

Department of Health

Environmental Health
108 Cherry Street – PO Box 70
Burlington, VT 05402-0070

HealthVermont.gov

Agency of Human Services

[phone] 802-863-7220
[fax] 802-863-7483
[toll free] 800-439-8550

Published April 2017; updated January 2023

The Vermont Department of Health (Health) develops and maintains two sets of chemical-specific soil guidance values, Residential Soil Values (RSVs) and Commercial Soil Values (CSVs), which may be used in the evaluation of potential exposure to chemicals in soil. RSVs are appropriate for consideration in the assessment of potential residential exposure to chemicals in soil. CSVs are appropriate for consideration in the assessment of potential exposure to chemicals in soil under a hypothetical nonresidential (e.g., commercial worker) exposure scenario.

The 2023 list of chemicals and associated soil values presented in Attachment 1 supplants any earlier guidance. Additional chemicals may be considered for evaluation and potential inclusion based on request by other State of Vermont offices, in response to public comments or as deemed appropriate by Health.

A summary of toxicity values and chemical-specific inputs is presented in Attachment 2, and exposure route and endpoint-specific risk-based concentrations are presented in Attachment 3a for RSVs and Attachment 3b for CSVs.

Each value is based upon the best available information at the time of derivation; therefore, is subject to change as updated information and risk assessment methodologies become available.

This memo provides an overview of the general approach used to develop the 2023 RSVs and CSVs and how these values are recommended to be applied.

RSV and CSV Derivation and Application

In general, RSVs are risk-based values generated by combining current toxicity values (e.g., oral reference doses, inhalation reference concentrations, oral cancer slope factors and inhalation unit risks) with a hypothetical residential exposure scenario using standard point estimate risk assessment procedures to derive an estimate of the concentration of each individual chemical (or in a few specific instances, a group of chemicals) in soil that corresponds to a fixed level of risk i.e., a target Hazard Quotient (THQ) of 1.0 for noncarcinogenic (systemic, threshold) effects or incremental lifetime cancer risk (ILCR) of one-in-one-million (1×10^{-6}). CSVs are similarly derived using a hypothetical nonresidential exposure scenario (i.e., commercial worker).

For each scenario, where a chemical is known to have both carcinogenic and noncarcinogenic effects and toxicity values are available, a risk-based value is derived based on each endpoint with the most conservative reported as the soil guidance value for that chemical under that specific scenario. Thus, the most conservative value derived for each chemical based on the residential scenario is reported as the RSV



for that chemical. Likewise, the most conservative value derived for each chemical based on the nonresidential scenario is reported as the CSV for that chemical.

Cumulative Risk Assessment

If more than one chemical is reported in excess of its analytical laboratory reporting limit, cumulative (total) cancer risk and noncancer hazard should be evaluated.

For carcinogens, it is recommended that a medium-specific cumulative (total) ILCR be estimated for each receptor, i.e., a Residential Soil or Nonresidential (Commercial) Soil cumulative ILCR, and compared to a target cumulative ILCR of 1×10^{-6} (one-in-one-million). This can be accomplished using the following approach where i represents the i^{th} such chemical reported in soil and, as described above, the target ICLR associated with each cancer risk-based value is 1×10^{-6} :

$$\text{Cumulative ILCR} = \sum_{i=1}^n (\text{SiteConcentration}_i)(1 \times 10^{-6}) / (\text{RSVi or CSV}_i)$$

A medium-specific cumulative ILCR in excess of 1×10^{-6} for a receptor suggests further consideration is warranted.

Similarly, for noncarcinogenic effects, it is recommended that a medium-specific total HQ, termed a Hazard Index (HI is the sum of two or more HQs), be estimated for each receptor (Residential and Nonresidential i.e., Commercial) and compared to a target Total HI=1.0. HQ are summed across all chemicals to develop the HI; chemicals are not segregated by critical effect. This can be accomplished using the following approach where i represents the i^{th} such chemical reported in soil and, as described above, the target HQ associated with each noncancer risk-based value is 1:

$$\text{Total Hazard Index} = \sum_{i=1}^n (\text{SiteConcentration}_i)(1) / (\text{RSVi or CSV}_i)$$

A medium-specific total HI in excess of 1 for a receptor suggests further consideration is warranted.

Soil Exposure Pathways

Direct exposure via incidental ingestion, dermal contact, inhalation of particles emitted from soil (i.e., fugitive dust) and, in the case of volatile compounds, inhalation due to volatilization from soil to ambient air were considered in the development of the 2023 values.

Exposure Assumptions

Several conservative assumptions are made in order to estimate the potential intake of a chemical in soil by a hypothetical receptor. In reality, the magnitude and frequency of exposure will vary depending on individual circumstances. The use of such health protective assumptions, which tend to represent reasonable upper bound estimates of exposure, adds additional conservatism to the risk-based soil values derived.

Overall, for the residential scenario a 70-year age-weighted approach (birth to age 70 years) is employed in the assessment of carcinogens while a hypothetical Young Child (birth to age 6 years) is the focus of assessment of noncarcinogenic (systemic, threshold) effects. Exposure is assumed to occur each day of the

year. The Commercial Worker scenario assumes a 30-year exposure duration as an adult. A duration weighted approach based on a lifetime of 70-years is employed in the assessment of carcinogens. The receptor is assumed to be on-site and exposed 250 days of the year for 10 hours per day. A summary of exposure assumptions and factors employed in the development of the 2023 RSVs and CSVs is presented in Attachments 4a and 4b.

Equations

Endpoint and exposure route specific equations used to develop residential and nonresidential risk-based values are presented in Attachments 5a and 5b. These equations combine chemical-specific toxicity information with scenario-specific exposure assumptions to generate a level in soil estimated to correspond to a fixed level of risk i.e., a THQ of 1.0 or ILCR of 1×10^{-6} for the receptor considered.

Chemical Specific Parameters

Volatility

For purpose of this effort, chemicals with a Henry's Law constant greater than 1×10^{-5} atmosphere-cubic meter/mole ($\text{atm}\cdot\text{m}^3/\text{mol}$) or a vapor pressure greater than 1 millimeter of mercury (mm Hg) at 25°C were deemed sufficiently volatile to warrant consideration of this route of exposure. These chemicals are flagged "v" in Attachments 2a and 2b. As recommended by the Vermont Department of Environmental Conservation, for this subset, chemical-specific volatilization factors based on a groundwater temperature of 15°C were developed using the Regional Screening Levels (RSL) on-line calculator. The RSL calculator was also used to generate chemical-specific soil saturation concentrations (C_{sat}) for members of this group that are not solid at ambient temperature. If the derived cancer or noncancer risk-based value for soil inhalation was greater than the corresponding C_{sat}, the C_{sat} was used in its place.

In addition, for "v" chemicals only, at the advice of the Vermont IRULE workgroup, a snow cover modification factor (SCMF) less than 1 was employed in the development of residential soil inhalation risk-based values (cancer and noncancer). An SCMF of 0.7342 was applied to generate a residential soil inhalation exposure frequency of 268 days. This reflects the estimated average number of days per year in Vermont with less than one inch of snow cover over the last decade (a ten-year average of approximately 97 days with at least one inch of snow cover was provided for consideration in this effort). At present, risk-based values for the soil inhalation pathway are developed by considering exposure via inhalation as volatile and as fugitive dust in a single equation. At the advice of the IRULE workgroup, the SCMF was applied to the entire equation, not just the volatile component, given the level of uncertainty regarding dissipation over time of "v" chemicals from dust. Such application implies that exposure to any residual particulate emissions indoors does not contribute significantly to the inhalation route of exposure. An SCMF of 1 was used for other residential soil routes of exposure and for all routes of exposure for chemicals not meeting the "v" criteria.

Physical Chemical Constants

Estimates of chemical-specific physical properties, such as Henry's Law constant, were primarily obtained from the Estimation Programs Interface (EPI) Suite™ which is a screening level tool developed by the U.S. Environmental Protection Agency (U.S. EPA) Office of Pollution Prevention and Toxics and Syracuse Research Corporation. In accordance with EPI Suite guidance, experimental (measured) values are used when both experimental and estimated values are available.

Absorption Factors

An oral relative bioavailability factor of 1 is employed in the development of soil guidance values (i.e., bioavailability is assumed to be similar between the medium of environmental exposure and that associated with the toxicity value).

Consistent with current guidance (EPA 2004), dermal exposure is evaluated using oral toxicity values adjusted for gastrointestinal absorption efficiency, where warranted, and incorporation of a dermal absorption fraction from soil. For purposes of this effort, a default gastrointestinal absorption efficiency of 1 is used in the evaluation of chemicals not listed in Exhibit 4-1 of said guidance. Chemical-specific dermal absorption fractions provided in Exhibit 3-4 are employed including the recommended default value of 0.1 for those semi-volatile organic compounds not called out by name in the list. In addition, Health employs a dermal absorption fraction of 0.01 for inorganics that are not listed in Exhibit 3-4 and a value of 0.03 in the evaluation of volatiles.

These evaluations do not account for existing background concentrations of naturally occurring inorganics. In some cases, it is possible that the theoretical derived value may be below naturally occurring levels. In such instances, the theoretical derived value may not be applicable, as clean-up below background is usually not advisable.

Toxicity Information Sources

Toxicity information and oral and inhalation toxicity values are obtained and reviewed from relevant and appropriate sources including:

- U.S. EPA Integrated Risk Information System
- U.S. EPA Office of Pesticide Programs
- U.S. EPA Office of Research and Development/National Center for Environmental Assessment/Superfund Health Risk Technical Support Center (STSC) Provisional Peer Reviewed Toxicity Values
- International Agency for Research on Cancer
- National Toxicology Program
- California EPA Office of Environmental Health Hazard Assessment
- California Department of Pesticide Regulation
- Agency for Toxic Substances and Disease Registry

Mutagenic Mode of Action

Consistent with U.S. EPA guidance (EPA 2005), multipliers termed Age Dependent Adjustment Factors (ADAFs) are used in the evaluation of carcinogens identified by U.S. EPA to operate via a mutagenic mode of action. Such chemicals are noted in the detailed Summary Table (Attachment 2a). Per the guidance, ADAFs "...reflect the potential for early-life exposure to make a greater contribution to the cancers appearing later in life."

Chemical-specific ADAFs are used if available.

Otherwise, the following non-chemical specific, default adjustments provided by U.S. EPA are used:

- A 10-fold increase for exposures between the day of birth up until the second birthday.
- A 3-fold increase for exposures between the second birthday up until the sixteenth birthday.
- No adjustment is made for exposures occurring after turning 16 years of age.

Toxicity Equivalence Factors and Relative Potency Factors

Some chemicals are members of the same family or group and have been shown to exhibit similar toxicological properties; however, each chemical may differ in the degree of toxicity (EPA 2019). In some such instances, a toxicity equivalency factor (TEF) or relative potency factor (RPF) must be applied to convert the reported concentration of each member of the group to a toxicity equivalent concentration (TEQ) relative to the toxicity of the index chemical for the group. The index chemical is assigned a TEF of 1. Total TEQ for a sample can then be compared to the value for the index chemical.

Dioxins, Furans and dioxin-like Polychlorinated Biphenyls (PCBs)

The index chemical for this group is 2,3,7,8-Tetrachlorodibenzo-*p*-dioxin (2,3,7,8-TCDD). As of this writing, Health recommends that the 2005 World Health Organization TEFs (Van den Berg et al., 2006) be employed in the evaluation of dioxins, furans and dioxin-like PCBs. These values are also presented in the May 2013 U.S. EPA fact sheet, "Use of Dioxin TEFs in Calculating Dioxin TEQs at CERCLA and RCRA Sites" which references the 2010 U.S. EPA report, "Recommended Toxicity Equivalency Factors (TEFs) for Human Health Risk Assessments of 2,3,7,8-Tetrachlorodibenzo-*p*-dioxin and Dioxin-Like Compounds" (EPA 2019). TEFs for Di-ortho PCBs may be obtained from Ahlborg, U.G. et al., 1994 (EPA 2019). TEFs may be applied to the ingestion, dermal (see EPA 2004) or inhalation routes of exposure and adjusted values may be used in the assessment of both cancer and noncarcinogenic effects (EPA 2013). The sum of adjusted concentrations is often referred to as 2,3,7,8-TCDD TEQ.

Dioxin Toxicity Equivalence Factors (EPA 2022)

CAS Registry Number	Compound	2,3,7,8-TCDD Toxicity Equivalence Factor
Chlorinated dibenzo-<i>p</i>-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
72918-21-9	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
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

CAS Registry Number	Compound	2,3,7,8-TCDD Toxicity Equivalence Factor
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

Polychlorinated biphenyls (PCBs)			
	IUPAC No.	Structure	2,3,7,8-TCDD Toxicity Equivalence Factor
<i>Non-ortho</i>			
32598-13-3	77	3,3',4,4'-TetraCB	0.0001
70362-50-4	81	3,4,4',5-TetraCB	0.0003
57465-28-8	126	3,3',4,4',5-PeCB	0.1
32774-16-6	169	3,3',4,4',5,5'-HxCB	0.03
<i>Mono-ortho</i>			
32598-14-4	105	2,3,3',4,4'-PeCB	0.00003
74472-37-0	114	2,3,4,4',5-PeCB	0.00003
31508-00-6	118	2,3',4,4',5-PeCB	0.00003
65510-44-3	123	2',3,4,4',5-PeCB	0.00003
38380-08-4	156	2,3,3',4,4',5-HxCB	0.00003
69782-90-7	157	2,3,3',4,4',5'-HxCB	0.00003
52663-72-6	167	2,3',4,4',5,5'-HxCB	0.00003
39635-31-9	189	2,3,3',4,4',5,5'-HpCB	0.00003
<i>Di-ortho*</i>			
35065-30-6	170	2,2',3,3',4,4',5-HpCB	0.0001
35065-29-3	180	2,2',3,4,4',5,5'-HpCB	0.00001

*Di-ortho values come from Ahlborg, U.g., et al (1994), which are the WHO 1994 values from Toxic equivalency factors for dioxin-like PCBs: Report on WHO-ECEH and IPCS consultation. December 1993. Chemosphere Volume 28, Issue 6. March 1994. Pages 1049-1067.

