335-77-5 | Imazethapyr | 3.2E+04 | N | 3.1E+02 | Calculated | Yes |
| 7553-56-2 | Iodine | 1.6E+02 | N | 4.2E+01 | Calculated | Yes |
| 36734-19-7 | Iprodione | 5.1E+02 | N | 1.7E+00 | Calculated | Yes |
| 7439-89-6 | Iron | 1.1E+04 | N | 1.5E+02 | 2L or IMAC | Yes |
| 78-83-1 | Isobutyl Alcohol | 1.6E+03 | N | 8.6E+00 | Calculated | Yes |
| 78-59-1 | Isophorone | 5.7E+02 | C | 2.6E-01 | 2L or IMAC | Yes |
| 33820-53-0 | Isopropalin | 2.3E+02 | N | 4.8E+01 | Calculated | Yes |
| 67-63-0 | Isopropanol | 1.2E+03 | N | 5.7E+01 | Calculated | Yes |
| 1832-54-8 | Isopropyl Methyl Phosphonic Acid | 1.3E+03 | N | 3.0E+00 | Calculated | Yes |
| 82558-50-7 | Isoxaben | 6.3E+02 | N | 1.9E+01 | Calculated | Yes |
| E1737665 | Jet propulsion fuel 7 (JP-7) | 3.7E+09 | N | | | Yes |
| 77501-63-4 | Lactofen | 1.0E+02 | N | 5.2E+01 | Calculated | Yes |
| 78-97-7 | Lactonitrile | 2.5E+00 | N | 5.7E-03 | Calculated | Yes |
| 7439-91-0 | Lanthanum | 7.8E-01 | N | | | Yes |
| 100587-90-4 | Lanthanum Acetate Hydrate | 2.6E-01 | N | | | Yes |
| 10025-84-0 | Lanthanum Chloride Heptahydrate | 2.9E-01 | N | | | Yes |
| 10099-58-8 | Lanthanum Chloride, Anhydrous | 4.4E-01 | N | | | Yes |
| 10277-43-7 | Lanthanum Nitrate Hexahydrate | 2.5E-01 | N | | | Yes |
| | Lead Compounds | | | | | |
| 7446-27-7 | ~Lead Phosphate | 8.2E+01 | C | | | |
| 301-04-2 | ~Lead acetate | 2.6E+00 | C | 6.7E-04 | Calculated | |
| 7439-92-1 | ~Lead and Compounds | 2.0E+02 | | 2.7E+02 | 2L or IMAC | |
| 7439-92-1b | ~Lead and Compounds (with other sources of lead present, see Guidance) | 1.0E+02 | | 2.7E+02 | 2L or IMAC | |

North Carolina Department of Environmental Quality

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|------------|---|---------------------------------------|-------|--|------------|---|
| 1335-32-6 | ~Lead subacetate | 1.4E+01 | C | 4.1E-03 | Calculated | |
| 78-00-2 | ~Tetraethyl Lead | 1.6E-03 | N | 4.9E-05 | Calculated | Yes |
| 541-25-3 | Lewisite | 7.8E-02 | N | 3.0E-04 | Calculated | Yes |
| 330-55-2 | Linuron | 9.7E+01 | N | 9.5E-01 | Calculated | Yes |
| 7439-93-2 | Lithium | 3.1E+01 | N | 4.2E+01 | Calculated | Yes |
| 94-74-6 | MCPA | 6.3E+00 | N | 1.8E-02 | Calculated | Yes |
| 94-81-5 | MCPB | 5.6E+02 | N | 2.4E+00 | Calculated | Yes |
| 93-65-2 | MCPP | 1.3E+01 | N | 4.2E-02 | Calculated | Yes |
| 121-75-5 | Malathion | 2.5E+02 | N | 7.4E-01 | Calculated | Yes |
| 108-31-6 | Maleic Anhydride | 1.3E+03 | N | 2.8E+00 | Calculated | Yes |
| 123-33-1 | Maleic Hydrazide | 6.3E+03 | N | 1.4E+01 | Calculated | Yes |
| 109-77-3 | Malononitrile | 1.3E+00 | N | 2.9E-03 | Calculated | Yes |
| 8018-01-7 | Mancozeb | 3.8E+02 | N | 5.9E+00 | Calculated | Yes |
| 12427-38-2 | Maneb | 6.3E+01 | N | 9.9E-01 | Calculated | Yes |
| 7439-96-5 | Manganese (Diet) | | | | | |
| 7439-96-5 | Manganese (Non-diet) | 3.8E+02 | N | 6.5E+01 | 2L or IMAC | Yes |
| 950-10-7 | Mephosfolan | 1.1E+00 | N | 1.9E-02 | Calculated | Yes |
| 24307-26-4 | Mepiquat Chloride | 3.8E+02 | N | 1.4E+00 | Calculated | Yes |
| 149-30-4 | Mercaptobenzothiazole, 2- | 4.9E+01 | C | 1.9E-01 | Calculated | Yes |
| | Mercury Compounds | | | | | |
| 7487-94-7 | ~Mercuric Chloride (and other Mercury salts) | 4.7E+00 | N | | | Yes |
| 7439-97-6 | ~Mercury (elemental) | 2.3E+00 | N | 1.0E+00 | 2L or IMAC | Yes |
| 22967-92-6 | ~Methyl Mercury | 1.6E+00 | N | 4.9E+01 | Calculated | Yes |
| 62-38-4 | ~Phenylmercuric Acetate | 1.0E+00 | N | 3.5E-03 | Calculated | Yes |
| 150-50-5 | Merphos | 4.7E-01 | N | 4.1E-01 | Calculated | Yes |
| 57837-19-1 | Metalaxyl | 7.6E+02 | N | 2.3E+00 | Calculated | Yes |
| 126-98-7 | Methacrylonitrile | 1.5E+00 | N | 3.2E-03 | Calculated | Yes |
| 10265-92-6 | Methamidophos | 6.3E-01 | N | 1.5E-03 | Calculated | Yes |
| 67-56-1 | Methanol | 2.5E+04 | N | 1.6E+01 | 2L or IMAC | Yes |
| 950-37-8 | Methidathion | 1.9E+01 | N | 5.1E-02 | Calculated | Yes |
| 16752-77-5 | Methomyl | 3.2E+02 | N | 7.7E-01 | Calculated | Yes |
| 99-59-2 | Methoxy-5-nitroaniline, 2- | 1.1E+01 | C | 4.9E-03 | Calculated | |
| 72-43-5 | Methoxychlor | 6.3E+01 | N | 4.3E+01 | 2L or IMAC | Yes |
| 110-49-6 | Methoxyethanol Acetate, 2- | 2.2E+01 | N | 2.3E-01 | Calculated | Yes |
| 109-86-4 | Methoxyethanol, 2- | 5.2E+01 | N | 1.4E-01 | Calculated | Yes |
| 79-20-9 | Methyl Acetate | 1.6E+04 | N | 2.9E+01 | Calculated | Yes |
| 96-33-3 | Methyl Acrylate | 3.1E+01 | N | | | Yes |
| 78-93-3 | Methyl Ethyl Ketone (2-Butanone) | 5.5E+03 | N | 1.7E+01 | 2L or IMAC | Yes |
| 60-34-4 | Methyl Hydrazine | 1.5E-01 | C | 3.2E-02 | Calculated | Yes |
| 108-10-1 | Methyl Isobutyl Ketone (4-methyl-2-pentanone) | 7.0E+03 | N | 4.5E-01 | 2L or IMAC | Yes |
| 624-83-9 | Methyl Isocyanate | 9.8E-01 | N | | | Yes |
| 80-62-6 | Methyl Methacrylate | 9.4E+02 | N | 1.1E-01 | 2L or IMAC | Yes |
| 298-00-0 | Methyl Parathion | 3.2E+00 | N | 5.8E-02 | Calculated | Yes |

