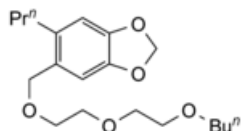




Edition: BP 2025 (Ph. Eur. 11.6 update)

Piperonyl Butoxide

[General Notices](#)



$C_{19}H_{30}O_5$ 338.4 51-03-6

Action and use

Insecticide.

DEFINITION

Piperonyl Butoxide is 5-[2-(2-butoxyethoxy)ethoxymethyl]-6-propyl-1,3-benzodioxole. It contains not less than 94.0% and not more than 102.0% of $C_{19}H_{30}O_5$.

CHARACTERISTICS

A colourless to light yellow, oily liquid.

Very slightly soluble in [water](#); miscible with [ethanol \(96%\)](#) with [ether](#) and with petroleum oils.

IDENTIFICATION

The [infrared absorption spectrum](#), [Appendix II A](#), is concordant with the *reference spectrum* of Piperonyl Butoxide ([RS 478](#)).

TESTS

[Refractive index](#)

1.494 to 1.504, [Appendix V E](#).

[Weight per mL](#)

1.050 to 1.065 g, [Appendix V G](#).

[Sulfated ash](#)

Related substances

Carry out the method for [gas chromatography, Appendix III B](#), using the following solutions in n-[hexane](#).

- (1) 1.2% w/v of the substance being examined.
- (2) Dilute 1 volume of solution (1) to 100 volumes and dilute 1 volume of the resulting solution to 10 volumes.
- (3) Dilute 1 volume of solution (2) to 20 volumes.
- (4) 0.16% w/v of [piperonyl butoxide impurity standard BPCRS](#).

CHROMATOGRAPHIC CONDITIONS

- (a) Use a fused silica capillary column (25 m × 0.32 mm) bonded with a 0.52-µm layer of 100% dimethylpolysiloxane (DB1 is suitable).
- (b) Use [helium](#) as the carrier gas at 1 mL per minute.
- (c) Use the temperature gradient described below.
- (d) Use an inlet temperature of 250°.
- (e) Use a flame ionisation detector at a temperature of 300°.
- (f) Use a split ratio of 1:10.
- (g) Inject 1 µL of each solution.

Time (Minutes)	Temperature °C	Comment
0-2	70	isothermal
2-45	70→285	linear gradient
45-60	285	isothermal

SYSTEM SUITABILITY

The test is not valid unless:

the chromatogram obtained with solution (4) closely resembles the reference chromatogram supplied with [piperonyl butoxide impurity standard BPCRS](#);

in the chromatogram obtained with solution (3), the [signal-to-noise ratio](#) of the principal peak is at least 10;

the [resolution](#) between piperonyl butoxide and dipiperonyl [methane](#) is at least 2.

LIMITS

In the chromatogram obtained with solution (1):

determine the amount of each [secondary peak](#) by [normalisation](#).

the area of any peak corresponding to dipiperonyl [methane](#) is not greater than 2.0%;

the area of any peak corresponding to diethylene glycol butyl ether is not greater than 2.0%;

the area of any peak corresponding to dipiperonyl ether is not greater than 1.5%;

the area of any peak corresponding to dihydrosafrole is not greater than 50 ppm;

the area of any other [secondary peak](#) is not greater than 0.5%;

the sum of the areas of any other [secondary peaks](#) is not greater than 2.5%.

Disregard any peak with an area less than principal peak in solution (2) (0.1%) with the exception of dihydrosafrole.

ASSAY

Carry out the method for [gas chromatography](#), [Appendix III B](#), using the following solutions in [hexane](#).

- (1) 0.16% w/v of the substance being examined.
- (2) 0.16% w/v of [piperonyl butoxide BPCRS](#).
- (3) 0.16% w/v of [piperonyl butoxide impurity standard BPCRS](#).

CHROMATOGRAPHIC CONDITIONS

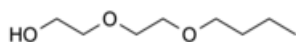
The chromatographic conditions described under Related substances may be used.

SYSTEM SUITABILITY

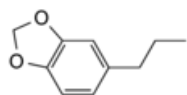
The test is not valid unless the chromatogram obtained with solution (3) closely resembles the reference chromatogram supplied with [piperonyl butoxide impurity standard BPCRS](#) and the [resolution](#) between piperonyl butoxide and dipiperonyl methane is at least 2.0.

Calculate the content of $C_{19}H_{30}O_5$ using the declared content of $C_{19}H_{30}O_5$ in [piperonyl butoxide BPCRS](#).

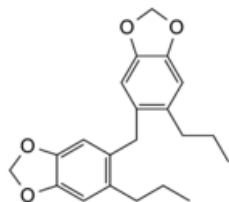
IMPURITIES



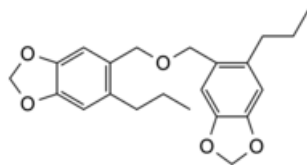
- A. 2-(2-butoxyethoxy)ethanol, (diethylene glycol butyl ether);



- B. 5-propyl-1,3-benzodioxole, 1,2-methylenedioxy-4-propylbenzene, (dihydrosafrole);



- C. [1,1'- di-(6-propyl-benzo-[1,3]-dioxol-5-yl)methane], (dipiperonyl methane);



- D. [di-((6-propyl-benzo-[1,3]dioxol-5-yl)methyl)ether] (dipiperonyl ether).