Carcinogenic Polycyclic Aromatic Hydrocarbons (cPAH)

Benzo(a)pyrene is the index chemical for this group of compounds. As of this writing, Health recommends that the following RPFs (EPA 1993) be employed in the evaluation of cPAH. Adjusted values are summed and expressed as Benz(a)pyrene toxic equivalents i.e., B(a)P-TE. B(a)P-TE is assessed for cancer risk only. Only B(a)P itself is assessed for potential noncarcinogenic effects.

Relative Potency Factors for Carcinogenic Polycyclic Aromatic Hydrocarbons

Compound	Benzo(a)pyrene Relative Potency Factor
Benzo(a)pyrene	1
Benzo(a)anthracene	0.1
Benzo(b)fluoranthene	0.1
Benzo(k)fluoranthene	0.01
Chrysene	0.001
Dibenzo(a,h)anthracene	1
Indeno(1,2,3cd)pyrene	0.1

Food Quality Protection Act Safety Factor for Pesticides

The Food Quality Protection Act (FQPA) of 1996 amended the Federal Insecticide, Fungicide and Rodenticide Act and the Federal Food, Drugs and Cosmetic Act and significantly revised the way in which pesticides are evaluated by the U.S. EPA.

The FQPA mandates that “in the case of threshold effects, an additional tenfold margin of safety for the pesticide chemical residue and other sources of exposure shall be applied for infants and children to take into account potential pre- and post-natal toxicity and completeness of data with respect to exposure and toxicity to infants and children. Notwithstanding such requirement for an additional margin of safety, the Administrator [of the U.S. EPA] may use a different margin of safety for the pesticide chemical residue only if, on the basis of reliable data, such margin will be safe for infants and children” (FQPA 1996).

Risk-based values derived for pesticides based on threshold type effects may reflect incorporation of a U.S. EPA derived FQPA Safety Factor (SF). Use of any FQPA SF greater than 1 is noted in the detailed Summary Table (Attachment 2).

References

EPA 1993. Provisional Guidance for Quantitative Risk Assessment of Polycyclic Aromatic Hydrocarbons. U.S. Environmental Protection Agency. Research Triangle Park, N.C. EPA/600/R-93/089, July 1993.

EPA 2004. Risk Assessment Guidance for Superfund Volume I: Human Health Evaluation Manual (Part E, Supplemental Guidance for Dermal Risk Assessment) Final. U.S. Environmental Protection Agency. Washington, D.C. OSWER 9285.7-02EP. July 2004.

EPA 2005. Supplemental Guidance for Assessing Susceptibility from Early-Life Exposure to Carcinogens. U.S. Environmental Protection Agency. Washington, D.C. EPA/630/R-03/003F. March 2005.

EPA 2008. Child-Specific Exposure Factor Handbook. U.S. Environmental Protection Agency. Office of Research and Development, National Center for Environmental Assessment. Washington, D.C. EPA/600/R-06/096F. September 2008.

EPA 2022 United States Environmental Protection Agency. Regional Screening Levels for Chemical Contaminants at Superfund Sites. User's Guide. November 2022 edition (accessed January, 2023).

FQPA 1996. Food Quality Protection Act of August 3, 1996 as amended. United States Public Law 104-170.

Van den Berg et al., 2006. The 2005 World Health Organization re-evaluation of human and mammalian toxic equivalency factors for dioxins and dioxin-like compounds. *Toxicol Sci* 93(2):223-241.

ATTACHMENT 1
VERMONT DEPARTMENT OF HEALTH
2023 RESIDENTIAL & COMMERCIAL SOIL VALUES (mg/kg)
INCIDENTAL INGESTION, DERMAL CONTACT AND INHALATION

Chemical Name	CAS No.	Residential Soil Value (RSV)		Commercial Soil Value (CSV)	
		RSV (mg/kg)	Endpoint	CSV (mg/kg)	Endpoint
Acetochlor	34256-82-1	1216	nc	14362	nc
Acetone	67-64-1	63079	nc	816169	nc
Alachlor	15972-60-8	5.0	nc	30	c
Aldrin	309-00-2	0.020	c	0.10	c
Aluminum	7429-90-5	72507	nc	941748	nc
Antimony	7440-36-0	29	nc	392	nc
Arsenic, Inorganic	7440-38-2	0.23	c	1.41	c
Barium	7440-39-3	11247	nc	127382	nc
Benomyl	17804-35-2	116	c	701	c
Benzene	71-43-2	0.70	c	4.19	c
Benzo(a)pyrene	50-32-8	0.073	c-mmoa(a)	1.54	c
Beryllium	7440-41-7	35	nc	289	nc
Bis(2-chloro-1-methyl ethyl)ether	108-60-1	2804	nc	36274	nc
Boron	7440-42-8	14650	nc	195698	nc
Bromate	15541-45-4	0.54	c	3.27	c
Bromochloromethane	74-97-5	193	nc	597	nc
Bromoxynil	1689-84-5	2.69	c	16.27	c
Butylbenzene, n-	104-51-8	3504	nc	45343	nc
Butylbenzene, sec-	135-98-8	7009	nc	90685	nc
Butylbenzene, tert-	98-06-6	7009	nc	90685	nc
Cadmium (food)	7440-43-9	6.86	nc	87.24	nc
Carbaryl	63-25-2	317	c	1915	c
Carbon Disulfide	75-15-0	608	nc	661	nc
Carbon tetrachloride	56-23-5	0.37	c	2.23	c
Chlorobenzene	108-90-7	267	nc	676	nc
Chromium (III) (insoluble salts)	16065-83-1	40223	nc	360223	nc
Chromium (VI)	18540-29-9	0.091	c-mmoa	1.75	c
Cobalt	7440-48-4	22	nc	291	nc
Copper	7440-50-8	10407	nc	139231	nc
Di (2-ethylhexyl) phthalate	117-81-7	20	c	120	c
Dibromochloropropane	96-12-8	0.0060	c-mmoa	0.061	c
Dibromoethane, 1,2-	106-93-4	0.023	c	0.14	c
Dichloroethane, 1,1-	75-34-3	2.10	c	12.59	c
Dichloroethane, 1,2-	107-06-2	0.29	c	1.71	c
Dichloroethylene, cis 1,2-	156-59-2	77	nc	352	nc
Dichloroethylene, trans 1,2-	156-60-5	108	nc	296	nc
Dichloropropane, 1,2-	78-87-5	1.51	c	9.06	c
Dioxane, 1,4-	123-91-1	2.78	c	16.89	c
Ethylbenzene	100-41-4	3.68	c	22.11	c
Fluoranthene	206-44-0	2301	nc	26371	nc
Fluorene	86-73-7	2301	nc	26371	nc
Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6	0.23	nc(c)	3	nc(c)
Hexachlorobenzene	118-74-1	0.13	c	0.69	c
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	4.60	c	28.03	c
Iron	7439-89-6	51302	nc	686351	nc
Isopropylbenzene (cumene)	98-82-8	256	nc	264	nc
Lead		41	nc(b)	41	nc(b)
Manganese (non-diet)	7439-96-5	1118	nc	11350	nc
Mercury (elemental)	7439-97-6	3.13	nc	3.13	nc
Methyl ethyl ketone	78-93-3	16952	nc	26991	nc
Methyl tert-butyl ether (MTBE)	1634-04-4	27	nc	165	c
Molybdenum	7439-98-7	366	nc	4900	nc
Naphthalene	91-20-3	1.2	c	7.19	c
Nickel	7440-02-0	940	nc	9707	nc
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0	3698	nc	49834	nc
Pentachlorophenol	87-86-5	0.48	c	2.90	c
Pentaerythritol tetranitrate (PETN)	78-11-5	65	nc	390	c
Perchlorate	14797-73-0	51	nc	686	nc
Perfluorobutane sulfonic acid (PFBS)	375-73-5	18	nc(c)	215	nc(c)
Perfluorohexane sulfonic acid (PFHxS)	355-46-4	1.2	nc(c)	14	nc(c)
Perfluorononanoic acid (PFNA)	375-95-1	0.18	nc(c)	2.2	nc(c)
Perfluorooctane sulfonic acid (PFOS)	1763-23-1	0.12	nc(c)	1.4	nc(c)

Chemical Name	CAS No.	Residential Soil Value (RSV)		Commercial Soil Value (CSV)	
		RSV (mg/kg)	Endpoint	CSV (mg/kg)	Endpoint
Perfluorooctanoic acid (PFOA)	335-67-1	0.18	nc(c)	2.2	nc(c)
Polychlorinated biphenyls (PCBs)	1336-36-3	0.11	c(d)	0.68	c(d)
Propoxur (Baygon)	114-26-1	79	c	476	c
Propyl benzene, n-	103-65-1	253	nc	261	nc
Selenium	7782-49-2	366	nc	4900	nc
Silver	7440-22-4	237	nc	2483	nc
Tetrachlorodibenzo-p-dioxin, 2,3,7,8- (TCDD)	1746-01-6	2.58E-06	c	1.57E-05	c
Tetrachloroethane, 1,1,1,2-	630-20-6	1.32	c	7.91	c
Tetrachloroethylene	127-18-4	2.38	c	14.29	c
Thallium (soluble Thallium)	7440-28-0	0.73	nc	9.8	nc
Toluene	108-88-3	705	nc	798	nc
Trichloroethylene (non-moa)	79-01-6	1.9	c/c--m-moa	6.47	c
Trichloropropane, 1,2,3-	96-18-4	0.0031	c-m-moa	0.071	c
Trimethylbenzene, 1,2,3-	526-73-8	206	nc	282	nc
Trimethylbenzene, 1,2,4-	95-63-6	166	nc	212	nc
Trimethylbenzene, 1,3,5-	108-67-8	144	nc(e)	177	nc(e)
Trinitrotoluene, 2,4,6- (TNT)	118-96-7	12	c	70	c
Uranium (soluble salts)	7440-61-1	44	nc	588	nc
Vanadium	7440-62-2	2.8	nc	27.22	nc
Vinyl chloride	75-01-4	0.10	c-m-moa	0.59	c
Xylenes	1330-20-7	252	nc	257	nc
Zinc	7440-66-6	21986	nc	294150	nc

Notes:

mg/kg - milligram per kilogram

c - RSV corresponds to a one-in-one-million incremental lifetime cancer risk

c-m-moa - carcinogen identified by U.S. EPA to operate via mutagenic mode of action

nc - RSV corresponds to a noncarcinogenic Hazard Quotient as indicated (HQ=1.0) based on Hypothetical Young Child Resident scenario

* - CAS Number is for Metallic Thallium

(a) Benzo(a)pyrene cancer-based value applicable to benzo(a)pyrene itself and to total benzo(a)pyrene toxic equivalents [B(a)P-TE]. Benzo(a)pyrene noncancer-based value applicable only to benzo(a)pyrene itself.

(b) Health used EPA's IEUBK software and Adult Lead Model to find the soil concentration that would result in 95% probability that no exposed child (or fetus) would exceed the BLL of < 3 ug/dL. Vermont DEC conducted a soil background study and determined the rural background level of lead is 41 mg/kg. Because the soil concentration that would result in a 95% probability that no exposed child (or fetus) would exceed the BLL < 3 ug/dL is less than the Vermont background level, the background level is recommended

(c) PFAS - RSV and CSV are based on direct contact with soil and do not represent concentrations protective for other potential exposure pathways such as leaching from soil to groundwater.

(d) PCBs- sum of all PCBs not to exceed 1.14E-01 mg/kg (IRIS high risk and persistence cancer toxicity values used in cancer assessment; oral reference dose for Aroclor 1254 used in noncancer assessment).

(e) Trimethyl benzenes -Sum of the three isomers not to exceed applicable resident or non-resident values, based on the most conservative value derived for an individual isomer.