North Carolina Department of Environmental Quality

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|--------------|---|--|--------------|---|--------------|--|
| 993-13-5 | Methyl Phosphonic Acid | 7.6E+02 | N | 1.7E+00 | Calculated | Yes |
| 25013-15-4 | Methyl Styrene (Mixed Isomers) | 6.5E+01 | N | 1.4E+00 | Calculated | Yes |
| 66-27-3 | Methyl methanesulfonate | 5.5E+00 | C | 1.5E-03 | Calculated | |
| 1634-04-4 | Methyl tert-Butyl Ether (MTBE) | 4.9E+01 | C | 9.0E-02 | 2L or IMAC | Yes |
| 615-45-2 | Methyl-1,4-benzenediamine dihydrochloride, 2- | 3.8E+00 | N | 2.5E-02 | Calculated | Yes |
| 108-11-2 | Methyl-2-Pentanol, 4- | 1.1E+04 | N | | | Yes |
| 99-55-8 | Methyl-5-Nitroaniline, 2- | 6.0E+01 | C | 4.3E-02 | Calculated | Yes |
| 70-25-7 | Methyl-N-nitro-N-nitrosoguanidine, N- | 6.5E-02 | C | 2.9E-05 | Calculated | |
| 636-21-5 | Methylaniline Hydrochloride, 2- | 4.2E+00 | C | 2.3E-03 | Calculated | |
| 124-58-3 | Methylarsonic acid | 1.3E+02 | N | 4.0E-01 | Calculated | Yes |
| 74612-12-7 | Methylbenzene, 1-4-diamine monohydrochloride, 2- | 2.5E+00 | N | | | Yes |
| 615-50-9 | Methylbenzene-1,4-diamine sulfate, 2- | 3.8E+00 | N | | | Yes |
| 56-49-5 | Methylcholanthrene, 3- | 5.5E-03 | C | 6.1E-02 | Calculated | |
| 108-87-2 | Methylcyclohexane | 2.1E+01 | N | | | Yes |
| 75-09-2 | Methylene Chloride | 5.8E+01 | C | 2.5E-02 | 2L or IMAC | Yes |
| 101-14-4 | Methylene-bis(2-chloroaniline), 4,4'- | 1.2E+00 | C | 8.1E-02 | Calculated | Yes |
| 101-61-1 | Methylene-bis(N,N-dimethyl) Aniline, 4,4'- | 1.2E+01 | C | 8.4E-02 | Calculated | |
| 101-77-9 | Methylenebisbenzenamine, 4,4'- | 3.4E-01 | C | 1.9E-03 | Calculated | Yes |
| 101-68-8 | Methylenediphenyl Diisocyanate | 7.4E+06 | N | | | Yes |
| 98-83-9 | Methylstyrene, Alpha- | 1.1E+03 | N | 1.6E+01 | Calculated | Yes |
| 51218-45-2 | Metolachlor | 1.9E+03 | N | 2.5E+01 | Calculated | Yes |
| 21087-64-9 | Metribuzin | 3.2E+02 | N | 1.1E+00 | Calculated | Yes |
| 74223-64-6 | Metsulfuron-methyl | 3.2E+03 | N | 1.3E+01 | Calculated | Yes |
| E1790669 | Midrange Aliphatic Hydrocarbon Streams | 6.9E-01 | C | 1.9E+01 | Calculated | Yes |
| 8012-95-1 | Mineral oils | 4.7E+04 | N | 1.6E+04 | Calculated | Yes |
| 2385-85-5 | Mirex | 3.6E-02 | C | 2.8E-02 | Calculated | Yes |
| 2212-67-1 | Molinate | 2.5E+01 | N | 1.6E-01 | Calculated | Yes |
| 7439-98-7 | Molybdenum | 7.8E+01 | N | 7.1E+00 | Calculated | Yes |
| 10599-90-3 | Monochloramine | 1.6E+03 | N | | | Yes |
| 100-61-8 | Monomethylaniline | 2.5E+01 | N | 1.0E-01 | Calculated | Yes |
| 88671-89-0 | Myclobutanil | 3.2E+02 | N | 4.3E+01 | Calculated | Yes |
| 74-31-7 | N,N'-Diphenyl-1,4-benzenediamine | 3.8E+00 | N | 4.4E+00 | Calculated | Yes |
| 300-76-5 | Naled | 3.1E+01 | N | 1.3E-01 | Calculated | Yes |
| 64742-95-6 | Naphtha, High Flash Aromatic (HFAN) | 4.7E+02 | N | | | Yes |

North Carolina Department of Environmental Quality

Soil Cleanup Goals (SCG)

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|------------|---|---------------------------------------|-------|--|------------|---|
| 91-59-8 | Naphthylamine, 2- | 3.0E-01 | C | 2.0E-03 | Calculated | |
| 15299-99-7 | Napropamide | 1.5E+03 | N | 1.1E+02 | Calculated | Yes |
| 373-02-4 | Nickel Acetate | 1.4E+02 | N | 3.1E-01 | Calculated | Yes |
| 3333-67-3 | Nickel Carbonate | 1.4E+02 | N | | | Yes |
| 13463-39-3 | Nickel Carbonyl | 1.7E+02 | N | | | Yes |
| 12054-48-7 | Nickel Hydroxide | 1.7E+02 | N | | | Yes |
| 1313-99-1 | Nickel Oxide | 1.7E+02 | N | | | Yes |
| E715532 | Nickel Refinery Dust | 1.7E+02 | N | 1.2E+02 | Calculated | Yes |
| 7440-02-0 | Nickel Soluble Salts | 3.1E+02 | N | 1.3E+02 | 2L or IMAC | Yes |
| 12035-72-2 | Nickel Subsulfide | 4.1E-01 | C | | | Yes |
| 1271-28-9 | Nickelocene | 6.0E-01 | C | | | Yes |
| 14797-55-8 | Nitrate (measured as nitrogen) | 2.5E+04 | N | | | Yes |
| E701177 | Nitrate + Nitrite (measured as nitrogen) | | | | | |
| 14797-65-0 | Nitrite (measured as nitrogen) | 1.6E+03 | N | | | Yes |
| 88-74-4 | Nitroaniline, 2- | 1.3E+02 | N | 5.9E-01 | Calculated | Yes |
| 100-01-6 | Nitroaniline, 4- | 2.7E+01 | C | 1.5E-02 | Calculated | Yes |
| 98-95-3 | Nitrobenzene | 5.5E+00 | C | 1.8E-01 | Calculated | Yes |
| 9004-70-0 | Nitrocellulose | 3.8E+07 | N | 9.2E+04 | Calculated | Yes |
| 67-20-9 | Nitrofurantoin | 8.8E+02 | N | 4.2E+00 | Calculated | Yes |
| 59-87-0 | Nitrofurazone | 4.2E-01 | C | 4.8E-04 | Calculated | |
| 55-63-0 | Nitroglycerin | 1.3E+00 | N | 6.0E-03 | Calculated | Yes |
| 556-88-7 | Nitroguanidine | 1.3E+03 | N | 3.4E+00 | Calculated | Yes |
| 75-52-5 | Nitromethane | 5.8E+00 | C | | | Yes |
| 79-46-9 | Nitropropane, 2- | 6.8E-02 | C | | | Yes |
| 759-73-9 | Nitroso-N-ethylurea, N- | 4.5E-03 | C | 6.3E-06 | Calculated | |
| 684-93-5 | Nitroso-N-methylurea, N- | 1.0E-03 | C | 1.3E-06 | Calculated | |
| 924-16-3 | Nitrosodibutylamine, N- | 1.0E-01 | C | 2.6E-04 | Calculated | |
| 1116-54-7 | Nitrosodiethanolamine, N- | 1.9E-01 | C | 5.1E-05 | Calculated | |
| 55-18-5 | Nitrosodiethylamine, N- | 8.1E-04 | C | 1.7E-06 | Calculated | |
| 62-75-9 | Nitrosodimethylamine, N- | 2.0E-03 | C | 3.4E-06 | 2L or IMAC | Yes |
| 86-30-6 | Nitrosodiphenylamine, N- | 1.1E+02 | C | 7.8E-01 | Calculated | |
| 621-64-7 | Nitrosodipropylamine, N- | 7.8E-02 | C | 7.5E-05 | Calculated | |
| 10595-95-6 | Nitrosomethylethylamine, N- | 2.0E-02 | C | 9.1E-06 | Calculated | |
| 59-89-2 | Nitrosomorpholine [N-] | 8.1E-02 | C | 2.6E-05 | Calculated | |
| 100-75-4 | Nitrosopiperidine [N-] | 5.8E-02 | C | 4.0E-05 | Calculated | |
| 930-55-2 | Nitrosopyrrolidine, N- | 2.6E-01 | C | 1.3E-04 | Calculated | |
| 99-08-1 | Nitrotoluene, m- | 1.3E+00 | N | 1.3E-02 | Calculated | Yes |
| 88-72-2 | Nitrotoluene, o- | 3.2E+00 | C | 3.0E-03 | Calculated | Yes |
| 99-99-0 | Nitrotoluene, p- | 3.4E+01 | C | 4.1E-02 | Calculated | Yes |
| 111-84-2 | Nonane, n- | 2.3E+00 | N | 5.8E-01 | Calculated | Yes |
| 27314-13-2 | Norflurazon | 1.9E+01 | N | 1.4E+00 | Calculated | Yes |
| 32536-52-0 | Octabromodiphenyl Ether | 3.8E+01 | N | 8.3E+01 | Calculated | Yes |
| 2691-41-0 | Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX) | 7.7E+02 | N | 8.8E+00 | Calculated | Yes |

