

Trade name : PIPERONYL BUTOXIDE ULTRA

This material complies with the WHO Specification 33/TC (September 2011)
and FAO Specification 33/TC (September 2011) for Piperonyl Butoxide

CHEMICAL IDENTIFICATION

Nomenclature

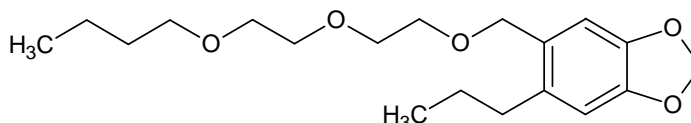
Common name : Piperonyl Butoxide
 IUPAC : 5-[2-(2-butoxyethoxy) ethoxymethyl]-6-propyl-1,3-benzodioxole
 CA name : 5-[[2-(2-Butoxyethoxy)ethoxy]methyl]-6-propyl-1,3-benzodioxole
 Others : (Butylcarbityl) (6-propyl-piperonyl) ether
 PBO

CAS number : 51-03-6

EC number : 200-076-7

Molecular formula : C₁₉H₃₀O₅

Molecular weight : 338.40

Structural formula : 

PRODUCT SPECIFICATIONS

Chemical assay (ENDQC-3 method by HRGC) :

Piperonyl Butoxide	:	94.0 % a/a min.*
Butyl Carbitol®	:	2.0 % a/a max.
Dipiperonyl methane	:	4.0 % a/a max.
Dipiperonyl ether	:	2.0 % a/a max.

Acidity number (ENDQC-6 method) : 0.1 mg KOH/g max.
Density (ENDQC-4 method) : 1.060 (±0.005) g/cm³ at 20°C
Colour (ENDQC-7 method) : 2.5 max (Gardner scale)
Appearance (visual method) : Pale yellow oily liquid

* The analytical result expressed in a/a is equivalent to the w/w analysis.

NOTE: Endura reserves the right to make amendments to the data reported in this sheet either in compliance to any official body's updating or in agreement with Internal Company decisions. The reported information is, to the best of our knowledge, as accurate and complete as possible and is given in good faith but without warranty from our part. Any use for Registration purposes must be authorised by Endura in advance.

TECHNICAL DATA SHEET



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ADDITIONAL INFORMATION

Odour : Mildly aromatic

Refractive index : 1.502 (± 0.005)

Boiling point : 203 °C at 278 Pa

Vapour pressure : 2.11×10^{-7} hPa at 60°C

Flash point : 179.2 °C

Solubility : Practically insoluble in water
Soluble in all common organic solvents including mineral oils and petroleum distillates

Partition coefficient (Log Pow) : 4.8

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