ATTACHMENT 2
VERMONT DEPARTMENT OF HEALTH
TOXICITY & CHEMICAL SPECIFIC VALUES

Chemical Name	CAS No.	Toxicity and Chemical Specific Information											
		CSFo (mg/kg-d) ¹	Source	IUR (µg/m ³) ¹	Source	RfDo mg/kg-d	Source	RfC mg/m ³	Source	Volatile	Mutagenic	ABSd	GIABS
Acetochlor	34256-82-1	NA	O	NA	O	2.00E-02	LO	NA	--	-	-	0.10	1
Acetone	67-64-1	NA	I	NA	I	9.00E-01	I	NA	A	V	-	0.03	1
Alachlor	15972-60-8	5.60E-02	CE	NA	O	1.00E-02	O	NA	--	-	-	0.10	1
Aldrin	309-00-2	1.70E+01	I	4.90E-03	I	4.00E-05	A	NA	--	V	-	0.03	1
Aluminum	7429-90-5	NA	P	NA	P	1.00E+00	P	5.00E-03	P	-	-	0.01	1
Antimony	7440-36-0	NA	I	NA	I	4.00E-04	I	3.00E-04	A	-	-	0.01	1
Arsenic, Inorganic	7440-38-2	1.50E+00	I	4.30E-03	I	3.00E-04	I	1.50E-05	CE	-	-	0.03	0.60
Barium	7440-39-3	NA	I	NA	I	2.00E-01	I	NA	-	-	-	0.01	0.07
Benenyl	17804-35-2	2.39E-03	O	NA	O	1.30E-02	O ^(b)	NA	--	-	-	0.10	1
Benzene	71-43-2	5.50E-02	(c)	7.80E-06	(d)	4.00E-03	I	3.00E-02	I	V	-	0.03	1
Benzo(a)pyrene	50-32-8	1.00E+00	I	6.00E-04	I	3.00E-04	I	2.00E-06	I	-	yes	0.13	1
Beryllium	7440-41-7	NA	I	2.40E-03	I	2.00E-03	I	2.00E-05	I	-	-	0.01	0.01
Bis(2-chloro-1-methyl ethyl)ether	108-60-1	NA	P	NA	P	4.00E-02	I	NA	I	V	-	0.03	1
Boron	7440-42-8	NA	I	NA	I	2.00E-01	I	2.00E-02	A	-	-	0.01	1
Bromate	15541-45-4	7.00E-01	I	NA	I	4.00E-03	I	NA	I	-	-	0.01	1
Bromochloromethane	74-97-5	NA	I	NA	I	1.30E-02	E	4.00E-02	P	V	-	0.03	1
Bromoxynil	1689-84-5	1.03E-01	O	NA	O	1.50E-02	O	NA	--	-	-	0.10	1
Butylbenzene, n-	104-51-8	NA	P	NA	P	5.00E-02	P	NA	--	V	-	0.03	1
Butylbenzene, sec-	135-98-8	NA	P	NA	P	1.00E-01	PA	NA	--	V	-	0.03	1
Butylbenzene, tert-	98-06-6	NA	P	NA	P	1.00E-01	PA	NA	--	V	-	0.03	1
Cadmium (food)	7440-43-9	NA	I	1.80E-03	I	1.00E-04	(e)	1.00E-05	A	-	-	0.001	1
Carbaryl	63-25-2	8.75E-04	O	NA	O	1.00E-01	I ^(f)	NA	--	-	-	0.10	1
Carbon Disulfide	75-15-0	NA	--	NA	--	1.00E-01	I	7.00E-01	I	V	-	0.03	1
Carbon tetrachloride	56-23-5	7.00E-02	I	6.00E-06	I	4.00E-03	I	1.00E-01	I	V	-	0.03	1
Chlorobenzene	108-90-7	NA	I	NA	I	7.00E-03	A	5.00E-02	P	V	-	0.03	1
Chromium (III) (insoluble salts)	16065-83-1	NA	I	NA	I	1.50E+00	I	NA	I	-	-	0.01	0.01
Chromium (VI)	18540-29-9	5.00E-01	(g)	8.40E-02	(h)	3.00E-03	I	1.00E-04	I	-	yes	0.01	0.03
Cobalt	7440-48-4	NA	P	9.00E-03	P	3.00E-04	P	6.00E-06	P	-	-	0.01	1
Copper	7440-50-8	NA	I	NA	I	1.42E-01	CE	NA	--	-	-	0.01	1
Di (2-ethylhexyl) phthalate	117-81-7	1.40E-02	I	2.40E-06	C	2.00E-02	I	NA	--	-	-	0.10	1
Dibromochloropropane	96-12-8	8.00E-01	P	6.00E-03	P	2.00E-04	P	2.00E-04	I	V	yes	0.03	1
Dibromomethane, 1,2-	106-93-4	2.00E+00	I	6.00E-04	I	9.00E-03	I	9.00E-03	I	V	-	0.03	1
Dichloroethane, 1,1-	75-34-3	5.70E-03	CE	1.60E-06	CE	2.00E-01	P	NA	--	V	-	0.03	1
Dichloroethane, 1,2-	107-06-2	9.10E-02	I	2.60E-05	I	2.00E-02	A	7.00E-03	P	V	-	0.03	1
Dichloroethylene, cis 1,2-	156-59-2	NA	I	NA	I	2.00E-03	I	4.00E-02	PA	V	-	0.03	1
Dichloroethylene, trans 1,2-	156-60-5	NA	I	NA	I	2.00E-02	I	4.00E-02	PA	V	-	0.03	1
Dichloropropane, 1,2-	78-87-5	3.70E-02	P	3.70E-06	P	7.00E-02	A	4.00E-03	I	V	-	0.03	1
Dioxane, 1,4-	123-91-1	1.00E-01	I	5.00E-06	I	3.00E-02	I	3.00E-02	I	V	-	0.03	1
Ethylbenzene	100-41-4	1.10E-02	CE	2.50E-06	CE	5.00E-02	U	2.60E-01	A	V	-	0.03	1
Fluoranthene	206-44-0	NA	I	NA	I	4.00E-02	I	NA	I	-	-	0.13	1
Fluorene	86-73-7	NA	I	NA	I	4.00E-02	I	NA	I	V	-	0.13	1
Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6	NA	NA	NA	NA	3.00E-06	E	NA	-	-	-	-	-
Hexachlorobenzene	118-74-1	1.60E+00	I	4.60E-04	I	1.00E-05	P	NA	I	V	-	0.03	1
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	8.00E-02	I	NA	I	4.00E-03	I	NA	I	-	-	0.02	1
Iron	7439-89-6	NA	P	NA	P	7.00E-01	P	NA	P	-	-	0.01	1
Isopropylbenzene (cumene)	98-82-8	NA	I	NA	I	1.00E-01	I	4.00E-01	I	V	-	-	1
Lead ^(b)	7439-92-1	NA	--	NA	--	NA	--	NA	--	-	-	-	-
Manganese (non-diet)	7439-96-5	NA	I	NA	I	2.40E-02	I	5.00E-05	I	-	-	0.01	0.04
Mercury (elemental)	7439-97-6	NA	I	NA	I	NA	I	3.00E-04	I	V	-	0.01	1
Methyl ethyl ketone	78-93-3	NA	I	NA	I	6.00E-01	I	5.00E+00	I	V	-	0.03	1
Methyl tert-butyl ether (MTBE)	1634-04-4	1.80E-03	CE	2.60E-07	CE	1.00E-02	VH	3.00E+00	I	V	-	0.03	1
Molybdenum	7439-98-7	NA	I	NA	I	5.00E-03	I	2.00E-03	A	-	-	0.01	1
Naphthalene	91-20-3	1.20E-01	CE	3.40E-05	CE	2.00E-02	I	3.00E-03	I	V	-	0.13	1
Nickel	7440-02-0	NA	I	2.60E-04	CE	2.00E-02	I ^(f)	9.00E-05	A	-	-	0.01	0.04
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0	NA	I	NA	I	5.00E-02	I	NA	--	-	-	0.01	1
Pentachlorophenol	87-86-5	4.00E-01	I	5.10E-06	C	5.00E-03	I	NA	I	-	-	0.25	1
Pentaerythritol tetranitrate (PETN)	78-11-5	4.30E-03	PA	NA	--	9.00E-03	P	NA	--	-	-	0.10	1
Perchlorate	14797-73-0	NA	I	NA	I	7.00E-04	VH	NA	I	-	-	0.01	1
Perfluorobutane sulfonic acid (PFBS)	375-73-5	NA	--	NA	--	2.00E-05	P	NA	--	-	-	0.10	1
Perfluorohexane sulfonic acid (PFHxS)	355-46-4	NA	--	NA	--	3.00E-06	A	NA	--	-	-	0.10	1
Perfluorononanoic acid (PFNA)	375-95-1	NA	--	NA	--	3.00E-06	A	NA	--	-	-	0.10	1
Perfluorooctane sulfonic acid (PFOS)	1763-23-1	NA	E	NA	E	2.00E-06	A	NA	E	-	-	0.10	1
Perfluorooctanoic acid (PFOA)	335-67-1	7.00E-02	E	NA	E	3.00E-06	A	NA	E	-	-	0.10	1
Polychlorinated biphenyls (PCBs)	1336-36-3	2.00E+00	I ^(k)	5.70E-04	I ^(k)	2.50E-05	I ^(b)	NA	I	V	-	0.14	1
Propoxur (Baygon)	114-26-1	3.52E-03	O	NA	O	4.00E-03	I ^(f)	NA	--	-	-	0.10	1
Propyl benzene, n-	103-65-1	NA	--	NA	--	1.00E-01	(l)	2.60E-01	(m)	V	-	-	1
Selenium	7782-49-2	NA	I	NA	I	5.00E-03	I	1.80E-03	VA ⁽ⁿ⁾	-	-	0.01	1
Silver	7440-22-4	NA	I	NA	I	5.00E-03	I	NA	--	-	-	0.01	0.04
Tetrachlorodibenzo-p-dioxin, 2,3,7,8- (TCDD)	1746-01-6	1.30E+05	CE	3.80E+01	CE	7.00E-10	I	4.00E-08	CE	V	-	0.03	1
Tetrachloroethane, 1,1,1,2-	630-20-6	2.60E-02	I	7.40E-06	I	3.00E-02	I	NA	--	V	-	0.03	1
Tetrachloroethylene	127-18-4	1.10E-02	(o)	1.60E-06	(p)	6.00E-03	I	4.00E-02	I	V	-	0.03	1
Thallium (soluble Thallium)	7440-28-0 (a)	NA	I	NA	I	1.00E-05	P	NA	I	-	-	0.01	1
Toluene	108-88-3	NA	I	NA	I	8.00E-02	I	5.00E+00	I	V	-	0.03	1
Trichloroethylene (non-mmoa / mmoa)	79-01-6	3.72E-02/9.3E-03	I	3.1E-06/1.0E-06	I	5.00E-04	I	2.00E-03	I	V	yes	0.03	1
Trichloropropane, 1,2,3-	96-18-4	3.00E+01	I	NA	I	1.00E-02	A	3.00E-04	I	V	yes	0.03	1
Trimethylbenzene, 1,2,3-	526-73-8	NA	I	NA	I	1.00E-02	I	6.00E-02	I	V	-	0.03	1
Trimethylbenzene, 1,2,4-	95-63-6	NA	I	NA	I	1.00E-02	I	6.00E-02	I	V	-	0.03	1
Trimethylbenzene, 1,3,5-	108-67-8	NA	I	NA	I	1.00E-02	I	6.00E-02	I	V	-	0.03	1
Trinitrotoluene, 2,4,6- (TNT)	118-96-7	3.00E-02	I	NA	I	5.00E-04	I	NA	--	-	-	0.03	1
Uranium (soluble salts)	NA	NA	(q)	NA	(q)	6.00E-04	(q)	NA	--	-	-	0.01	1
Vanadium	7440-62-2	NA	P	NA	P	5.00E-03	U	1.00E-04	A	-	-	0.01	0.03
Vinyl chloride	75-01-4	1.40E+00	I	8.80E-06	I	3.00E-03	I	8.00E-02	A	V	yes	0.03	1
Xylenes	1330-20-7	NA	I	NA	I	2.00E-01	I	1.00E-01	I	V	-	0.03	1
Zinc	7440-66-6	NA	I	NA	I	3.00E-01	I	NA	I	-	-	0.01	1

Abbreviations:

Sources: A=ATSDR, CE=CAL EPA, E=U.S. Environmental Protection Agency, H=HEAST 1997, I=IRIS, N=NCEA, O=OPP, P=PPRTV, PA=PPRTV Appendix, VA=Vermont Air Pollution Control Regulations, VH=Vermont mg/kg-day - milligram per kilogram-day, µg/m³ - microgram per cubic meter, mg/m³ milligram per cubic meter
CSF₀ - Oral Cancer Slope Factor
IUR - Inhalation Unit Risk

ATTACHMENT 2

VERMONT DEPARTMENT OF HEALTH
TOXICITY & CHEMICAL SPECIFIC VALUES

NA - Not Available/Not Applicable

RfC - Inhalation Reference Concentration

RfD_o - Oral Reference Dose

V - Denotes chemical with a Henry's Law constant greater than 1×10^{-5} atm-m³/mole or a vapor pressure greater than 1 mm Hg

ABSd - Dermal adsorption fraction from soil EPA 2004 Exhibit 3- 4 recommended default values for chemicals not listed 0.1 for semivolatiles, 0.01 for inorganics and 0.03 for volatiles.

GIABS - Gastrointestinal absorption value EPA 2004 Exhibit 4- 1, If ABSGI >50%, 1

Notes:

(a) - CAS Number is for Metallic Thallium

(b) Benomyl - Reflects incorporation of EPA Office of Pesticide Programs Food Quality Protection Act Safety Factor = 10 (for children and females 15-50 years).

(c) Benzene - Most conservative end of range of oral CSFs presented in IRIS citation dated 1/19/00 [1.5E-02 to 5.5E-02 (mg/kg-d)-1].

(d) Benzene - Most conservative end of range of IURs presented in IRIS citation dated 1/19/00 [2.2E-06 to 7.8E-06 (µg/m³)-1].

(e) Cadmium -IRIS,1992 RfD (food) adjusted with Vermont Department of Health Modifying Factor of 10 to account for unquantified carcinogenic potential.

(f) Carbaryl and Propoxur - EPA Office of Pesticide Programs does not employ a long-term or chronic oral RfD as they contend there is rapid recovery from acetylcholinesterase inhibition, because there are multiple other effects and potential concerns regarding extended exposure, assessment of chronic exposure is appropriate.

(g) Chromium (VI) - Considers information presented in OPP (CARC), 2008; IRIS, 2010 Draft; CAL EPA PHG, 2011 (mnoa); NJ DEP, 2009.

(h) Chromium (VI) - The IRIS 9/3/98 IUR of 1.2E-2 (ug/m³)-1 assumes a ratio of Cr(VI) to Cr(III) of 1:6. To obtain a Cr(VI) specific IUR, this value was multiplied by 7.

(i) IRIS citation for Nickel, soluble salts.

(k) Polychlorinated Biphenyls -IRIS high risk and persistence cancer toxicity values employed; noncancer assessment of Total PCBs based on RfD and VF for Aroclor 1254.