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(based on May 2024 USEPA Regional Screening Tables)

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|------------|---|---------------------------------------|-------|--|------------|---|
| 152-16-9 | Octamethylpyrophosphoramidate | 2.5E+01 | N | 6.7E-02 | Calculated | Yes |
| 19044-88-3 | Oryzalin | 7.0E+01 | C | 1.7E-01 | Calculated | Yes |
| 19666-30-9 | Oxadiazon | 6.3E+01 | N | 7.1E+00 | Calculated | Yes |
| 23135-22-0 | Oxamyl | 3.2E+02 | N | 8.8E-01 | 2L or IMAC | Yes |
| 42874-03-3 | Oxyfluorfen | 7.4E+00 | C | 7.7E-01 | Calculated | Yes |
| 76738-62-0 | Paclobutrazol | 1.6E+02 | N | 3.7E+00 | Calculated | Yes |
| 1910-42-5 | Paraquat Dichloride | 5.7E+01 | N | 8.7E+00 | Calculated | Yes |
| 56-38-2 | Parathion | 7.6E+01 | N | 4.2E+00 | Calculated | Yes |
| 1114-71-2 | Pebulate | 7.8E+02 | N | 5.6E+00 | Calculated | Yes |
| 40487-42-1 | Pendimethalin | 3.8E+03 | N | 4.8E+02 | Calculated | Yes |
| 32534-81-9 | Pentabromodiphenyl Ether | 3.1E+01 | N | 1.2E+01 | Calculated | Yes |
| 60348-60-9 | Pentabromodiphenyl ether, 2,2',4,4',5- (BDE-99) | 1.3E+00 | N | 6.1E-01 | Calculated | Yes |
| 608-93-5 | Pentachlorobenzene | 1.3E+01 | N | 8.5E-01 | Calculated | Yes |
| 76-01-7 | Pentachloroethane | 7.7E+00 | C | 3.7E-03 | Calculated | |
| 82-68-8 | Pentachloronitrobenzene | 2.7E+00 | C | 3.3E-02 | Calculated | Yes |
| 87-86-5 | Pentachlorophenol | 1.0E+00 | C | 8.3E-03 | 2L or IMAC | Yes |
| 78-11-5 | Pentaerythritol tetranitrate (PETN) | 1.1E+02 | N | 2.4E-01 | Calculated | Yes |
| 10159-46-3 | Pentamethylphosphoramidate (PMPA) | 1.3E+00 | N | 2.9E-03 | Calculated | Yes |
| 109-66-0 | Pentane, n- | 1.7E+02 | N | | | Yes |
| | Per- and Polyfluoroalkyl Substances (PFAS) | | | | | |
| 62037-80-3 | ~Ammonium perfluoro-2-methyl-3-oxahexanoate | 3.8E-02 | N | | | Yes |
| 10495-86-0 | ~Ammonium perfluorobutanoate | 1.6E+01 | N | | | Yes |
| 21615-47-4 | ~Ammonium perfluorohexanoate | 6.3E+00 | N | | | Yes |
| 3825-26-1 | ~Ammonium perfluorooctanoate | 1.9E-05 | C | | | Yes |
| 82113-65-3 | ~Bis(trifluoromethylsulfonyl)amine (TFSI) | 4.7E+00 | N | | | Yes |
| 13252-13-6 | ~Hexafluoropropylene oxide dimer acid (HFPO-DA) | 4.7E-02 | N | | | Yes |
| 90076-65-6 | ~Lithium bis[(trifluoromethyl)sulfonyl]azide | 4.7E+00 | N | | | Yes |
| 45187-15-3 | ~Perfluorobutanesulfonate | 3.8E+00 | N | | | Yes |
| 375-73-5 | ~Perfluorobutanesulfonic acid (PFBS) | 3.8E+00 | N | | | Yes |
| 45048-62-2 | ~Perfluorobutanoate | 1.6E+01 | N | | | Yes |
| 375-22-4 | ~Perfluorobutanoic acid (PFBA) | 1.6E+01 | N | | | Yes |

