

1-(2-tert.-Butyl cyclohexyloxy)-2-butanol

| CAS-No.: | 139504-68-0 The scope of this Standard includes, but is not limited to the CAS number(s) indicated above; any other CAS number(s) used to identify this fragrance ingredient should be considered in scope as well. |
|-----------|--|
| Synonyms: | 1-(2-tertButyl cyclohexyloxy)-2-butanol 1-(2-tert-Butylcyclohexyl)oxybutan-2-ol 1-[(2-tert-Butylcyclohexyl)oxy]butan-2-ol 1-(2-tButylcyclohexyl)oxybutan-2-ol 1-[(2-t-Butylcyclohexyl)oxy]butan-2-ol Amber Core (commercial name) Coramber (commercial name) |

| History: F | Publication date: | 2023 (Amendment 51) | Previous Publications: | Not applicable |
|------------|-------------------|---------------------|---------------------------|----------------|
|------------|-------------------|---------------------|---------------------------|----------------|

| Implementation | For new creation*: | March 30, 2024 |
|----------------|---|---------------------------------|
| dates: | For existing creation*: | October 30, 2025 |
| | *These dates apply to the supply of fragrance mixtu | res (formulas) only, not to the |
| | finished consumer products in the marketplace. | |

RECOMMENDATION:

RESTRICTION

| MAXIMUM ACCEPTABLE CONCENTRATIONS IN THE FINISHED PRODUCT (%): | | | | |
|--|---------|--------------|--------|--|
| Category 1 | 0.24 % | Category 7A | 2.7 % | |
| Category 2 | 0.071 % | Category 7B | 2.7 % | |
| Category 3 | 1.4 % | Category 8 | 0.11 % | |
| Category 4 | 1.3 % | Category 9 | 2.6 % | |
| Category 5A | 0.34 % | Category 10A | 4.5 % | |
| Category 5B | 0.34 % | Category 10B | 9.3 % | |
| Category 5C | 0.34 % | Category 11A | 0.11 % | |
| Category 5D | 0.11 % | Category 11B | 0.11 % | |

1-(2-tert.-Butyl cyclohexyloxy)-2-butanol

| Category 6 | 0.75 % | Category 12 | No restriction |
|-----------------------------------|-----------------|---|----------------|
| | | | |
| FLAVOR REQUIREMEN | ITS: | Due to the possible ingestion of small amounts of fragrance ingredients from their use in products in Categories 1 and 6, materials must not only comply with IFRA Standards but must also be recognized as safe as a flavoring ingredient as defined by the IOFI Code of Practice (www.iofi.org). For more details see chapter 1 of the Guidance for the use of IFRA Standards. | |
| | | | |
| CONTRIBUTIONS FROM OTHER SOURCES: | | NONE TO CONSIDER BEYOND TRACES (SEE ALSO THE SECTION ON CONTRIBUTIONS FROM OTHER SOURCES IN CHAPTER 1 OF THE GUIDANCE FOR THE USE OF IFRA STANDARDS) | |
| | | | |
| INTRINSIC PROPER MANAGEMENT: | TY DRIVING RISK | DERMAL SENSITIZATION | N AND SYSTEMIC |

RIFM SUMMARIES:

Maximum acceptable concentrations are based on a comprehensive safety assessment, considering various endpoints. Depending on the outcome of the safety assessment, it might be one or more endpoint(s) that will drive the derivation of the concentration levels. If more than one endpoint is of relevance, the maximum acceptable concentrations for each product category are derived from comparing maximum permitted level per endpoint consideration (e.g. dermal sensitization and/or systemic toxicity). Such maximum acceptable concentrations correspond to the lowest level obtained per category.

Additional information is available in the RIFM safety assessment for 1-(2-tert.-Butyl cyclohexyloxy)-2butanol, which can be downloaded from the RIFM Fragrance Material Safety Assessment Center: http://fragrancematerialsafetyresource.elsevier.com/.

EXPERT PANEL FOR FRAGRANCE SAFETY RATIONALE / CONCLUSION:

The Expert Panel for Fragrance Safety reviewed all the available data for 1-(2-tert.-Butyl cyclohexyloxy)-2butanol and recommends the concentrations for the 12 different product categories, which are the maximum acceptable concentrations of 1-(2-tert.-Butyl cyclohexyloxy)-2-butanol in the various product categories.

REFERENCES:

The IFRA Standard on 1-(2-tert.-Butyl cyclohexyloxy)-2-butanol is based on at least one of the following publications:

• The RIFM Safety Assessment on 1-(2-tert.-Butyl cyclohexyloxy)-2-butanol if available at the RIFM Fragrance Material Safety Assessment Center: http://fragrancematerialsafetyresource.elsevier.com

• Api A.M., Belsito D., Bruze M., Cadby P., Calow P., Dagli M. L., Dekant W., Dent M., Ellis G., Fryer A. D.,

1-(2-tert.-Butyl cyclohexyloxy)-2-butanol

Fukayama M., Griem P., Hickey C., Kromidas L., Lalko J., Liebler D.C., Miyachi Y., Politano V.T., RenskersK., Ritacco G., Salvito D., Schultz T.W., Sipes I. G., Smith B., Vitale D., Wilcox D.K. (2015). Criteria for theResearch Institute for Fragrance Materials, Inc. (RIFM) safety evaluation process for fragrance ingredients.FoodChemToxicol.2015Aug;82Suppl:S1-S19(http://fragrancematerialsafetyresource.elsevier.com/sites/default/files/Criteria_Document_Final.pdf).

• Salvito D.T., Senna R. J., Federle T.W. (2002). A framework for prioritizing fragrance materials for aquatic risk assessment. Environ Toxicol Chem. 2002;21:1301-1308 (https://www.ncbi.nlm.nih.gov/pubmed/12069318).

Additional information on the application of IFRA Standards is available in the Guidance for the use of IFRA Standards, publicly available at www.ifrafragrance.org.