(l) n- Propylbenzene - VDH employs IRIS1987 RfD for ethylbenzene as surrogate. PPRTV2009 developed screening chronic provisional values using ethylbenzene value as surrogate.

(m) n-Propylbenzene - VDH employs ATSDR₀₁₀ inhalation chronic MRL for ethylbenzene as a surrogate. PPRTV₀₀₆ develops provisional chronic RfC using older IRIS₉₉₁ RfC for ethylbenzene as surrogate.

(n) Selenium - Consistent with Vermont Air Pollution Control Regulations, Hazardous Ambient Air Standard (HAAS), route to route extrapolation of IRIS oral RfD with 70 kg body weight, inhalation rate of 20 m³/d and an uncertainty factor of 10.

(o) Tetrachloroethylene - Geometric mean of oral CSFs noted in IRIS citation dated 2/10/12 [2.1E-3 (mg/kg-d)-1 based on hepatocellular adenoma/carcinoma & 6E-2 (mg/kg-d)-1 based on mononuclear cell leukemia]

(p) Tetrachloroethylene -Geometric mean of IURs presented in IRIS citation dated 2/10/12 [2.6E-07 (µg/m³)-1 based on hepatocellular adenoma/carcinoma & 1E-05 (µg/m³)-1 based on mononuclear cell leukemia]

(q) Uranium assessment based on non-radiogenic effects only. Assessment employs RfD presented in U.S. EPA National Primary Drinking Water Regulations; Radionuclides; Final Rule 12/7/00.

(u) Health used EPA's IEUBK software and Adult Lead Model to find the soil concentration that would result in 95% probability that no exposed child (or fetus) would exceed the B11 of < 3 ug/dL. Vermont DEC conducted a soil background study and determined the rural background level of lead is 41 mg/kg. Because the soil concentration that would result in a 95% probability that no exposed child (or fetus) would exceed the B11 < 3 ug/dL is less than the Vermont background level, the background level is recommended.

ATTACHMENT 3a
SUMMARY TABLE
2023 RESIDENTIAL SOIL VALUES (RSV) (mg/kg)
INCIDENTAL INGESTION, DERMAL CONTACT AND INHALATION

Chemical Name	CAS No.	Resident Soil Values (mg/kg)									
		Cancer Target Risk = 1x10 ⁻⁶				Noncancer Hazard Quotient = 1				HQ=1 & ILCR=1E ⁻⁶	
		Ingestion	Dermal	Inhalation	Combined	Ingestion	Dermal	Inhalation	Combined	Minimum RSV mg/kg	Endpoint
Acetochlor	34256-82-1	-	-	-	-	1.50E+03	6.42E+03	-	1.22E+03	1.22E+03	nc
Acetone	67-64-1	-	-	-	-	6.75E+04	9.63E+05	-	6.31E+04	6.31E+04	nc
Alachlor	15972-60-8	6.97E+00	1.71E+01	-	4.95E+00	7.50E+01	3.21E+02	-	6.08E+01	4.95E+00	c
Aldrin	309-00-2	2.30E-02	5.64E-02	1.61E+00	1.62E-02	3.00E+00	1.28E+01	-	2.43E+00	1.62E-02	c
Aluminum	7429-90-5	-	-	-	-	7.50E+04	3.21E+06	6.80E+06	7.25E+04	7.25E+04	nc
Antimony	7440-36-0	-	-	-	-	3.00E+01	1.28E+03	4.08E+05	2.93E+01	2.93E+01	nc
Arsenic, Inorganic	7440-38-2	2.60E-01	2.13E+00	3.16E+02	2.32E-01	2.25E+01	3.21E+02	2.04E+04	2.10E+01	2.32E-01	c
Barium	7440-39-3	-	-	-	-	1.50E+04	4.49E+04	-	1.12E+04	1.12E+04	nc
Benomyl	17804-35-2	1.63E+02	4.01E+02	-	1.16E+02	9.75E+02	4.17E+03	-	7.90E+02	1.16E+02	c
Benzene	71-43-2	7.10E+00	5.81E+01	7.84E-01	6.98E-01	3.00E+02	4.28E+03	1.83E+02	1.11E+02	6.98E-01	c
Benzo(a)pyrene	50-32-8	1.02E-01	2.55E-01	1.37E+03	7.28E-02	2.25E+01	7.41E+01	2.72E+03	1.72E+01	7.28E-02	c
Beryllium	7440-41-7	-	-	5.67E+02	5.67E+02	1.50E+02	4.49E+01	2.72E+04	3.45E+01	3.45E+01	nc
Bis(2-chloro-1-methyl ethyl)ethe:	108-60-1	-	-	-	-	3.00E+03	4.28E+04	-	2.80E+03	2.80E+03	nc
Boron	7440-42-8	-	-	-	-	1.50E+04	6.42E+05	2.72E+07	1.46E+04	1.46E+04	nc
Bromate	15541-45-4	5.58E-01	1.37E+01	-	5.36E-01	3.00E+02	1.28E+04	-	2.93E+02	5.36E-01	c
Bromochloromethane	74-97-5	-	-	-	-	9.75E+02	1.39E+04	2.45E+02	1.93E+02	1.93E+02	nc
Bromoxynil	1689-84-5	3.79E+00	9.31E+00	-	2.69E+00	1.13E+03	4.82E+03	-	9.12E+02	2.69E+00	c
Butylbenzene, n-	104-51-8	-	-	-	-	3.75E+03	5.35E+04	-	3.50E+03	3.50E+03	nc
Butylbenzene, sec-	135-98-8	-	-	-	-	7.50E+03	1.07E+05	-	7.01E+03	7.01E+03	nc
Butylbenzene, tert-	98-06-6	-	-	-	-	7.50E+03	1.07E+05	-	7.01E+03	7.01E+03	nc
Cadmium (food)	7440-43-9	-	-	7.56E+02	7.56E+02	7.50E+00	8.03E+01	1.36E+04	6.86E+00	6.86E+00	nc
Carbaryl	63-25-2	4.46E+02	1.10E+03	-	3.17E+02	7.50E+03	3.21E+04	-	6.08E+03	3.17E+02	c
Carbon Disulfide	75-15-0	-	-	-	-	7.50E+03	1.07E+05	6.66E+02	6.08E+02	6.08E+02	nc
Carbon tetrachloride	56-23-5	5.58E+00	4.56E+01	4.02E-01	3.72E-01	3.00E+02	4.28E+03	2.41E+02	1.30E+02	3.72E-01	c
Chlorobenzene	108-90-7	-	-	-	-	5.25E+02	7.49E+03	5.87E+02	2.67E+02	2.67E+02	nc
Chromium (III) (insoluble salts)	16065-83-1	-	-	-	-	1.13E+05	6.26E+04	-	4.02E+04	4.02E+04	nc
Chromium (VI)	18540-29-9	2.04E-01	1.66E-01	9.77E+00	9.06E-02	2.25E+02	2.41E+02	1.36E+05	1.16E+02	9.06E-02	c
Cobalt	7440-48-4	-	-	1.51E+02	1.51E+02	2.25E+01	9.63E+02	8.16E+03	2.19E+01	2.19E+01	nc
Copper	7440-50-8	-	-	-	-	1.07E+04	4.56E+05	-	1.04E+04	1.04E+04	nc
Di (2-ethylhexyl) phthalate	117-81-7	2.79E+01	6.85E+01	5.67E+05	1.98E+01	1.50E+03	6.42E+03	-	1.22E+03	1.98E+01	c
Dibromochloropropane	96-12-8	1.27E-01	1.38E+00	6.33E-03	6.00E-03	1.50E+01	2.14E+02	1.26E+01	6.63E+00	6.00E-03	c
Dibromoethane, 1,2-	106-93-4	1.95E-01	1.60E+00	2.61E-02	2.27E-02	6.75E+02	9.63E+03	1.41E+02	1.15E+02	2.27E-02	c
Dichloroethane, 1,1-	75-34-3	6.85E+01	5.61E+02	2.17E+00	2.10E+00	1.50E+04	2.14E+05	-	1.40E+04	2.10E+00	c
Dichloroethane, 1,2-	107-06-2	4.29E+00	3.51E+01	3.08E-01	2.85E-01	1.50E+03	2.14E+04	5.61E+01	5.39E+01	2.85E-01	c
Dichloroethylene, cis 1,2-	156-59-2	-	-	-	-	1.50E+02	2.14E+03	1.70E+02	7.68E+01	7.68E+01	nc
Dichloroethylene, trans 1,2-	156-60-5	-	-	-	-	1.50E+03	2.14E+04	1.17E+02	1.08E+02	1.08E+02	nc
Dichloropropane, 1,2-	78-87-5	1.06E+01	8.64E+01	1.79E+00	1.51E+00	3.00E+03	4.28E+04	2.65E+01	2.63E+01	1.51E+00	c
Dioxane, 1,4-	123-91-1	3.90E+00	3.20E+01	1.39E+01	2.78E+00	2.25E+03	3.21E+04	2.09E+03	1.05E+03	2.78E+00	c
Ethylbenzene	100-41-4	3.55E+01	2.90E+02	4.16E+00	3.68E+00	3.75E+03	5.35E+04	4.75E+02	4.18E+02	3.68E+00	c
Fluoranthene	206-44-0	-	-	-	-	3.00E+03	9.88E+03	-	2.30E+03	2.30E+03	nc
Fluorene	86-73-7	-	-	-	-	3.00E+03	9.88E+03	-	2.30E+03	2.30E+03	nc
Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6	-	-	-	-	2.25E-01	-	-	2.25E-01	2.25E-01	nc
Hexachlorobenzene	118-74-1	2.44E-01	5.99E-01	3.32E-01	1.14E-01	7.50E-01	3.21E+00	-	6.08E-01	1.14E-01	c
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	4.88E+00	7.99E+01	-	4.60E+00	3.00E+02	8.56E+03	-	2.90E+02	4.60E+00	c
Iron	7439-89-6	-	-	-	-	5.25E+04	2.25E+06	-	5.13E+04	5.13E+04	nc
Isopropylbenzene (cumene)	98-82-8	-	-	-	-	7.50E+03	-	2.65E+02	2.56E+02	2.56E+02	nc
Manganese (non-diet)	7439-96-5	-	-	-	-	1.80E+03	3.08E+03	6.80E+04	1.12E+03	1.12E+03	nc
Mercury (elemental)	7439-97-6	-	-	-	-	-	-	3.13E+00	3.13E+00	3.13E+00	nc

ATTACHMENT 3a
SUMMARY TABLE
2023 RESIDENTIAL SOIL VALUES (RSV) (mg/kg)
INCIDENTAL INGESTION, DERMAL CONTACT AND INHALATION