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|-------------|---|---------------------------------------|-------|--|------------|---|
| 307-55-1 | ~Perfluorododecanoic acid (PFDoDA) | 6.3E-01 | N | | | Yes |
| 108427-53-8 | ~Perfluorohexanesulfonate | 2.5E-01 | N | | | Yes |
| 355-46-4 | ~Perfluorohexanesulfonic acid (PFHxS) | 2.5E-01 | N | | | Yes |
| 92612-52-7 | ~Perfluorohexanoate | 6.3E+00 | N | | | Yes |
| 307-24-4 | ~Perfluorohexanoic acid (PFHxA) | 6.3E+00 | N | | | Yes |
| 72007-68-2 | ~Perfluorononanoate | 3.8E-02 | N | | | Yes |
| 375-95-1 | ~Perfluorononanoic acid (PFNA) | 3.8E-02 | N | | | Yes |
| 16517-11-6 | ~Perfluorooctadecanoic acid (PFODA) | 5.1E+02 | N | | | Yes |
| 45298-90-6 | ~Perfluorooctanesulfonate | 1.3E-03 | N | | | Yes |
| 1763-23-1 | ~Perfluorooctanesulfonic acid (PFOS) | 1.3E-03 | N | | | Yes |
| 45285-51-6 | ~Perfluorooctanoate | 1.9E-05 | C | | | Yes |
| 335-67-1 | ~Perfluorooctanoic acid (PFOA) | 1.9E-05 | C | | | Yes |
| 422-64-0 | ~Perfluoropropanoic acid (PFPrA) | 7.8E+00 | N | | | Yes |
| 376-06-7 | ~Perfluorotetradecanoic acid (PFTetDA) | 1.3E+01 | N | | | Yes |
| 2058-94-8 | ~Perfluoroundecanoic acid (PFUDA) | 3.8E+00 | N | | | Yes |
| 2966-54-3 | ~Potassium heptafluorobutanoate | 3.1E+01 | N | | | Yes |
| 29420-49-3 | ~Potassium perfluorobutanesulfonate | 3.8E+00 | N | | | Yes |
| 2795-39-3 | ~Potassium perfluorooctanesulfonate | 1.3E-03 | N | | | Yes |
| 2218-54-4 | ~Sodium perfluorobutanoate | 1.6E+01 | N | | | Yes |
| 2923-26-4 | ~Sodium perfluorohexanoate | 6.3E+00 | N | | | Yes |
| | Perchlorates | | | | | |
| 7790-98-9 | ~Ammonium Perchlorate | 1.1E+01 | N | | | Yes |
| 7791-03-9 | ~Lithium Perchlorate | 1.1E+01 | N | | | Yes |
| 14797-73-0 | ~Perchlorate and Perchlorate Salts | 1.1E+01 | N | | | Yes |
| 7778-74-7 | ~Potassium Perchlorate | 1.1E+01 | N | | | Yes |
| 7601-89-0 | ~Sodium Perchlorate | 1.1E+01 | N | | | Yes |
| 52645-53-1 | Permethrin | 6.3E+02 | N | 1.7E+03 | Calculated | Yes |
| 62-44-2 | Phenacetin | 2.5E+02 | C | 9.0E-02 | Calculated | |
| 13684-63-4 | Phenmedipham | 3.0E+03 | N | 1.8E+02 | Calculated | Yes |
| 108-95-2 | Phenol | 3.8E+03 | N | 3.4E-01 | 2L or IMAC | Yes |
| 114-26-1 | Phenol, 2-(1-methylethoxy)-, methylcarbamate | 5.1E+01 | N | 1.8E-01 | Calculated | Yes |
| 92-84-2 | Phenothiazine | 6.3E+00 | N | 2.2E-01 | Calculated | Yes |

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|--------------|---|--|--------------|---|--------------|--|
| 103-72-0 | Phenyl Isothiocyanate | 3.1E+00 | N | 1.8E-02 | Calculated | Yes |
| 108-45-2 | Phenylenediamine, m- | 7.6E+01 | N | 2.2E-01 | Calculated | Yes |
| 95-54-5 | Phenylenediamine, o- | 4.5E+00 | C | 1.6E-03 | Calculated | Yes |
| 106-50-3 | Phenylenediamine, p- | 1.3E+01 | N | 3.7E-02 | Calculated | Yes |
| 90-43-7 | Phenylphenol, 2- | 2.8E+02 | C | 4.9E+00 | Calculated | |
| 298-02-2 | Phorate | 2.5E+00 | N | 2.2E-02 | 2L or IMAC | Yes |
| 75-44-5 | Phosgene | 6.5E-02 | N | | | Yes |
| 732-11-6 | Phosmet | 2.5E+02 | N | 6.2E-01 | Calculated | Yes |
| | Phosphates, Inorganic | | | | | |
| 13776-88-0 | ~Aluminum metaphosphate | 4.6E+04 | N | | | Yes |
| E524680405 | ~Aluminum salts of inorganic phosphates | 4.7E+03 | N | | | Yes |
| 7758-11-4 | ~Dipotassium phosphate | 1.6E+04 | N | | | Yes |
| 7558-79-4 | ~Disodium phosphate | 1.6E+04 | N | | | Yes |
| 13530-50-2 | ~Monoaluminum phosphate | 5.5E+04 | N | | | Yes |
| 7778-77-0 | ~Monopotassium phosphate | 1.6E+04 | N | | | Yes |
| 7558-80-7 | ~Monosodium phosphate | 1.6E+04 | N | | | Yes |
| 7784-30-7 | ~Phosphoric acid, aluminum salt (1:1) [aluminum phosphate] | 1.7E+04 | N | | | Yes |
| 7785-88-8 | ~Phosphoric acid, aluminum sodium salt (1:X:X) [sodium aluminum phosphate acidic (acidic SALP)] | 6.7E+04 | N | | | Yes |
| 8017-16-1 | ~Polyphosphoric acid | 1.6E+04 | N | | | Yes |
| E524680403 | ~Potassium salts of inorganic phosphates | 1.6E+04 | N | | | Yes |
| 13845-36-8 | ~Potassium tripolyphosphate | 1.6E+04 | N | | | Yes |
| 10279-59-1 | ~Sodium aluminum phosphate (anhydrous) | 7.8E+04 | N | | | Yes |
| 10305-76-7 | ~Sodium aluminum phosphate (tetrahydrate) | 5.5E+04 | N | | | Yes |
| 10124-56-8 | ~Sodium hexametaphosphate | 1.6E+04 | N | | | Yes |
| 68915-31-1 | ~Sodium polyphosphate | 1.6E+04 | N | | | Yes |
| 7758-16-9 | ~Sodium pyrophosphate | 1.6E+04 | N | | | Yes |
| E524680404 | ~Sodium salts of inorganic phosphates | 1.6E+04 | N | | | Yes |
| 7785-84-4 | ~Sodium trimetaphosphate | 1.6E+04 | N | | | Yes |
| 7758-29-4 | ~Sodium tripolyphosphate | 1.6E+04 | N | | | Yes |
| 7320-34-5 | ~Tetrapotassium phosphate | 1.6E+04 | N | | | Yes |
| 7722-88-5 | ~Tetrasodium pyrophosphate | 1.6E+04 | N | | | Yes |
| 15136-87-5 | ~Trialuminum sodium tetra decahydrogenoctaorthophosphate (dihydrate) | 5.1E+04 | N | | | Yes |
| 13939-25-8 | ~Triphosphoric acid, aluminum salt (1:1) [aluminum triphosphate] | 4.0E+04 | N | | | Yes |

