

Benzoyl Peroxide

Other names:

- Abcure S-40-25
- Acne-Aid Cream
- Acnegel
- Akneroxid 5
- Akneroxid L
- Akneroxide L
- Asidopan
- Aztec BPO
- BPO
- BZF-60
- Benzac
- Benzac W
- Benzagel 10
- Benzaknen
- Benzaknew
- Benzoic acid, peroxide
- Benzol peroxide
- Benzoyl peroxyde
- Benzoylperoxid
- Brevoxyl
- Cadet
- Cadet BPO 78W
- Cadox
- Cadox 40E
- Cadox B
- Cadox BS
- Clear By Design
- Clearasil BP acne treatment
- Clearasil benzoyl peroxide lotion
- Cuticura acne cream
- Debroxide
- Dermoxyl
- Desanden
- Desquam E
- Dibenzoylperoxid
- Dibenzoylperoxyde
- Diphenylglyoxal peroxide
- Dry and Clear
- Duresthin 5
- Eloxyl

Epi-Clear
Fostex
G20
Garox
Incidol
Loroxide
Lucidol
Lucidol (peroxide)
Lucidol 75FP
Lucidol 78
Lucidol B 50
Lucidol G 20
Luperco
Luperco AA
Luperco AST
Luperox FL
Luzidol
Mytolac
NA 2085 (DOT)
NSC 675
Nayper B and BO
Nayper BO
Nericur
Norox bzp-250
Norox bzp-C-35
Novadelox
Nyper B
Nyper BMT
OXY WASH
OXY-10
Oxy 5
Oxy-L
Oxylite
Panoxyl
Perkadox 20S
Perossido di benzoile
Peroxide, dibenzoyl
Peroxyde de benzoyle
Peroxyderm
Peroxydex
Persa-Gel
Persadox
Preoxydex

Quinolor compound

Resdan Akne

Sanoxit

Superox

Theraderm

Topex

Triaz

UN 2085 (DOT)

UN 2086

UN 2088

Vanoxide

W 75

Xerac

Xerac BP 10

Xerac BP 5

acetoxyl

benoxyl

benzoperoxide

benzoyl superoxide

dibenzoyl peroxide

Inchi: InChI=1S/C14H10O4/c15-13(11-7-3-1-4-8-11)17-18-14(16)12-9-5-2-6-10-12/h1-10H

InchiKey: OMPJBNCRMGITSC-UHFFFAOYSA-N

Formula: C14H10O4

SMILES: O=C(OOC(=O)c1cccc1)c1cccc1

Mol. weight [g/mol]: 242.23

CAS: 94-36-0

Physical Properties

Property code	Value	Unit	Source
chs	-6568.70 ± 4.60	kJ/mol	NIST Webbook
chs	-6568.90 ± 8.40	kJ/mol	NIST Webbook
chs	-6546.30 ± 6.30	kJ/mol	NIST Webbook
gf	-176.02	kJ/mol	Joback Method
hf	-281.70 ± 6.20	kJ/mol	NIST Webbook
hf	-272.00 ± 5.40	kJ/mol	NIST Webbook
hfs	-392.00 ± 6.30	kJ/mol	NIST Webbook
hfs	-369.40 ± 8.40	kJ/mol	NIST Webbook
hfs	-369.00 ± 4.60	kJ/mol	NIST Webbook
hfus	25.67	kJ/mol	Joback Method
hsub	97.90 ± 2.50	kJ/mol	NIST Webbook

hsub	98.00 ± 3.00	kJ/mol	NIST Webbook
hvap	69.62	kJ/mol	Joback Method
log10ws	-3.76		Crippen Method
logp	2.615		Crippen Method
mcvol	175.480	ml/mol	McGowan Method
pc	3059.17	kPa	Joback Method
tb	725.66	K	Joback Method
tc	970.13	K	Joback Method
tf	444.70	K	Joback Method
tt	377.75	K	Thermal behavior of benzoyl peroxide mixed with NaOH solution
vc	0.651	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	514.89	J/mol×K	970.13	Joback Method
cpg	454.39	J/mol×K	725.66	Joback Method
cpg	467.35	J/mol×K	766.40	Joback Method
cpg	479.12	J/mol×K	807.15	Joback Method
cpg	489.73	J/mol×K	847.89	Joback Method
cpg	499.20	J/mol×K	888.64	Joback Method
cpg	507.58	J/mol×K	929.38	Joback Method
dvisc	0.0001199	Paxs	725.66	Joback Method
dvisc	0.0009967	Paxs	444.70	Joback Method
dvisc	0.0005919	Paxs	491.53	Joback Method
dvisc	0.0003849	Paxs	538.35	Joback Method
dvisc	0.0002681	Paxs	585.18	Joback Method
dvisc	0.0001971	Paxs	632.01	Joback Method
dvisc	0.0001511	Paxs	678.83	Joback Method
hfust	22.59	kJ/mol	378.00	NIST Webbook
hsubt	89.70 ± 4.20	kJ/mol	303.00	NIST Webbook

Sources

Thermal behavior of benzoyl peroxide mixed with NaOH solution:
Joback Method:

<https://www.doi.org/10.1016/j.tca.2018.10.003>

https://en.wikipedia.org/wiki/Joback_method

McGowan Method:

<http://link.springer.com/article/10.1007/BF02311772>

NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=C94360&Units=SI>
Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>
Crippen Method: https://www.cheméo.com/doc/models/crippen_log10ws

Legend

chs: Standard solid enthalpy of combustion
cpg: Ideal gas heat capacity
dvisc: Dynamic viscosity
gf: Standard Gibbs free energy of formation
hf: Enthalpy of formation at standard conditions
hfs: Solid phase enthalpy of formation at standard conditions
hfus: Enthalpy of fusion at standard conditions
hfust: Enthalpy of fusion at a given temperature
hsub: Enthalpy of sublimation at standard conditions
hsubt: Enthalpy of sublimation at a given temperature
hvap: Enthalpy of vaporization at standard conditions
log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
mcvol: McGowan's characteristic volume
pc: Critical Pressure
tb: Normal Boiling Point Temperature
tc: Critical Temperature
tf: Normal melting (fusion) point
tt: Triple Point Temperature
vc: Critical Volume

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