Chemical Name	CAS No.	Resident Soil Values (mg/kg)									
		Cancer Target Risk = 1x10 ⁻⁶				Noncancer Hazard Quotient = 1				HQ=1 & ILCR=1E ⁻⁶	
		Ingestion	Dermal	Inhalation	Combined	Ingestion	Dermal	Inhalation	Combined	Minimum RSV mg/kg	Endpoint
Methyl ethyl ketone	78-93-3	-	-	-	-	4.50E+04	6.42E+05	2.84E+04	1.70E+04	1.70E+04	nc
Methyl tert-butyl ether (MTBE)	1634-04-4	2.17E+02	1.78E+03	3.19E+01	2.74E+01	7.50E+02	1.07E+04	8.79E+03	6.49E+02	2.74E+01	c
Molybdenum	7439-98-7	-	-	-	-	3.75E+02	1.61E+04	2.72E+06	3.66E+02	3.66E+02	nc
Naphthalene	91-20-3	3.25E+00	6.15E+00	2.72E+00	1.19E+00	1.50E+03	4.94E+03	2.78E+02	2.24E+02	1.19E+00	c
Nickel	7440-02-0	-	-	5.23E+03	5.23E+03	1.50E+03	2.57E+03	1.22E+05	9.40E+02	9.40E+02	nc
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0	-	-	-	-	3.75E+03	2.68E+05	-	3.70E+03	3.70E+03	nc
Perfluorobutane sulfonic acid (PFBS)	375-73-5	-	-	-	-	2.25E+01	9.63E+01	-	1.82E+01	1.82E+01	nc
Pentachlorophenol	87-86-5	9.76E-01	9.59E-01	2.67E+05	4.84E-01	3.75E+02	6.42E+02	-	2.37E+02	4.84E-01	c
Pentaerythritol tetranitrate (PETN)	78-11-5	9.08E+01	2.23E+02	-	6.45E+01	6.75E+02	2.89E+03	-	5.47E+02	6.45E+01	c
Perchlorate	14797-73-0	-	-	-	-	5.25E+01	2.25E+03	-	5.13E+01	5.13E+01	nc
Perfluorohexane sulfonic acid (PFHxS)	355-46-4	-	-	-	-	1.50E+00	6.42E+00	-	1.22E+00	1.22E+00	nc
Perfluorononanoic acid (PFNA)	375-95-1	-	-	-	-	1.50E+00	6.42E+00	-	1.22E+00	1.22E+00	nc
Perfluorooctane sulfonic acid (PFOS)	1763-23-1	-	-	-	-	1.50E+00	6.42E+00	-	1.22E+00	1.22E+00	nc
Perfluorooctanoic acid (PFOA)	335-67-1	5.58E+00	1.37E+01	-	3.96E+00	1.50E+00	6.42E+00	-	1.22E+00	1.22E+00	nc
Polychlorinated biphenyls (PCBs)	1336-36-3	1.95E-01	3.42E-01	1.30E+00	1.14E-01	1.88E+00	5.73E+00	-	1.41E+00	1.14E-01	c
Propoxur (Baygon)	114-26-1	1.11E+02	2.72E+02	-	7.88E+01	3.00E+02	1.28E+03	-	2.43E+02	7.88E+01	c
Propyl benzene, n-	103-65-1	-	-	-	-	7.50E+03	-	2.62E+02	2.53E+02	2.53E+02	nc
Selenium	7782-49-2	-	-	-	-	3.75E+02	1.61E+04	2.45E+06	3.66E+02	3.66E+02	nc
Silver	7440-22-4	-	-	-	-	3.75E+02	6.42E+02	-	2.37E+02	2.37E+02	nc
Tetrachlorodibenzo-p-dioxin, 2,3,7,8- (TCDD)	1746-01-6	3.00E-06	2.46E-05	7.02E-05	2.58E-06	5.25E-05	7.49E-04	1.07E-01	4.90E-05	2.58E-06	c
Tetrachloroethane, 1,1,1,2-	630-20-6	1.50E+01	1.23E+02	1.46E+00	1.32E+00	2.25E+03	3.21E+04	-	2.10E+03	1.32E+00	c
Tetrachloroethylene	127-18-4	3.55E+01	2.90E+02	2.57E+00	2.38E+00	4.50E+02	6.42E+03	1.55E+02	1.13E+02	2.38E+00	c
Thallium (soluble Thallium)	7440-28-0	-	-	-	-	7.50E-01	3.21E+01	-	7.33E-01	7.33E-01	nc
Toluene	108-88-3	-	-	-	-	6.00E+03	8.56E+04	8.07E+02	7.05E+02	7.05E+02	nc
Trichloroethylene (non-moa)	79-01-6	1.10E+01	1.19E+02	2.28E+00	1.86E+00	3.75E+01	5.35E+02	7.55E+00	6.21E+00	1.86E+00	c
Trichloropropane, 1,2,3-	96-18-4	3.40E-03	3.64E-02	-	3.11E-03	7.50E+02	1.07E+04	8.95E+00	8.84E+00	3.11E-03	c
Trimethylbenzene, 1,2,3-	526-73-8	-	-	-	-	7.50E+02	1.07E+04	2.91E+02	2.06E+02	2.06E+02	nc
Trimethylbenzene, 1,2,4-	95-63-6	-	-	-	-	7.50E+02	1.07E+04	2.17E+02	1.66E+02	1.66E+02	nc
Trimethylbenzene, 1,3,5-	108-67-8	-	-	-	-	7.50E+02	1.07E+04	1.81E+02	1.44E+02	1.44E+02	nc
Trinitrotoluene, 2,4,6- (TNT)	118-96-7	1.30E+01	9.99E+01	-	1.15E+01	3.75E+01	5.02E+02	-	3.49E+01	1.15E+01	c
Uranium (soluble salts)	7440-61-1	-	-	-	-	4.50E+01	1.93E+03	-	4.40E+01	4.40E+01	nc
Vanadium	7440-62-2	-	-	-	-	5.25E+00	5.84E+00	1.36E+05	2.77E+00	2.77E+00	nc
Vinyl chloride	75-01-4	2.79E-01	2.28E+00	1.63E-01	9.83E-02	2.25E+02	3.21E+03	1.09E+00	1.08E+00	9.83E-02	c
Xylenes	1330-20-7	-	-	-	-	1.50E+04	2.14E+05	2.57E+02	2.52E+02	2.52E+02	nc
Zinc	7440-66-6	-	-	-	-	2.25E+04	9.63E+05	-	2.20E+04	2.20E+04	nc

Notes:

mg/kg-day - milligram per kilogram-day, µg/m³ - microgram per cubic meter, mg/m³ milligram per cubic meter

c - RSV corresponds to a one in one million Incremental Lifetime Cancer Risk (ILCR)

mmoa - carcinogen identified by U.S. EPA to operate via mutagenic mode of action

nc - RSV corresponds to a noncarcinogenic Hazard Quotient as indicated (HQ=1) based on Hypothetical Young Child Resident scenario

Csat substitution used if soil inhalation screening value greater than Csat. Csat values derived using May 2022 Regional Screening Level Calculator

All cancer-based soil inhalation values were less than respective Csat thus no substitutions

Noncancer-based soil inhalation value above respective Csat thus Csat substitution employed for the following:

Acetone, Carbon Disulfide, Ethylbenzene, Isopropylbenzene (cumene), Mercury (elemental), Methyl ethyl ketone, Methyl tert-butyl ether, n-Propyl benzene, Tetrachloroethylene, Toluene, Trichloroethylene, Trimethyl benzenes, Xylenes.

ATTACHMENT 3b
SUMMARY TABLE
2023 COMMERCIAL SOIL VALUES (CSV) (mg/kg)
INCIDENTAL INGESTION, DERMAL CONTACT AND INHALATION

Chemical Name	CAS No.	Commercial Worker Soil Values (mg/kg)									
		Cancer Target Risk = 1x10 ⁻⁶				Noncancer Hazard Quotient = 1				HQ=1 & ILCR=1E ⁻⁶	
		Ingestion	Dermal	Inhalation	Combined	Ingestion	Dermal	Inhalation	Combined	Minimum CSV mg/kg	Endpoint
Acetochlor	34256-82-1	-	-	-	-	2.04E+04	4.83E+04	-	1.44E+04	1.44E+04	nc
Acetone	67-64-1	-	-	-	-	9.20E+05	7.24E+06	-	8.16E+05	8.16E+05	nc
Alachlor	15972-60-8	4.26E+01	1.01E+02	-	2.99E+01	1.02E+03	2.41E+03	-	7.18E+02	2.99E+01	c
Aldrin	309-00-2	1.40E-01	3.31E-01	9.65E+00	9.76E-02	4.09E+01	9.66E+01	-	2.87E+01	9.76E-02	c
Aluminum	7429-90-5	-	-	-	-	1.02E+06	2.41E+07	2.38E+07	9.42E+05	9.42E+05	nc
Antimony	7440-36-0	-	-	-	-	4.09E+02	9.66E+03	1.43E+06	3.92E+02	3.92E+02	nc
Arsenic, Inorganic	7440-38-2	1.59E+00	1.25E+01	2.59E+03	1.41E+00	3.07E+02	2.41E+03	7.15E+04	2.71E+02	1.41E+00	c
Barium	7440-39-3	-	-	-	-	2.04E+05	3.38E+05	-	1.27E+05	1.27E+05	nc
Benomyl	17804-35-2	9.98E+02	2.36E+03	-	7.01E+02	1.33E+04	3.14E+04	-	9.34E+03	7.01E+02	c
Benzene	71-43-2	4.34E+01	3.41E+02	4.71E+00	4.19E+00	4.09E+03	3.22E+04	4.72E+02	4.18E+02	4.19E+00	c
Benzo(a)pyrene	50-32-8	2.38E+00	4.33E+00	1.85E+04	1.54E+00	3.07E+02	5.57E+02	9.53E+03	1.94E+02	1.54E+00	c
Beryllium	7440-41-7	-	-	4.63E+03	4.63E+03	2.04E+03	3.38E+02	9.53E+04	2.89E+02	2.89E+02	nc
Bis(2-chloro-1-methyl ethyl)ether	108-60-1	-	-	-	-	4.09E+04	3.22E+05	-	3.63E+04	3.63E+04	nc
Boron	7440-42-8	-	-	-	-	2.04E+05	4.83E+06	9.53E+07	1.96E+05	1.96E+05	nc
Bromate	15541-45-4	3.41E+00	8.05E+01	-	3.27E+00	4.09E+03	9.66E+04	-	3.92E+03	3.27E+00	c
Bromochloromethane	74-97-5	-	-	-	-	1.33E+04	1.05E+05	6.29E+02	5.97E+02	5.97E+02	nc
Bromoxynil	1689-84-5	2.32E+01	5.47E+01	-	1.63E+01	1.53E+04	3.62E+04	-	1.08E+04	1.63E+01	c
Butylbenzene, n-	104-51-8	-	-	-	-	5.11E+04	4.02E+05	-	4.53E+04	4.53E+04	nc
Butylbenzene, sec-	135-98-8	-	-	-	-	1.02E+05	8.05E+05	-	9.07E+04	9.07E+04	nc
Butylbenzene, tert-	98-06-6	-	-	-	-	1.02E+05	8.05E+05	-	9.07E+04	9.07E+04	nc
Cadmium (food)	7440-43-9	-	-	6.18E+03	6.18E+03	1.02E+02	6.04E+02	4.77E+04	8.72E+01	8.72E+01	nc
Carbaryl	63-25-2	2.73E+03	6.44E+03	-	1.91E+03	1.02E+05	2.41E+05	-	7.18E+04	1.91E+03	c
Carbon Disulfide	75-15-0	-	-	-	-	1.02E+05	8.05E+05	6.66E+02	6.61E+02	6.61E+02	nc
Carbon tetrachloride	56-23-5	3.41E+01	2.68E+02	2.41E+00	2.23E+00	4.09E+03	3.22E+04	3.99E+02	3.59E+02	2.23E+00	c
Chlorobenzene	108-90-7	-	-	-	-	7.15E+03	5.63E+04	7.56E+02	6.76E+02	6.76E+02	nc
Chromium (III) (insoluble salts)	16065-83-1	-	-	-	-	1.53E+06	4.71E+05	-	3.60E+05	3.60E+05	nc
Chromium (VI)	18540-29-9	4.77E+00	2.82E+00	1.32E+02	1.75E+00	3.07E+03	1.81E+03	4.77E+05	1.14E+03	1.75E+00	c
Cobalt	7440-48-4	-	-	1.24E+03	1.24E+03	3.07E+02	7.24E+03	2.86E+04	2.91E+02	2.91E+02	nc
Copper	7440-50-8	-	-	-	-	1.45E+05	3.43E+06	-	1.39E+05	1.39E+05	nc
Di (2-ethylhexyl) phthalate	117-81-7	1.70E+02	4.02E+02	4.63E+06	1.20E+02	2.04E+04	4.83E+04	-	1.44E+04	1.20E+02	c
Dibromochloropropane	96-12-8	2.98E+00	2.35E+01	6.30E-02	6.15E-02	2.04E+02	1.61E+03	3.24E+01	2.75E+01	6.15E-02	c
Dibromoethane, 1,2-	106-93-4	1.19E+00	9.39E+00	1.57E-01	1.36E-01	9.20E+03	7.24E+04	3.63E+02	3.47E+02	1.36E-01	c
Dichloroethane, 1,1-	75-34-3	4.18E+02	3.29E+03	1.30E+01	1.26E+01	2.04E+05	1.61E+06	-	1.81E+05	1.26E+01	c
Dichloroethane, 1,2-	107-06-2	2.62E+01	2.06E+02	1.85E+00	1.71E+00	2.04E+04	1.61E+05	1.44E+02	1.43E+02	1.71E+00	c
Dichloroethylene, cis 1,2-	156-59-2	-	-	-	-	2.04E+03	1.61E+04	4.37E+02	3.52E+02	3.52E+02	nc
Dichloroethylene, trans 1,2-	156-60-5	-	-	-	-	2.04E+04	1.61E+05	3.01E+02	2.96E+02	2.96E+02	nc
Dichloropropane, 1,2-	78-87-5	6.45E+01	5.08E+02	1.08E+01	9.06E+00	4.09E+04	3.22E+05	6.83E+01	6.81E+01	9.06E+00	c
Dioxane, 1,4-	123-91-1	2.38E+01	1.88E+02	8.37E+01	1.69E+01	3.07E+04	2.41E+05	5.38E+03	4.49E+03	1.69E+01	c
Ethylbenzene	100-41-4	2.17E+02	1.71E+03	2.50E+01	2.21E+01	5.11E+04	4.02E+05	4.75E+02	4.70E+02	2.21E+01	c
Fluoranthene	206-44-0	-	-	-	-	4.09E+04	7.43E+04	-	2.64E+04	2.64E+04	nc
Fluorene	86-73-7	-	-	-	-	4.09E+04	7.43E+04	-	2.64E+04	2.64E+04	nc
Hexachlorobenzene	118-74-1	1.49E+00	3.52E+00	1.99E+00	6.86E-01	1.02E+01	2.41E+01	-	7.18E+00	6.86E-01	c
Hexafluoropropylene oxide dimer acid (HFPO-DA)	13252-13-6	-	-	-	-	3.07E+00	-	-	3.07E+00	3.07E+00	nc
Hexahydro-1,3,5-trinitro-1,3,5-triazine (RDX)	121-82-4	2.98E+01	4.70E+02	-	2.80E+01	4.09E+03	6.44E+04	-	3.84E+03	2.80E+01	c
Iron	7439-89-6	-	-	-	-	7.15E+05	1.69E+07	-	6.86E+05	6.86E+05	nc
Isopropylbenzene (cumene)	98-82-8	-	-	-	-	1.02E+05	-	2.65E+02	2.64E+02	2.64E+02	nc
Manganese (non-diet)	7439-96-5	-	-	-	-	2.45E+04	2.32E+04	2.38E+05	1.14E+04	1.14E+04	nc
Mercury (elemental)	7439-97-6	-	-	-	-	-	-	3.13E+00	3.13E+00	3.13E+00	nc
Methyl ethyl ketone	78-93-3	-	-	-	-	6.13E+05	4.83E+06	2.84E+04	2.70E+04	2.70E+04	nc
Methyl tert-butyl ether (MTBE)	1634-04-4	1.32E+03	1.04E+04	1.92E+02	1.65E+02	1.02E+04	8.05E+04	8.79E+03	4.46E+03	1.65E+02	c
Molybdenum	7439-98-7	-	-	-	-	5.11E+03	1.21E+05	9.53E+06	4.90E+03	4.90E+03	nc

ATTACHMENT 3b
SUMMARY TABLE
2023 COMMERCIAL SOIL VALUES (CSV) (mg/kg)
INCIDENTAL INGESTION, DERMAL CONTACT AND INHALATION