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|--------------|---|--|--------------|---|--------------|--|
| 7778-53-2 | ~Tripotassium phosphate | 1.6E+04 | N | | | Yes |
| 7601-54-9 | ~Trisodium phosphate | 1.6E+04 | N | | | Yes |
| 7803-51-2 | Phosphine | 4.7E+00 | N | | | Yes |
| 7664-38-2 | Phosphoric Acid | 1.6E+04 | N | | | Yes |
| 7723-14-0 | Phosphorus, White | 3.1E-01 | N | 3.4E-03 | Calculated | Yes |
| | Phthalates | | | | | |
| 117-81-7 | ~Bis(2-ethylhexyl)phthalate | 3.9E+01 | C | 1.4E+01 | 2L or IMAC | Yes |
| 85-68-7 | ~Butyl Benzyl Phthalate | 2.9E+02 | C | 2.9E+02 | 2L or IMAC | Yes |
| 85-70-1 | ~Butylphthalyl Butylglycolate | 1.3E+04 | N | 3.2E+03 | Calculated | Yes |
| 84-74-2 | ~Dibutyl Phthalate | 1.3E+03 | N | 3.5E+01 | 2L or IMAC | Yes |
| 84-66-2 | ~Diethyl Phthalate | 1.0E+04 | N | 4.9E+01 | 2L or IMAC | Yes |
| 120-61-6 | ~Dimethylterephthalate | 1.6E+03 | N | 3.7E+00 | Calculated | Yes |
| 117-84-0 | ~Octyl Phthalate, di-N- | 1.3E+02 | N | 5.6E+02 | 2L or IMAC | Yes |
| 100-21-0 | ~Phthalic Acid, p- | 6.3E+03 | N | 2.5E+01 | Calculated | Yes |
| 85-44-9 | ~Phthalic Anhydride | 2.5E+04 | N | 6.2E+01 | Calculated | Yes |
| 1918-02-1 | Picloram | 8.8E+02 | N | 2.7E+00 | Calculated | Yes |
| 96-91-3 | Picramic Acid (2-Amino-4,6-dinitrophenol) | 1.3E+00 | N | 9.1E-03 | 2L or IMAC | Yes |
| 88-89-1 | Picric Acid (2,4,6-Trinitrophenol) | 2.5E+01 | N | 1.3E+00 | Calculated | Yes |
| 29232-93-7 | Pirimiphos, Methyl | 9.2E+00 | N | 9.7E-02 | Calculated | Yes |
| 36355-01-8 | Polybrominated Biphenyls | 1.8E-02 | C | | | Yes |
| | Polychlorinated Biphenyls (PCBs) | | | | | |
| 12674-11-2 | ~Aroclor 1016 | 8.2E-01 | N | 9.4E-01 | Calculated | Yes |
| 11104-28-2 | ~Aroclor 1221 | 2.0E-01 | C | 5.9E-03 | Calculated | |
| 11141-16-5 | ~Aroclor 1232 | 1.7E-01 | C | 5.9E-03 | Calculated | |
| 53469-21-9 | ~Aroclor 1242 | 2.3E-01 | C | 5.5E-02 | Calculated | |
| 12672-29-6 | ~Aroclor 1248 | 2.3E-01 | C | 5.4E-02 | Calculated | |
| 11097-69-1 | ~Aroclor 1254 | 2.3E-01 | N | 9.1E-02 | Calculated | Yes |
| 11096-82-5 | ~Aroclor 1260 | 2.4E-01 | C | 2.4E-01 | Calculated | |
| 11126-42-4 | ~Aroclor 5460 | 7.0E+00 | N | 1.4E+01 | Calculated | Yes |
| 39635-31-9 | ~Heptachlorobiphenyl, 2,3,3',4,4',5,5'- (PCB 189) | 1.3E-01 | C | 1.3E-01 | Calculated | Yes |
| 52663-72-6 | ~Hexachlorobiphenyl, 2,3',4,4',5,5'- (PCB 167) | 1.2E-01 | C | 7.5E-02 | Calculated | Yes |
| 69782-90-7 | ~Hexachlorobiphenyl, 2,3,3',4,4',5'- (PCB 157) | 1.2E-01 | C | 7.7E-02 | Calculated | Yes |
| 38380-08-4 | ~Hexachlorobiphenyl, 2,3,3',4,4',5- (PCB 156) | 1.2E-01 | C | 7.7E-02 | Calculated | Yes |
| 32774-16-6 | ~Hexachlorobiphenyl, 3,3',4,4',5,5'- (PCB 169) | 1.2E-04 | C | 7.5E-05 | Calculated | Yes |
| 65510-44-3 | ~Pentachlorobiphenyl, 2',3,4,4',5- (PCB 123) | 1.2E-01 | C | 4.7E-02 | Calculated | Yes |
| 31508-00-6 | ~Pentachlorobiphenyl, 2,3',4,4',5- (PCB 118) | 1.2E-01 | C | 4.6E-02 | Calculated | Yes |

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|------------|---|---------------------------------------|-------|--|------------|---|
| 32598-14-4 | ~Pentachlorobiphenyl, 2,3,3',4,4'- (PCB 105) | 1.2E-01 | C | 4.7E-02 | Calculated | Yes |
| 74472-37-0 | ~Pentachlorobiphenyl, 2,3,4,4',5- (PCB 114) | 1.2E-01 | C | 4.7E-02 | Calculated | Yes |
| 57465-28-8 | ~Pentachlorobiphenyl, 3,3',4,4',5- (PCB 126) | 3.6E-05 | C | 1.4E-05 | Calculated | Yes |
| 1336-36-3 | ~Polychlorinated Biphenyls (high risk) | 2.3E-01 | C | 5.5E-02 | Calculated | |
| 1336-36-3 | ~Polychlorinated Biphenyls (low risk) | | | | | |
| 1336-36-3 | ~Polychlorinated Biphenyls (lowest risk) | | | | | |
| 32598-13-3 | ~Tetrachlorobiphenyl, 3,3',4,4'- (PCB 77) | 3.8E-02 | C | 8.4E-03 | Calculated | Yes |
| 70362-50-4 | ~Tetrachlorobiphenyl, 3,4,4',5- (PCB 81) | 1.2E-02 | C | 2.8E-03 | Calculated | Yes |
| 9016-87-9 | Polymeric Methylene Diphenyl Diisocyanate (PMDI) | 7.4E+06 | N | | | Yes |
| | Polynuclear Aromatic Hydrocarbons (PAHs) | | | | | |
| 83-32-9 | ~Acenaphthene | 7.2E+02 | N | 1.6E+01 | 2L or IMAC | Yes |
| 120-12-7 | ~Anthracene | 3.6E+03 | N | 1.3E+03 | 2L or IMAC | Yes |
| 56-55-3 | ~Benz[a]anthracene | 1.1E+00 | C | 3.5E-01 | 2L or IMAC | |
| 50-32-8 | ~Benzo[a]pyrene | 1.1E-01 | C | 1.2E-01 | 2L or IMAC | Yes |
| 205-99-2 | ~Benzo[b]fluoranthene | 1.1E+00 | C | 1.2E+00 | 2L or IMAC | |
| 192-97-2 | ~Benzo[e]pyrene | 1.1E+00 | N | 1.5E+01 | Calculated | Yes |
| 205-82-3 | ~Benzo[j]fluoranthene | 4.2E-01 | C | 7.0E-01 | Calculated | |
| 207-08-9 | ~Benzo[k]fluoranthene | 1.1E+01 | C | 1.2E+01 | 2L or IMAC | |
| 91-58-7 | ~Chloronaphthalene, Beta- | 9.6E+02 | N | 5.8E+01 | Calculated | Yes |
| 218-01-9 | ~Chrysene | 1.1E+02 | C | 3.6E+01 | 2L or IMAC | |
| 53-70-3 | ~Dibenz[a,h]anthracene | 1.1E-01 | C | 3.8E-01 | 2L or IMAC | |
| 192-65-4 | ~Dibenzo[a,e]pyrene | 4.2E-02 | C | 7.6E-01 | Calculated | |
| 57-97-6 | ~Dimethylbenz[a]anthracene, 7,12- | 4.6E-04 | C | 2.8E-03 | Calculated | |
| 206-44-0 | ~Fluoranthene | 4.8E+02 | N | 6.7E+02 | 2L or IMAC | Yes |
| 86-73-7 | ~Fluorene | 4.8E+02 | N | 1.1E+02 | 2L or IMAC | Yes |
| 193-39-5 | ~Indeno[1,2,3-cd]pyrene | 1.1E+00 | C | 3.9E+00 | 2L or IMAC | |
| 90-12-0 | ~Methylnaphthalene, 1- | 3.9E-02 | N | 1.1E-01 | 2L or IMAC | Yes |
| 91-57-6 | ~Methylnaphthalene, 2- | 4.8E+01 | N | 3.1E+00 | 2L or IMAC | Yes |
| 91-20-3 | ~Naphthalene | 2.1E+00 | C | 3.9E-01 | 2L or IMAC | Yes |
| 57835-92-4 | ~Nitropyrene, 4- | 4.2E-01 | C | 1.0E-01 | Calculated | |
| 198-55-0 | ~Perylene | 1.1E+00 | N | 1.5E+01 | Calculated | Yes |
| 129-00-0 | ~Pyrene | 3.6E+02 | N | 4.4E+02 | 2L or IMAC | Yes |
| 67747-09-5 | Prochloraz | 3.6E+00 | C | 2.4E-02 | Calculated | Yes |
| 26399-36-0 | Profluralin | 9.4E+01 | N | 5.1E+01 | Calculated | Yes |
| 1610-18-0 | Prometon | 1.9E+02 | N | 1.0E+00 | Calculated | Yes |
| 7287-19-6 | Prometryn | 5.1E+02 | N | 8.5E+00 | Calculated | Yes |