Chemical Name	CAS No.	Commercial Worker Soil Values (mg/kg)									
		Cancer Target Risk = 1x10 ⁻⁶				Noncancer Hazard Quotient = 1				HQ=1 & ILCR=1E ⁻⁶	
		Ingestion	Dermal	Inhalation	Combined	Ingestion	Dermal	Inhalation	Combined	Minimum CSV mg/kg	Endpoint
Naphthalene	91-20-3	1.99E+01	3.61E+01	1.64E+01	7.19E+00	2.04E+04	3.71E+04	7.15E+02	6.78E+02	7.19E+00	c
Nickel	7440-02-0	-	-	4.28E+04	4.28E+04	2.04E+04	1.93E+04	4.29E+05	9.71E+03	9.71E+03	nc
Octahydro-1,3,5,7-tetranitro-1,3,5,7-tetrazocine (HMX)	2691-41-0	-	-	-	-	5.11E+04	2.01E+06	-	4.98E+04	4.98E+04	nc
Pentachlorophenol	87-86-5	5.96E+00	5.63E+00	2.18E+06	2.90E+00	5.11E+03	4.83E+03	-	2.48E+03	2.90E+00	c
Pentaerythritol tetranitrate (PETN)	78-11-5	5.55E+02	1.31E+03	-	3.90E+02	9.20E+03	2.17E+04	-	6.46E+03	3.90E+02	c
Perchlorate	14797-73-0	-	-	-	-	7.15E+02	1.69E+04	-	6.86E+02	6.86E+02	nc
Perfluorobutane sulfonic acid (PFBS)	375-73-5	-	-	-	-	3.07E+02	7.24E+02	-	2.15E+02	2.15E+02	nc
Perfluorohexane sulfonic acid (PFHxS)	355-46-4	-	-	-	-	2.04E+01	4.83E+01	-	1.44E+01	1.44E+01	nc
Perfluorononanoic acid (PFNA)	375-95-1	-	-	-	-	2.04E+01	4.83E+01	-	1.44E+01	1.44E+01	nc
Perfluorooctane sulfonic acid (PFOS)	1763-23-1	-	-	-	-	2.04E+01	4.83E+01	-	1.44E+01	1.44E+01	nc
Perfluorooctanoic acid (PFOA)	335-67-1	3.41E+01	8.05E+01	-	2.39E+01	2.04E+01	4.83E+01	-	1.44E+01	1.44E+01	nc
Polychlorinated biphenyls (PCBs)	1336-36-3	1.19E+00	2.01E+00	7.83E+00	6.83E-01	2.56E+01	4.31E+01	-	1.60E+01	6.83E-01	c
Propoxur (Baygon)	114-26-1	6.77E+02	1.60E+03	-	4.76E+02	4.09E+03	9.66E+03	-	2.87E+03	4.76E+02	c
Propyl benzene, n-	103-65-1	-	-	-	-	1.02E+05	-	2.62E+02	2.61E+02	2.61E+02	nc
Selenium	7782-49-2	-	-	-	-	5.11E+03	1.21E+05	8.58E+06	4.90E+03	4.90E+03	nc
Silver	7440-22-4	-	-	-	-	5.11E+03	4.83E+03	-	2.48E+03	2.48E+03	nc
Tetrachlorodibenzo-p-dioxin, 2,3,7,8- (TCDD)	1746-01-6	1.83E-05	1.44E-04	4.21E-04	1.57E-05	7.15E-04	5.63E-03	2.74E-01	6.33E-04	1.57E-05	c
Tetrachloroethane, 1,1,1,2-	630-20-6	9.17E+01	7.22E+02	8.76E+00	7.91E+00	3.07E+04	2.41E+05	-	2.72E+04	7.91E+00	c
Tetrachloroethylene	127-18-4	2.17E+02	1.71E+03	1.54E+01	1.43E+01	6.13E+03	4.83E+04	1.55E+02	1.51E+02	1.43E+01	c
Thallium (soluble Thallium)	7440-28-0	-	-	-	-	1.02E+01	2.41E+02	-	9.81E+00	9.81E+00	nc
Toluene	108-88-3	-	-	-	-	8.18E+04	6.44E+05	8.07E+02	7.98E+02	7.98E+02	nc
Trichloroethylene (non-moa)	79-01-6	6.41E+01	5.05E+02	7.31E+00	6.47E+00	5.11E+02	4.02E+03	1.94E+01	1.86E+01	6.47E+00	c
Trichloropropane, 1,2,3-	96-18-4	7.95E-02	6.26E-01	-	7.05E-02	1.02E+04	8.05E+04	2.30E+01	2.30E+01	7.05E-02	c
Trimethylbenzene, 1,2,3-	526-73-8	-	-	-	-	1.02E+04	8.05E+04	2.91E+02	2.82E+02	2.82E+02	nc
Trimethylbenzene, 1,2,4-	95-63-6	-	-	-	-	1.02E+04	8.05E+04	2.17E+02	2.12E+02	2.12E+02	nc
Trimethylbenzene, 1,3,5-	108-67-8	-	-	-	-	1.02E+04	8.05E+04	1.81E+02	1.77E+02	1.77E+02	nc
Trinitrotoluene, 2,4,6- (TNT)	118-96-7	7.95E+01	5.87E+02	-	7.00E+01	5.11E+02	3.77E+03	-	4.50E+02	7.00E+01	c
Uranium (soluble salts)	7440-61-1	-	-	-	-	6.13E+02	1.45E+04	-	5.88E+02	5.88E+02	nc
Vanadium	7440-62-2	-	-	-	-	7.15E+01	4.39E+01	4.77E+05	2.72E+01	2.72E+01	nc
Vinyl chloride	75-01-4	1.70E+00	1.34E+01	9.76E-01	5.93E-01	3.07E+03	2.41E+04	2.80E+00	2.79E+00	5.93E-01	c
Xylenes	1330-20-7	-	-	-	-	2.04E+05	1.61E+06	2.57E+02	2.57E+02	2.57E+02	nc
Zinc	7440-66-6	-	-	-	-	3.07E+05	7.24E+06	-	2.94E+05	2.94E+05	nc

Notes:

mg/kg-day - milligram per kilogram-day, µg/m³ - microgram per cubic meter, mg/m³ milligram per cubic meter

c - CSV corresponds to a one in one million Incremental Lifetime Cancer Risk (ILCR)

nc - CSV corresponds to a noncarcinogenic Hazard Quotient as indicated (HQ=1) based on an Adult Worker scenario.

Csat substitution used if soil inhalation screening value greater than Csat. Csat values derived using May 2022 Regional Screening Level Calculator.

All cancer-based soil inhalation values were less than respective Csat thus no substitutions.

Noncancer-based soil inhalation value above respective Csat thus Csat substitution employed for the following:

Acetone, Carbon Disulfide, Ethylbenzene, Isopropylbenzene (cumene), Mercury (elemental), Methyl ethyl ketone, Methyl tert-butyl ether, n-Propyl benzene, Tetrachloroethylene, Toluene, Trichloroethylene, Trimethyl benzenes, Xylenes.

ATTACHMENT 4a
VERMONT DEPARTMENT OF HEALTH
EXPOSURE ASSUMPTIONS, PARAMETER VALUES AND FACTORS
2023 RESIDENTIAL SOIL VALUES

SYMBOL	DEFINITION (units)	VALUE	REFERENCE
RSV	Residential Soil Value (mg/kg)	Chemical-Specific	Attachments 1, 3, 5
RSV _{nc-ing}	Resident, Soil, Noncancer, Ingestion (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{nc-der}	Resident, Soil, Noncancer, Dermal (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{nc-inh}	Resident, Soil, Noncancer, Inhalation (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{nc-comb}	Resident, Soil, Noncancer, Combined Routes of Exposure (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{ca-ing}	Resident, Soil, Cancer, Ingestion (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{ca-der}	Resident, Soil, Cancer, Dermal (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{ca-inh}	Resident, Soil, Cancer, Inhalation (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{ca-comb}	Resident, Soil, Cancer, Combined Routes of Exposure (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{m-ing}	Resident, Soil, Mutagenic, Ingestion (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{m-der}	Resident, Soil, Mutagenic, Dermal (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{m-inh}	Resident, Soil, Mutagenic, Inhalation (mg/kg)	Chemical-Specific	Attachment 5a
RSV _{m-comb}	Resident, Soil, Mutagenic, Combined Routes of Exposure (mg/kg)	Chemical-Specific	Attachment 5a
RfD _o	Chronic Oral Reference Dose (mg/kg-d)	Chemical-Specific	Attachment 2
RfC	Chronic Inhalation Reference Concentration (mg/m ³)	Chemical-Specific	Attachment 2
CSF _o	Oral Cancer Slope Factor (mg/kg-d) ⁻¹	Chemical-Specific	Attachment 2
IUR	Inhalation Unit Risk (μg/m ³) ⁻¹	Chemical-Specific	Attachment 2
THQ	Target Hazard Quotient (unitless)	1.0	See Text
TR	Target Incremental Lifetime Cancer Risk (unitless)	1x10 ⁻⁶	See Text
LT	Lifetime (yrs)	70	EPA 1989
AT _{R-nc}	Averaging Time, Resident, Noncancer (days)	365 x ED _{Yc} = 2190	Calculated
AT _{R-ca}	Averaging Time, Resident, Cancer (days)	365 x ED _{LT} = 25550	Calculated
IR _{Yc}	Soil Ingestion Rate, Young Child _{Birth-<6yrs} (mg/day)	200	EPA 1991
IR _{Oc}	Soil Ingestion Rate, Older Child _{6-<18yrs} (mg/day)	100	EPA 1991
IR _{Birth-<2 yr}	Soil Ingestion Rate, Fine Age Range Child _{Birth-<2yrs} (mg/day)	200	EPA 1991
IR _{2-<6yr}	Soil Ingestion Rate, Fine Age Range Child _{2-<6yrs} (mg/day)	200	EPA 1991
IR _{6-<16yr}	Soil Ingestion Rate, Fine Age Range Child _{6-<16yrs} (mg/day)	100	EPA 1991
IR _{16-<18yr}	Soil Ingestion Rate, Fine Age Range Child _{16-<18yrs} (mg/day)	100	EPA 1991
IR _A	Soil Ingestion Rate, Adult (mg/day)	100	EPA 1991
IFSR _{adj}	Resident Soil Ingestion Rate Factor, Age-adjusted (mg/kg)	65,439	Attachment 5a
IFSM _{R-adj}	Resident Mutagenic Soil Ingestion Rate Factor, Age-adjusted (mg/kg)	250,620	Attachment 5a
SA _{Yc}	Skin Surface Area, Young Child _{Birth-<6yrs} (cm ²)	2336	(a)
SA _{Oc}	Skin Surface Area, Older Child _{6-<18yrs} (cm ²)	4591	(a)
SA _{Birth-<2 yr}	Skin Surface Area, Fine Age Range Child _{Birth-<2yrs} (cm ²)	2028	(a)
SA _{2-<6yr}	Skin Surface Area, Fine Age Range Child _{2-<6yrs} (cm ²)	2490	(a)
SA _{6-<16yr}	Skin Surface Area, Fine Age Range Child _{6-<16yrs} (cm ²)	4407	(a)
SA _{16-<18yr}	Skin Surface Area, Fine Age Range Child _{16-<18yrs} (cm ²)	5512	(a)
SA _A	Skin Surface Area, Adult (cm ²)	6034	(a)
DFSR _{adj}	Soil Dermal Contact Factor, Age-adjusted (mg/kg)	266,522	Attachment 5a

SYMBOL	DEFINITION (units)	VALUE	REFERENCE
DFSM _{R-adj}	Mutagenic Soil Dermal Contact Factor, Age-adjusted (mg/kg)	770,281	Attachment 5a
AD _c	Soil on Skin Adherence Factor, Child (mg/cm ²)	0.2	EPA 2002
AD _A	Soil on Skin Adherence Factor, Adult (mg/cm ²)	0.07	EPA 2002
BW _{YC}	Body Weight, Young Child _{Birth-<6yrs} (kg)	15	(b)
BW _{OC}	Body Weight, Older Child _{6-<18yrs} (kg)	48	(b)
BW _{Birth-<2yr}	Body Weight, Fine Age Range, Child _{Birth-<2yrs} (kg)	10	(b)
BW _{2-<6yr}	Body Weight, Fine Age Range, Child _{2-<6yrs} (kg)	17	(b)
BW _{6-<16yr}	Body Weight, Fine Age Range, Child _{6-<16yrs} (kg)	44	(b)
BW _{16-<18yr}	Body Weight, Fine Age Range, Child _{16-<18yrs} (kg)	67	(b)
BW _A	Body Weight, Adult (kg)	70	EPA 1991
ABS _d	Fraction of chemical absorbed from soil due to dermal contact (unitless)	Chemical-specific	EPA 2004 (Exhibit 3-4); See Text
ABS _{GI}	Fraction of chemical absorbed in gastrointestinal tract (unitless). If ABS _{GI} >50%, a value of 1 (100%) used.	Chemical-specific	EPA 2004 (Exhibit 4-1)
EF _{YC}	Exposure Frequency, Young Child _{Birth-<6yrs} (days/yr)	365	See Text
EF _{OC}	Exposure Frequency, Older Child _{6-<18yrs} (days/yr)	365	See Text
EF _{Birth-<2yr}	Exposure Frequency, Fine Age Range Child _{Birth-<2yrs} (days/yr)	365	See Text
EF _{2-<6yr}	Exposure Frequency, Fine Age Range Child _{2-<6yrs} (days/yr)	365	See Text
EF _{6-<16yr}	Exposure Frequency, Fine Age Range Child _{6-<16yrs} (days/yr)	365	See Text
EF _{16-<18yr}	Exposure Frequency, Fine Age Range Child _{16-<18yrs} (days/yr)	365	See Text
EF _A	Exposure Frequency, Adult (days/yr)	365	See Text
ED _{YC}	Exposure Duration, Young Child _{Birth-<6yrs} (yrs)	6	EPA 1991
ED _{OC}	Exposure Duration, Older Child _{6-<18yrs} (yrs)	12	Calculated
ED _{Birth-<2yr}	Exposure Duration, Fine Age Range Child _{Birth-<2yrs} (yrs)	2	Calculated
ED _{2-<6yr}	Exposure Duration, Fine Age Range Child _{2-<6yrs} (yrs)	4	Calculated
ED _{6-<16yr}	Exposure Duration, Fine Age Range Child _{6-<16yrs} (yrs)	10	Calculated
ED _{16-<18yr}	Exposure Duration, Fine Age Range, Child _{16-<18yrs} (yrs)	2	Calculated
ED _A	Exposure Duration, Adult (yrs)	52	Calculated
ET _{YC}	Exposure Time, Young Child _{Birth-<6yrs} (hours/day)	24	EPA 2022
ET _{OC}	Exposure Time, Older Child _{6-<18yrs} (hours/day)	24	EPA 2022
ET _{Birth-<2yr}	Exposure Time, Fine Age Range Child _{Birth-<2yrs} (hours/day)	24	EPA 2022
ET _{2-<6yr}	Exposure Time, Fine Age Range Child _{2-<6yrs} (hours/day)	24	EPA 2022
ET _{6-<16yr}	Exposure Time, Fine Age Range Child _{6-<16yrs} (hours/day)	24	EPA 2022
ET _{16-<18yr}	Exposure Time, Fine Age Range Child _{16-<18yrs} (hours/day)	24	EPA 2022
ET _A	Exposure Time, Adult (hours/day)	24	EPA 2022
InFSM _{R-adj}	Mutagenic Soil Inhalation Factor, Age-adjusted (days)	42,340	Attachment 5
PEF	Particulate Emission Factor (wind-driven) (m ³ /kg)	1.36 x 10 ⁹	EPA 2022 ^(c)
VF	Volatilization Factor (m ³ /kg)	Chemical-Specific	EPA 2022 ^(d)
RBA	Relative Bioavailability (unitless)	1	See Text
SCMF	Snow Cover Modification Factor (unitless)	(e)	See Text