North Carolina Department of Environmental Quality

Soil Cleanup Goals (SCG)

October 2024

(based on May 2024 USEPA Regional Screening Tables)

This table must be used with the SCG Notes/ See Compound-Specific Notes for chemicals in *blue*.

| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|---|--|--------------|---|--------------|--|
| 23950-58-5 | Pronamide | 9.5E+02 | N | 1.1E+01 | Calculated | Yes |
| 1918-16-7 | Propachlor | 1.6E+02 | N | 1.1E+00 | Calculated | Yes |
| 709-98-8 | Propanil | 6.3E+01 | N | 3.9E-01 | Calculated | Yes |
| 2312-35-8 | Propargite | 2.8E+00 | C | 2.7E-01 | Calculated | Yes |
| 107-19-7 | Propargyl Alcohol | 3.1E+01 | N | 5.7E-02 | Calculated | Yes |
| 139-40-2 | Propazine | 2.5E+02 | N | 2.5E+00 | Calculated | Yes |
| 122-42-9 | Propham | 2.5E+02 | N | 1.8E+00 | Calculated | Yes |
| 60207-90-1 | Propiconazole | 1.3E+03 | N | 4.6E+01 | Calculated | Yes |
| 123-38-6 | Propionaldehyde | 1.6E+01 | N | | | Yes |
| 103-65-1 | Propyl benzene | 7.8E+02 | N | 2.6E+00 | 2L or IMAC | Yes |
| 115-07-1 | Propylene | 4.7E+02 | N | | | Yes |
| 57-55-6 | Propylene Glycol | 2.5E+05 | N | 4.0E+02 | 2L or IMAC | Yes |
| 6423-43-4 | Propylene Glycol Dinitrate | 3.4E+06 | N | | | Yes |
| 107-98-2 | Propylene Glycol Monomethyl Ether | 8.3E+03 | N | 2.0E+01 | Calculated | Yes |
| 75-56-9 | Propylene Oxide | 2.1E+00 | C | 6.1E-04 | Calculated | Yes |
| 110-86-1 | Pyridine | 1.6E+01 | N | 4.8E-02 | Calculated | Yes |
| 13593-03-8 | Quinalphos | 6.3E+00 | N | 6.0E-01 | Calculated | Yes |
| 91-22-5 | Quinoline | 1.8E-01 | C | 7.7E-04 | Calculated | |
| 76578-14-8 | Quizalofop-ethyl | 1.1E+02 | N | 2.0E+01 | Calculated | Yes |
| E715557 | Refractory Ceramic Fibers (units in fibers) | | | | | |
| 10453-86-8 | Resmethrin | 3.8E+02 | N | 2.6E+03 | Calculated | Yes |
| 299-84-3 | Ronnel | 7.8E+02 | N | 6.4E+01 | Calculated | Yes |
| 83-79-4 | Rotenone | 5.1E+01 | N | 2.9E+02 | Calculated | Yes |
| 94-59-7 | Safrole | 5.5E-01 | C | 2.0E-03 | Calculated | |
| 7783-00-8 | Selenious Acid | 7.8E+01 | N | | | Yes |
| 7782-49-2 | Selenium | 7.8E+01 | N | 2.1E+00 | 2L or IMAC | Yes |
| 7446-34-6 | Selenium Sulfide | 7.8E+01 | N | | | Yes |
| 74051-80-2 | Sethoxydim | 1.8E+03 | N | 1.8E+02 | Calculated | Yes |
| 7631-86-9 | Silica (crystalline, respirable) | 3.7E+07 | N | | | Yes |
| 7440-22-4 | Silver | 7.8E+01 | N | 3.4E+00 | 2L or IMAC | Yes |
| 122-34-9 | Simazine | 4.5E+00 | C | 3.9E-02 | 2L or IMAC | Yes |
| 62476-59-9 | Sodium Acifluorfen | 1.6E+02 | N | 1.4E+01 | Calculated | Yes |
| 26628-22-8 | Sodium Azide | 6.3E+01 | N | | | Yes |
| 148-18-5 | Sodium Diethyldithiocarbamate | 2.0E+00 | C | 1.6E-03 | Calculated | Yes |
| 7681-49-4 | Sodium Fluoride | 7.8E+02 | N | 5.3E+02 | Calculated | Yes |
| 62-74-8 | Sodium Fluoroacetate | 2.5E-01 | N | 5.7E-04 | Calculated | Yes |
| 13718-26-8 | Sodium Metavanadate | 1.6E+01 | N | | | Yes |
| 13472-45-2 | Sodium Tungstate | 1.3E+01 | N | | | Yes |
| 10213-10-2 | Sodium Tungstate Dihydrate | 1.3E+01 | N | | | Yes |
| 961-11-5 | Stirofos (Tetrachlorovinphos) | 2.3E+01 | C | 8.6E-02 | Calculated | Yes |
| 7440-24-6 | Strontium, Stable | 9.4E+03 | N | 1.4E+03 | 2L or IMAC | Yes |
| 57-24-9 | Strychnine | 3.8E+00 | N | 4.6E-01 | Calculated | Yes |
| 100-42-5 | Styrene | 1.2E+03 | N | 1.5E+00 | 2L or IMAC | Yes |

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|--|--|--------------|---|--------------|--|
| 57964-39-3 | Styrene-Acrylonitrile (SAN) Trimer (THNA isomer) | 3.8E+01 | N | | | Yes |
| 57964-40-6 | Styrene-Acrylonitrile (SAN) Trimer (THNP isomer) | 3.8E+01 | N | | | Yes |
| 126-33-0 | Sulfolane | 1.3E+01 | N | 3.1E-02 | Calculated | Yes |
| 80-07-9 | Sulfonylbis(4-chlorobenzene), 1,1'- | 1.0E+01 | N | 6.6E-01 | Calculated | Yes |
| 7446-11-9 | Sulfur Trioxide | 1.2E+07 | N | | | Yes |
| 7664-93-9 | Sulfuric Acid | 1.2E+07 | N | | | Yes |
| 140-57-8 | Sulfurous acid, 2-chloroethyl 2-[4-(1,1-dimethylethyl)phenoxy]-1-methylethyl ester | 2.2E+01 | C | 3.2E-01 | Calculated | Yes |
| 34014-18-1 | Tebuthiuron | 8.8E+02 | N | 2.8E+00 | Calculated | Yes |
| 3383-96-8 | Temephos | 2.5E+02 | N | 5.3E+02 | Calculated | Yes |
| 5902-51-2 | Terbacil | 1.6E+02 | N | 5.5E-01 | Calculated | Yes |
| 13071-79-9 | Terbufos | 3.9E-01 | N | 7.7E-03 | Calculated | Yes |
| 886-50-0 | Terbutryn | 1.3E+01 | N | 2.0E-01 | Calculated | Yes |
| 540-88-5 | Tert-Butyl Acetate | 8.6E+00 | C | 3.2E-02 | Calculated | |
| 5436-43-1 | Tetrabromodiphenyl ether, 2,2',4,4'- (BDE-47) | 1.3E+00 | N | 3.7E-01 | Calculated | Yes |
| 95-94-3 | Tetrachlorobenzene, 1,2,4,5- | 4.7E-01 | N | 1.9E-01 | 2L or IMAC | Yes |
| 630-20-6 | Tetrachloroethane, 1,1,1,2- | 2.1E+00 | C | 7.6E-03 | 2L or IMAC | Yes |
| 79-34-5 | Tetrachloroethane, 1,1,2,2- | 6.4E-01 | C | 1.6E-03 | 2L or IMAC | Yes |
| 127-18-4 | Tetrachloroethylene | 1.7E+01 | N | 6.3E-03 | 2L or IMAC | Yes |
| 58-90-2 | Tetrachlorophenol, 2,3,4,6- | 3.8E+02 | N | 3.0E+00 | 2L or IMAC | Yes |
| 5216-25-1 | Tetrachlorotoluene, p- alpha, alpha, alpha- | 4.3E-02 | C | 1.5E-04 | Calculated | Yes |
| 3689-24-5 | Tetraethyl Dithiopyrophosphate | 6.3E+00 | N | 5.1E-02 | Calculated | Yes |
| 811-97-2 | Tetrafluoroethane, 1,1,1,2- | 2.2E+04 | N | | | Yes |
| 16853-36-4 | Tetramethylphosphoramidate, -N,N,N',N" (TMPA) | 1.3E+00 | N | | | Yes |
| 479-45-8 | Tetryl (Trinitrophenylmethylnitramine) | 3.1E+01 | N | 2.6E+00 | Calculated | Yes |
| 1314-32-5 | Thallic Oxide | 3.1E-01 | N | | | Yes |
| 10102-45-1 | Thallium (I) Nitrate | 1.6E-01 | N | | | Yes |
| 7440-28-0 | Thallium (Soluble Salts) | 1.6E-01 | N | 2.8E+00 | 2L or IMAC | Yes |
| 563-68-8 | Thallium Acetate | 1.6E-01 | N | 2.8E-04 | Calculated | Yes |
| 6533-73-9 | Thallium Carbonate | 2.5E-01 | N | 2.9E-04 | Calculated | Yes |
| 7791-12-0 | Thallium Chloride | 1.6E-01 | N | | | Yes |
| 12039-52-0 | Thallium Selenite | 1.6E-01 | N | | | Yes |
| 7446-18-6 | Thallium Sulfate | 3.1E-01 | N | | | Yes |
| 79277-27-3 | Thifensulfuron-methyl | 5.4E+02 | N | 1.8E+00 | Calculated | Yes |
| 28249-77-6 | Thiobencarb | 1.3E+02 | N | 4.8E+00 | Calculated | Yes |

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|---|--|--------------|---|--------------|--|
| 21564-17-0 | Thiocyanic acid, (2-benzothiazolylthio)methyl ester (TCMTB) | 3.8E+02 | N | 2.9E+01 | Calculated | Yes |
| 111-48-8 | Thiodiglycol | 1.1E+03 | N | 2.0E+00 | Calculated | Yes |
| 39196-18-4 | Thiofanox | 3.8E+00 | N | 1.4E-02 | Calculated | Yes |
| 23564-05-8 | Thiophanate, Methyl | 4.7E+01 | C | 5.2E-02 | Calculated | Yes |
| 137-26-8 | Thiram | 1.9E+02 | N | 3.0E+00 | Calculated | Yes |
| 7440-31-5 | Tin | 9.4E+03 | N | 1.0E+04 | 2L or IMAC | Yes |
| 7550-45-0 | Titanium Tetrachloride | 1.2E+06 | N | | | Yes |
| 108-88-3 | Toluene | 9.9E+02 | N | 8.3E+00 | 2L or IMAC | Yes |
| 584-84-9 | Toluene-2,4-diisocyanate | 1.4E+00 | N | 2.7E-01 | Calculated | Yes |
| 91-08-7 | Toluene-2,6-diisocyanate | 1.1E+00 | N | 2.8E-01 | Calculated | Yes |
| 2687-25-4 | Toluenediamine, 2,3- | 1.3E+00 | N | 4.4E-03 | Calculated | Yes |
| 95-70-5 | Toluenediamine, 2,5- | 2.5E+00 | N | 1.2E-03 | Calculated | Yes |
| 496-72-0 | Toluenediamine, 3,4- | 1.3E+00 | N | 4.4E-03 | Calculated | Yes |
| 99-94-5 | Toluic Acid, p- | 6.3E+01 | N | 1.8E-01 | 2L or IMAC | Yes |
| 95-53-4 | Toluidine, o- (Methylaniline, 2-) | 3.4E+01 | C | 1.9E-02 | Calculated | |
| 106-49-0 | Toluidine, p- | 1.8E+01 | C | 9.9E-03 | Calculated | Yes |
| E1790670 | Total Petroleum Hydrocarbons (Aliphatic High) | 4.7E+04 | N | | | Yes |
| E1790666 | Total Petroleum Hydrocarbons (Aliphatic Low) | 5.1E+01 | N | 5.6E+00 | 2L or IMAC | Yes |
| E1790668 | Total Petroleum Hydrocarbons (Aliphatic Medium) | 2.0E+01 | N | 1.9E+02 | 2L or IMAC | Yes |
| E1790676 | Total Petroleum Hydrocarbons (Aromatic High) | 3.6E+00 | N | | | Yes |
| E1790674 | Total Petroleum Hydrocarbons (Aromatic Medium) | 6.2E+01 | N | | | Yes |
| 8001-35-2 | Toxaphene | 4.9E-01 | C | 9.3E-02 | 2L or IMAC | Yes |
| E1841606 | Toxaphene, Weathered | 3.8E-01 | N | 6.5E-01 | Calculated | Yes |
| 66841-25-6 | Tralomethrin | 9.5E+01 | N | 4.0E+02 | Calculated | Yes |
| 688-73-3 | Tri-n-butyltin | 4.7E+00 | N | 9.1E-01 | Calculated | Yes |
| 102-76-1 | Triacetin | 1.0E+06 | N | 3.2E+03 | Calculated | Yes |
| 43121-43-3 | Triadimefon | 4.3E+02 | N | 3.8E+00 | Calculated | Yes |
| 2303-17-5 | Triallate | 9.7E+00 | C | 2.2E-02 | Calculated | Yes |
| 82097-50-5 | Triasulfuron | 1.3E+02 | N | 1.5E+00 | Calculated | Yes |
| 101200-48-0 | Tribenuron-methyl | 1.0E+02 | N | 4.4E-01 | Calculated | Yes |
| 615-54-3 | Tribromobenzene, 1,2,4- | 7.8E+01 | N | 1.0E+00 | Calculated | Yes |
| 118-79-6 | Tribromophenol, 2,4,6- | 1.1E+02 | N | 2.3E+00 | Calculated | Yes |
| 78-48-8 | Tribufos | 2.5E+00 | N | 1.4E-01 | Calculated | Yes |
| 126-73-8 | Tributyl Phosphate | 6.0E+01 | C | 3.8E-01 | Calculated | Yes |
| E1790679 | Tributyltin Compounds | 3.8E+00 | N | | | Yes |
| 56-35-9 | Tributyltin Oxide | 3.8E+00 | N | 2.2E+03 | Calculated | Yes |