Notes:

(a) Surface areas derived using information presented in EPA 2011 and Boniol et al., 2007 for sexes combined. Mean of 50th percentile (consistent with EPA 1989 p. 3-39) Total Body Surface Area for each age range of interest developed. Head, hands, forearms, lower legs and feet considered in contact/exposed for all Child age ranges. Consistent with EPA 2004 (p. 3-10), head, hands, forearms and lower legs considered for Adult. Percent of Total Surface Area represented by body parts considered in contact/exposed was calculated (mean across age range of interest).

(b) Average mean annual Body Weight for age range of interest (based on both sexes) derived using information presented in Portier, et al., 2007.

(c) Default value employed in U.S. EPA Regional Screening Levels for Chemical Contaminants at Superfund Sites. (accessed June 2022).

(d) Chemical-specific Volatilization Factors from U.S. EPA Regional Screening Levels for Chemical Contaminants at Superfund Sites. (accessed June 2022).

(e) Snow Cover Modification Factor (SCMF) of 0.7342 applied only to soil inhalation route and only for chemicals that meet “v” criteria (effectively yields exposure frequency of 268 days per yr for this route of exposure for this receptor). SCMF of 1 employed for all other routes and for chemicals that do not meet “v” criteria.

References:

Boniol, et al, 2007. Proportion of skin surface area of children and young adults from 2 to 18 yrs old. J Investig Dermatol 128(2):461-464.

EPA 1989. Risk Assessment Guidance for Superfund. Volume I: Human Health Evaluation Manual (Part A). Interim Final. United States Environmental Protection Agency. Office of Emergency and Remedial Response. Washington, D.C. EPA/540/1-89/002. December 1989.

EPA 1991. Human Health Evaluation Manual, Supplemental Guidance: “Standard Default Exposure Factors”. United States Environmental Protection Agency. . Office of Solid Waste and Emergency Response. OSWER Publication 9285.6-03. March 1991.

EPA 2002. Supplemental Guidance for Developing Soil Screening Levels for Superfund Sites. U.S. Environmental Protection Agency. Washington, D.C. OSWER 9355.4-24-02. December 2002.

EPA 2004. Risk Assessment Guidance for Superfund Volume I: Human Health Evaluation Manual (Part E, Supplemental Guidance for Dermal Risk Assessment) Final. U.S. Environmental Protection Agency. Washington, D.C. OSWER 9285.7-02EP. July 2004.

EPA 2011. Exposure Factors Handbook 2011 Edition (Final). U.S. Environmental Protection Agency. Office of Research and Development. Washington, D.C. EPA/600/R-090/052F. September 2011.

EPA 2022. United States Environmental Protection Agency. Regional Screening Levels for Chemical Contaminants at Superfund Sites. User’s Guide. (accessed November 2022).

Portier et al., 2007. Body weight distributions for risk assessment. Risk Anal 27(1):11-26.

ATTACHMENT 4b
VERMONT DEPARTMENT OF HEALTH
EXPOSURE ASSUMPTIONS, PARAMETER VALUES AND FACTORS
2023 COMMERCIAL WORKER SOIL VALUES

SYMBOL	DEFINITION (units)	VALUE	REFERENCE
CSV	Commercial Worker Soil Value (mg/kg)	Chemical-Specific	Attachments 1, 3, 5b
CSV _{nc-ing}	Commercial Worker, Soil, Noncancer, Ingestion (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{nc-der}	Commercial Worker, Soil, Noncancer, Dermal (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{nc-inh}	Commercial Worker, Soil, Noncancer, Inhalation (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{nc-comb}	Commercial Worker, Soil, Noncancer, Combined Routes of Exposure (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{ca-ing}	Commercial Worker, Soil, Cancer, Ingestion (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{ca-der}	Commercial Worker, Soil, Cancer, Dermal (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{ca-inh}	Commercial Worker, Soil, Cancer, Inhalation (mg/kg)	Chemical-Specific	Attachment 5b
CSV _{ca-comb}	Commercial Worker, Soil, Cancer, Combined Routes of Exposure (mg/kg)	Chemical-Specific	Attachment 5b
RfD _o	Chronic Oral Reference Dose (mg/kg-d)	Chemical-Specific	Attachment 2
RfC	Chronic Inhalation Reference Concentration (mg/m ³)	Chemical-Specific	Attachment 2
CSF _o	Oral Cancer Slope Factor (mg/kg-d) ⁻¹	Chemical-Specific	Attachment 2
IUR	Inhalation Unit Risk (µg/m ³) ⁻¹	Chemical-Specific	Attachment 2
THQ	Target Hazard Quotient (unitless)	1.0	See Text
TR	Target Incremental Lifetime Cancer Risk (unitless)	1x10 ⁻⁶	See Text
LT	Lifetime (yrs)	70	EPA 1989
AT _{R-nc}	Averaging Time, Commercial Worker, Noncancer (days)	365 x ED _w = 9125	Calculated
AT _{R-ca}	Averaging Time, Commercial Worker, Cancer (days)	365 x ED _{LT} = 25550	Calculated
IR _w	Soil Ingestion Rate, Commercial Worker (mg/day)	100	EPA 1991
SA _w	Skin Surface Area, Adult (cm ²)	3527	(a)
AD _w	Soil on Skin Adherence Factor, Adult (mg/cm ²)	0.12	EPA 2014
BW _w	Body Weight, Adult (kg)	70	EPA 1991
ABS _d	Fraction of chemical absorbed from soil due to dermal contact (unitless)	Chemical-specific	EPA 2004 (Exhibit 3-4); See Text
ABS _{GI}	Fraction of chemical absorbed in gastrointestinal tract (unitless). If ABS _{GI} >50%, a value of 1 (100%) used.	Chemical-specific	EPA 2004 (Exhibit 4-1)
EF _w	Exposure Frequency, Ingestion & Dermal Commercial Worker (days/yr)	250	EPA 1991
ET _w	Exposure Time, Adult (hours/day)	10	BLS, 2016
PEF	Particulate Emission Factor (wind-driven) (m ³ /kg)	1.36 x 10 ⁹	EPA 2022 ^(c)
VF	Volatilization Factor (m ³ /kg)	Chemical-Specific	EPA 2022 ^(d)
RBA	Relative Bioavailability (unitless)	1	See Text

Notes:

(a) Surface areas derived using information presented in EPA 2011, Table 7-2; weighted average of mean values for head, hands, and forearms (male and female, 21+yrs)

(b) Average mean annual Body Weight for age range of interest (based on both sexes) derived using information presented in Portier, et al., 2007.

(c) Default value employed in U.S. EPA Regional Screening Levels for Chemical Contaminants at Superfund Sites (accessed June 2022).

(d) Chemical-specific Volatilization Factors from U.S. EPA Regional Screening Levels for Chemical Contaminants at Superfund Sites (accessed June 2022).

References:

BLS, 2016. United States Bureau of Labor Statistics. Division of Labor Force Statistics. Labor Force Statistics from Current Population Survey. Household Data. Annual Average. Last modified February 8, 2017 (accessed 3/28/2017) <https://www.bls.gov/cps/cpsaat19.htm>.

EPA 1989. Risk Assessment Guidance for Superfund. Volume I: Human Health Evaluation Manual (Part A). Interim Final. United States Environmental Protection Agency. Office of Emergency and Remedial Response. Washington, D.C. EPA/540/1-89/002. December 1989.

EPA 1991. Human Health Evaluation Manual, Supplemental Guidance: "Standard Default Exposure Factors". United States Environmental Protection Agency. . Office of Solid Waste and Emergency Response. OSWER Publication 9285.6-03. March 1991.

EPA 2002. Supplemental Guidance for Developing Soil Screening Levels for Superfund Sites. U.S. Environmental Protection Agency. Washington, D.C. OSWER 9355.4-24-02. December 2002.

EPA 2004. Risk Assessment Guidance for Superfund Volume I: Human Health Evaluation Manual (Part E, Supplemental Guidance for Dermal Risk Assessment) Final. U.S. Environmental Protection Agency. Washington, D.C. OSWER 9285.7-02EP. July 2004.

EPA 2011. Exposure Factors Handbook 2011 Edition (Final). U.S. Environmental Protection Agency. Office of Research and Development. Washington, D.C. EPA/600/R-090/052F. September 2011.

EPA 2022. United States Environmental Protection Agency. Regional Screening Levels for Chemical Contaminants at Superfund Sites. User's Guide. May and November 2018 editions. (accessed November 2022).

Portier et al., 2007. Body weight distributions for risk assessment. Risk Anal 27(1):11-26.

ATTACHMENT 5a
VERMONT DEPARTMENT OF HEALTH
ENDPOINT AND PATHWAY SPECIFIC EQUATIONS
2023 RESIDENTIAL SOIL VALUES

Noncancer Residential Soil Value Equations

Ingestion

$$RSV_{nc-ing}(\text{mg/kg}) = \frac{THQ * AT_{R-nc} \left(\frac{365 \text{ days}}{\text{year}} * ED_{YC}(6 \text{ years}) \right) * BW_{YC}(15 \text{ kg})}{EF_{YC} \left(\frac{365 \text{ days}}{\text{year}} \right) * ED_{YC}(6 \text{ years}) * \frac{RBA}{RfD_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)} * IR_{YC} \left(\frac{200 \text{ mg}}{\text{day}} \right) * \frac{10^{-6} \text{ kg}}{1 \text{ mg}}}$$

Dermal

$$RSV_{nc-der}(\text{mg/kg}) = \frac{THQ * AT_{R-nc} \left(\frac{365 \text{ days}}{\text{year}} * ED_{YC}(6 \text{ years}) \right) * BW_{YC}(15 \text{ kg})}{EF_{YC} \left(\frac{365 \text{ days}}{\text{year}} \right) * ED_{YC}(6 \text{ years}) * \frac{1}{\left(RfD_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right) * ABS_{GI} \right)} * SA_{YC} \left(\frac{2336 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right) * ABS_d * \frac{10^{-6} \text{ kg}}{1 \text{ mg}}}$$

Inhalation

$$RSV_{nc-inh}(\text{mg/kg}) = \frac{THQ * AT_{R-nc} \left(\frac{365 \text{ days}}{\text{yr}} * ED_{YC}(6 \text{ years}) \right)}{EF_{YC} \left(\frac{365 \text{ days}}{\text{yr}} \right) * SCMF * ED_{YC}(6 \text{ yrs}) * ET_{YC} \left(\frac{24 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * \frac{1}{RfC \left(\frac{\text{mg}}{\text{m}^3} \right)} * \left(\frac{1}{VF \left(\frac{\text{m}^3}{\text{kg}} \right)} + \frac{1}{PEF \left(\frac{\text{m}^3}{\text{kg}} \right)} \right)}$$

Combined Routes of Exposure

$$RSV_{nc-comb}(\text{mg/kg}) = \frac{1}{\frac{1}{RSV_{nc-ing}} + \frac{1}{RSV_{nc-der}} + \frac{1}{RSV_{nc-inh}}}$$

Carcinogenic Residential Soil Value Equations

Ingestion

$$RSV_{ca-ing}(mg/kg) = \frac{TR * AT_{R-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{CSF_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)^{-1} * RBA * IFS_{R-adj} \left(\frac{65,439 \text{ mg}}{\text{kg}} \right) * \frac{10^{-6} \text{ kg}}{\text{mg}}}$$

Where:

$$IFS_{R-adj} \left(\frac{65,439 \text{ mg}}{\text{kg}} \right) = \frac{EF_{YC} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{YC}(6 \text{ yrs}) * IRS_{YC} \left(\frac{200 \text{ mg}}{\text{day}} \right)}{BW_{YC}(15 \text{ kg})} + \frac{EF_{OC} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{OC}(12 \text{ yrs}) * IRS_{OC} \left(\frac{100 \text{ mg}}{\text{day}} \right)}{BW_{OC}(48 \text{ kg})} + \frac{EF_A \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_A(52 \text{ yrs}) * IRS_A \left(\frac{100 \text{ mg}}{\text{day}} \right)}{BW_A(70 \text{ kg})}$$

Dermal

$$RSV_{ca-der}(mg/kg) = \frac{TR * AT_{R-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{\left(\frac{CSF_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)^{-1}}{ABS_{GI}} \right) * DFS_{R-adj} \left(\frac{266,522 \text{ mg}}{\text{kg}} \right) * ABS_d * \left(\frac{10^{-6} \text{ kg}}{\text{mg}} \right)}$$

Where:

$$DFS_{R-adj} \left(\frac{266,522 \text{ mg}}{\text{kg}} \right) = \frac{EF_{YC} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{YC}(6 \text{ yrs}) * SA_{YC} \left(\frac{2336 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right)}{BW_{YC}(15 \text{ kg})} + \frac{EF_{OC} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{OC}(12 \text{ yrs}) * SA_{OC} \left(\frac{4591 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right)}{BW_{OC}(48 \text{ kg})} + \frac{EF_A \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_A(52 \text{ yrs}) * SA_A \left(\frac{6034 \text{ cm}^2}{\text{day}} \right) * AD_A \left(\frac{0.07 \text{ mg}}{\text{cm}^2} \right)}{BW_A(70 \text{ kg})}$$

Inhalation

$RSV_{ca-inh}(mg/kg)$

$$= \frac{TR * AT_{R-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{IUR(\mu\text{g}/\text{m}^3)^{-1} * \left(\frac{1000 \mu\text{g}}{\text{mg}} \right) * EF_R \left(\frac{365 \text{ days}}{\text{yr}} \right) * SCMF * \left(\frac{1}{VF \left(\frac{\text{m}^3}{\text{kg}} \right)} + \frac{1}{PEF \left(\frac{\text{m}^3}{\text{kg}} \right)} \right) * ED_R(70 \text{ yrs}) * ET_R \left(\frac{24 \text{ hrs}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hrs}} \right)}$$

Combined Routes of Exposure

$$RSV_{ca-comb}(mg/kg) = \frac{1}{\frac{1}{RSV_{ca-ing}} + \frac{1}{RSV_{ca-der}} + \frac{1}{RSV_{ca-inh}}}$$

Carcinogenic via Mutagenic Mode of Action Residential Soil Value Equations

Ingestion

$$RSV_{m-ing}(mg/kg) = \frac{TR * AT_{R-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{CSF_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)^{-1} * RBA * IFSM_{R-adj} \left(\frac{250,620 \text{ mg}}{\text{kg}} \right) * \frac{10^{-6} \text{ kg}}{\text{mg}}}$$

Where:

$$IFSM_{R-adj} \left(\frac{250,620 \text{ mg}}{\text{kg}} \right) = \frac{EF_{\text{Birth-}<2\text{yr}} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{\text{Birth-}<2\text{yr}}(2 \text{ yrs}) * IR_{\text{Birth-}<2\text{yr}} \left(\frac{200 \text{ mg}}{\text{day}} \right) * 10}{BW_{\text{Birth-}<2\text{yrs}}(10 \text{ kg})} + \frac{EF_{2-<6\text{yr}} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{2-<6\text{yr}}(4 \text{ yrs}) * IR_{2-<6\text{yr}} \left(\frac{200 \text{ mg}}{\text{day}} \right) * 3}{BW_{2-<6\text{yrs}}(17 \text{ kg})} + \frac{EF_{6-<16\text{yr}} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{6-<16\text{yr}}(10 \text{ yrs}) * IR_{6-<16\text{yr}} \left(\frac{100 \text{ mg}}{\text{day}} \right) * 3}{BW_{6-<16\text{yr}}(44 \text{ kg})} + \frac{EF_{16-<18\text{yr}} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{16-<18\text{yr}}(2 \text{ yrs}) * IR_{16-<18\text{yr}} \left(\frac{100 \text{ mg}}{\text{day}} \right) * 1}{BW_{16-<18\text{yr}}(67 \text{ kg})} + \frac{EF_A \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_A(52 \text{ yrs}) * IR_A \left(\frac{100 \text{ mg}}{\text{day}} \right) * 1}{BW_A(70 \text{ kg})}$$

Dermal

$$RSV_{m-der}(mg/kg) = \frac{TR * AT_{R-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{\left(\frac{CSF_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)^{-1}}{ABS_{GI}} \right) * DFSM_{R-adj} \left(\frac{770,281 \text{ mg}}{\text{kg}} \right) * ABS_d * \left(\frac{10^{-6} \text{ kg}}{\text{mg}} \right)}$$

Where:

$$DFSM_{R-adj} \left(\frac{770,281 \text{ mg}}{\text{kg}} \right) = \frac{EF_{\text{Birth-}<2\text{yr}} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{\text{Birth-}<2\text{yr}}(2 \text{ yrs}) * SA_{\text{Birth-}<2\text{yr}} \left(\frac{2028 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right) * 10}{BW_{\text{Birth-}<2\text{yrs}}(10 \text{ kg})} + \frac{EF_{2-<6\text{yr}} \left(\frac{365 \text{ days}}{\text{yr}} \right) * ED_{2-<6\text{yr}}(4 \text{ yrs}) * SA_{2-<6\text{yr}} \left(\frac{2490 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right) * 3}{BW_{2-<6\text{yr}}(17 \text{ kg})}$$

$$\begin{aligned}
& + \frac{EF_{6-<16yr} \left(\frac{365 \text{ days}}{yr} \right) * ED_{6-<16yr} (10 \text{ yrs}) * SA_{6-<16yr} \left(\frac{4407 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right) * 3}{BW_{6-<16yr} (44 \text{ kg})} \\
& + \frac{EF_{16-<18yr} \left(\frac{365 \text{ days}}{yr} \right) * ED_{16-<18yr} (2 \text{ yrs}) * SA_{16-<18yr} \left(\frac{5512 \text{ cm}^2}{\text{day}} \right) * AD_C \left(\frac{0.2 \text{ mg}}{\text{cm}^2} \right) * 1}{BW_{16-<18yr} (67 \text{ kg})} \\
& + \frac{EF_A \left(\frac{365 \text{ days}}{yr} \right) * ED_A (52 \text{ yrs}) * SA_A \left(\frac{6034 \text{ cm}^2}{\text{day}} \right) * AD_A \left(\frac{0.07 \text{ mg}}{\text{cm}^2} \right) * 1}{BW_A (70 \text{ kg})}
\end{aligned}$$

Inhalation

$$RSV_{m-inh} (\text{mg/kg}) = \frac{TR * AT_{R-ca} \left(\frac{365 \text{ days}}{yr} * LT (70 \text{ yrs}) \right)}{IUR (\mu\text{g}/\text{m}^3)^{-1} * \left(\frac{1000 \mu\text{g}}{\text{mg}} \right) * SCMF * \left(\frac{1}{VF \left(\frac{\text{m}^3}{\text{kg}} \right)} + \frac{1}{PEF \left(\frac{\text{m}^3}{\text{kg}} \right)} \right) * InFSM_{R-adj} (42,340 \text{ days})}$$

Where:

$$\begin{aligned}
& InFSM_{R-adj} (42,340 \text{ days}) = \\
& [ET_{\text{Birth-<2yr}} \left(\frac{24 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * EF_{\text{Birth-<2yr}} \left(\frac{365 \text{ days}}{yr} \right) * ED_{\text{Birth-<2yr}} (2 \text{ yrs}) * 10] \\
& + \\
& [ET_{2-<6yr} \left(\frac{24 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * EF_{2-<6yr} \left(\frac{365 \text{ days}}{yr} \right) * ED_{2-<6yr} (4 \text{ yrs}) * 3] + \\
& [ET_{6-<16yr} \left(\frac{24 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * EF_{6-<16yr} \left(\frac{365 \text{ days}}{yr} \right) * ED_{6-<16yr} (10 \text{ yrs}) * 3] + \\
& [ET_{16-<18yr} \left(\frac{24 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * EF_{16-<18yr} \left(\frac{365 \text{ days}}{yr} \right) * ED_{16-<18yr} (2 \text{ yrs}) * 1] + \\
& [ET_A \left(\frac{24 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * EF_A \left(\frac{365 \text{ days}}{yr} \right) * ED_A (52 \text{ yrs}) * 1]
\end{aligned}$$

Combined Routes of Exposure

$$RSV_{m-comb} (\text{mg/kg}) = \frac{1}{\frac{1}{RSV_{m-ing}} + \frac{1}{RSV_{m-der}} + \frac{1}{RSV_{m-inh}}}$$

ATTACHMENT 5b
VERMONT DEPARTMENT OF HEALTH
ENDPOINT AND PATHWAY SPECIFIC EQUATIONS
2023 COMMERCIAL WORKER SOIL VALUES

Noncarcinogenic Commercial Worker Soil Value Equations

Ingestion

$$CSV_{nc-ing}(mg/kg) = \frac{THQ * AT_{W-nc} \left(\frac{365 \text{ days}}{yr} * ED_W(30 \text{ yrs}) \right) * BW_W(70 \text{ kg})}{EF_W \left(\frac{250 \text{ days}}{yr} \right) * ED_W(30 \text{ yrs}) * \frac{RBA}{RfD_0 \left(\frac{mg}{kg-day} \right)} * IR_W \left(\frac{100 \text{ mg}}{day} \right) * \frac{10^{-6} \text{ kg}}{1 \text{ mg}}}$$

Dermal

$$CSV_{nc-der}(mg/kg) = \frac{THQ * AT_{W-nc} \left(\frac{365 \text{ days}}{yr} * ED_W(30 \text{ yrs}) \right) * BW_W(70 \text{ kg})}{EF_W \left(\frac{250 \text{ days}}{yr} \right) * ED_W(30 \text{ yrs}) * \frac{1}{\left(RfD_0 \left(\frac{mg}{kg-day} \right) * ABS_{GI} \right)} * SA_W \left(\frac{3527}{day} \right) * AD_W \left(\frac{0.12 \text{ mg}}{cm^2} \right) * ABS_d * \frac{10^{-6} \text{ kg}}{1 \text{ mg}}}$$

Inhalation

$$CSV_{nc-inh}(mg/kg) = \frac{THQ * AT_{W-nc} \left(\frac{365 \text{ days}}{yr} * ED_W(30 \text{ yrs}) \right)}{EF_W \left(\frac{250 \text{ days}}{yr} \right) * ED_W(30 \text{ yrs}) * ET_W \left(\frac{10 \text{ hours}}{day} * \frac{1 \text{ day}}{24 \text{ hours}} \right) * \frac{1}{RfC \left(\frac{mg}{m^3} \right)} * \left(\frac{1}{VF \left(\frac{m^3}{kg} \right)} + \frac{1}{PEF \left(\frac{m^3}{kg} \right)} \right)}$$

Combined Routes of Exposure

$$CSV_{nc-comb}(mg/kg) = \frac{1}{\frac{1}{CSV_{nc-ing}} + \frac{1}{CSV_{nc-der}} + \frac{1}{CSV_{nc-inh}}}$$

Carcinogenic Commercial Worker Soil Value Equations

Ingestion

$$CSV_{ca-ing}(mg/kg) = \frac{TR * AT_{W-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{CSF_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)^{-1} * RBA * \left(\frac{EF_W \left(\frac{250 \text{ days}}{\text{yr}} \right) * ED_W(30 \text{ yrs}) * IRS_W \left(\frac{100 \text{ mg}}{\text{day}} \right)}{BW(70 \text{ kg})} \right) * \frac{10^{-6} \text{ kg}}{\text{mg}}}$$

Dermal

$$CSV_{ca-der}(mg/kg) = \frac{TR * AT_{W-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{\left(\frac{CSF_0 \left(\frac{\text{mg}}{\text{kg} - \text{day}} \right)^{-1}}{ABS_{GI}} \right) * \left(\frac{EF_W \left(\frac{250 \text{ days}}{\text{yr}} \right) * ED_W(30 \text{ yrs}) * SA_W \left(\frac{3527 \text{ cm}^2}{\text{day}} \right) * AD_W \left(\frac{0.12 \text{ mg}}{\text{cm}^2} \right)}{BW(70 \text{ kg})} \right) * ABS_d * \left(\frac{10^{-6} \text{ kg}}{\text{mg}} \right)}$$

Inhalation

$$CSV_{ca-inh}(mg/kg) = \frac{TR * AT_{W-ca} \left(\frac{365 \text{ days}}{\text{yr}} * LT(70 \text{ yrs}) \right)}{IUR \left(\frac{\mu\text{g}}{\text{m}^3} \right)^{-1} * \left(\frac{1000 \mu\text{g}}{\text{mg}} \right) * EF_W \left(\frac{250 \text{ days}}{\text{yr}} \right) * \left(\frac{1}{VF \left(\frac{\text{m}^3}{\text{kg}} \right)} + \frac{1}{PEF \left(\frac{\text{m}^3}{\text{kg}} \right)} \right) * ED_W(30 \text{ yrs}) * ET_W \left(\frac{10 \text{ hours}}{\text{day}} * \frac{1 \text{ day}}{24 \text{ hours}} \right)}$$

Combined Routes of Exposure

$$CSV_{ca-comb}(mg/kg) = \frac{1}{\frac{1}{CSV_{ca-ing}} + \frac{1}{CSV_{ca-der}} + \frac{1}{CSV_{ca-inh}}}$$