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|--------------|---|--|--------------|---|--------------|--|
| 10025-85-1 | Trichloramine | | | | | |
| 76-13-1 | Trichloro-1,2,2-trifluoroethane, 1,1,2- | 1.4E+03 | N | 9.8E+03 | 2L or IMAC | Yes |
| 76-03-9 | Trichloroacetic Acid | 7.8E+00 | C | 2.1E-03 | Calculated | Yes |
| 33663-50-2 | Trichloroaniline HCl, 2,4,6- | 1.9E+01 | C | 6.6E-02 | Calculated | |
| 634-93-5 | Trichloroaniline, 2,4,6- | 3.8E-01 | N | 3.8E-02 | Calculated | Yes |
| 87-61-6 | Trichlorobenzene, 1,2,3- | 1.3E+01 | N | 3.3E-01 | Calculated | Yes |
| 120-82-1 | Trichlorobenzene, 1,2,4- | 1.2E+01 | N | 4.1E+00 | 2L or IMAC | Yes |
| 71-55-6 | Trichloroethane, 1,1,1- | 1.7E+03 | N | 1.4E+00 | 2L or IMAC | Yes |
| 79-00-5 | Trichloroethane, 1,1,2- | 3.2E-01 | N | 3.9E-03 | 2L or IMAC | Yes |
| 79-01-6 | Trichloroethylene | 8.7E-01 | N | 2.1E-02 | 2L or IMAC | Yes |
| 75-69-4 | Trichlorofluoromethane | 4.7E+03 | N | 2.5E+01 | 2L or IMAC | Yes |
| 95-95-4 | Trichlorophenol, 2,4,5- | 1.3E+03 | N | 4.3E+00 | 2L or IMAC | Yes |
| 88-06-2 | Trichlorophenol, 2,4,6- | 1.3E+01 | N | 7.7E-02 | 2L or IMAC | Yes |
| 93-76-5 | Trichlorophenoxyacetic Acid, 2,4,5- | 1.3E+02 | N | 5.8E-01 | Calculated | Yes |
| 93-72-1 | Trichlorophenoxypropionic acid, -2,4,5 | 1.0E+02 | N | 5.5E-01 | 2L or IMAC | Yes |
| 598-77-6 | Trichloropropane, 1,1,2- | 7.8E+01 | N | 2.7E-01 | Calculated | Yes |
| 96-18-4 | Trichloropropane, 1,2,3- | 5.1E-03 | C | 4.3E-05 | 2L or IMAC | Yes |
| 96-19-5 | Trichloropropene, 1,2,3- | 1.6E-01 | N | 2.1E-01 | Calculated | Yes |
| 1330-78-5 | Tricresyl Phosphate (TCP) | 2.5E+02 | N | 2.6E+02 | Calculated | Yes |
| 58138-08-2 | Tridiphanne | 3.8E+01 | N | 3.0E+00 | Calculated | Yes |
| 121-44-8 | Triethylamine | 2.5E+01 | N | | | Yes |
| 112-27-6 | Triethylene Glycol | 2.5E+04 | N | 6.2E+01 | Calculated | Yes |
| 420-46-2 | Trifluoroethane, 1,1,1- | 3.1E+03 | N | | | Yes |
| 1582-09-8 | Trifluralin | 9.0E+01 | C | 3.0E+00 | Calculated | Yes |
| 512-56-1 | Trimethyl Phosphate | 2.7E+01 | C | 7.7E-03 | Calculated | Yes |
| 526-73-8 | Trimethylbenzene, 1,2,3- | 7.0E+01 | N | 2.1E+00 | Calculated | Yes |
| 95-63-6 | Trimethylbenzene, 1,2,4- | 6.3E+01 | N | 1.2E+01 | 2L or IMAC | Yes |
| 108-67-8 | Trimethylbenzene, 1,3,5- | 5.6E+01 | N | 1.1E+01 | 2L or IMAC | Yes |
| 25167-70-8 | Trimethylpentene, 2,4,4- | 1.6E+02 | N | 4.7E+00 | Calculated | Yes |
| 99-35-4 | Trinitrobenzene, 1,3,5- | 4.5E+02 | N | 1.5E+01 | Calculated | Yes |
| 118-96-7 | Trinitrotoluene, 2,4,6- | 7.3E+00 | N | 1.4E-01 | Calculated | Yes |
| 791-28-6 | Triphenylphosphine Oxide | 2.5E+02 | N | 1.2E+01 | Calculated | Yes |
| 13674-87-8 | Tris(1,3-Dichloro-2-propyl) Phosphate | 2.5E+02 | N | 6.3E+01 | Calculated | Yes |
| 13674-84-5 | Tris(1-chloro-2-propyl)phosphate | 1.3E+02 | N | 4.8E+00 | Calculated | Yes |
| 126-72-7 | Tris(2,3-dibromopropyl)phosphate | 2.8E-01 | C | 6.0E-03 | Calculated | |
| 115-96-8 | Tris(2-chloroethyl)phosphate | 2.7E+01 | C | 3.4E-02 | Calculated | Yes |
| 78-42-2 | Tris(2-ethylhexyl)phosphate | 1.7E+02 | C | 1.1E+03 | Calculated | Yes |
| 7440-33-7 | Tungsten | 1.3E+01 | N | 8.4E+00 | Calculated | Yes |
| 7440-61-1 | Uranium | 3.1E+00 | N | 6.3E+00 | Calculated | Yes |
| 51-79-6 | Urethane | 1.2E-01 | C | 1.6E-04 | Calculated | |

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| CAS # | Chemical Name (See companion notes for shaded chemicals) | Residential Health Based PSRG (mg/kg) | Basis | Protection of Groundwater PSRG (mg/kg) | Basis | Does Chemical Have a Non-Cancer Effect? |
|------------|---|---------------------------------------|-------|--|------------|---|
| 1314-62-1 | Vanadium Pentoxide | 1.4E+02 | N | | | Yes |
| 7440-62-2 | Vanadium and Compounds | 7.9E+01 | N | 1.4E+02 | 2L or IMAC | Yes |
| 1929-77-7 | Vernolate | 1.6E+01 | N | 1.1E-01 | Calculated | Yes |
| 50471-44-8 | Vinclozolin | 1.5E+01 | N | 1.3E-01 | Calculated | Yes |
| 108-05-4 | Vinyl Acetate | 1.9E+02 | N | 3.7E-01 | 2L or IMAC | Yes |
| 593-60-2 | Vinyl Bromide | 2.7E-01 | C | | | Yes |
| 75-01-4 | Vinyl Chloride | 6.1E-02 | C | 2.1E-04 | 2L or IMAC | Yes |
| 81-81-2 | Warfarin | 3.8E+00 | N | 4.4E-02 | Calculated | Yes |
| 108-38-3 | Xylene, m- | 1.2E+02 | N | 9.8E+00 | 2L or IMAC | Yes |
| 95-47-6 | Xylene, o- | 1.4E+02 | N | 9.8E+00 | 2L or IMAC | Yes |
| 106-42-3 | Xylene, p- | 1.2E+02 | N | 9.8E+00 | 2L or IMAC | Yes |
| 1330-20-7 | Xylenes | 1.2E+02 | N | 9.9E+00 | 2L or IMAC | Yes |
| 1314-84-7 | Zinc Phosphide | 4.7E+00 | N | | | Yes |
| 7440-66-6 | Zinc and Compounds | 4.7E+03 | N | 1.2E+03 | 2L or IMAC | Yes |
| 12122-67-7 | Zineb | 6.3E+02 | N | 2.0E+01 | Calculated | Yes |
| 7440-67-7 | Zirconium | 1.3E+00 | N | 1.7E+01 | Calculated | Yes |