

Ketoprofen

Other names:

(+)«alpha»-(3-benzoylphenyl)propionic acid ((+)-ketoprofen)
(+)«alpha»-(3-benzoylphenyl)propionic acid ((+)-ketoprofen)
(. +/-.)-3-Benzoyl-«alpha»-methylbenzeneacetic acid
(. +/-.)-3-Benzoyl-«alpha»-methylbenzeneacetic acid
(. +/-.)-m-Benzoylhydratropic acid
19583 RP
2-(3-Benzoylphenyl)propionic acid
2-(m-Benzoylphenyl)propionic acid
3-Benzoyl-«alpha»-methylbenzeneacetic acid
3-Benzoyl-«alpha»-methylbenzeneacetic acid
3-Benzoylhydratropic acid
Alrheumat
Alrheumun
Aneol
Benzeneacetic acid, 3-benzoyl-«alpha»-methyl-
Benzeneacetic acid, 3-benzoyl-«alpha»-methyl-
Capisten
Dexal
Epatec
Fastum
Iso-K
Kefenid
Ketoprofene
Ketopron
Lertus
Menamin
Meprofen
Orudis
Orugesic
Oruvail
Oscorel
Profenid
Propionic acid, 2-(3-benzoylphenyl)-
RP-19583
RU 4733
Racemic ketoprofen
Toprec
Toprek
m-Benzoylhydratropic acid

Inchi:

InChI=1S/C16H14O3/c1-11(16(18)19)13-8-5-9-14(10-13)15(17)12-6-3-2-4-7-12/h2-11H,1

InchiKey: DKYWVDODHFEZIM-UHFFFAOYSA-N
Formula: C₁₆H₁₄O₃
SMILES: CC(C(=O)O)c1cccc(C(=O)c2ccccc2)c1
Mol. weight [g/mol]: 254.28
CAS: 22071-15-4

Physical Properties

Property code	Value	Unit	Source
gf	-98.07	kJ/mol	Joback Method
hf	-294.65	kJ/mol	Joback Method
hfus	28.65	kJ/mol	Joback Method
hvap	86.21	kJ/mol	Joback Method
log10ws	-3.21		Aqueous Solubility Prediction Method
log10ws	-3.45		Aqueous Solubility Prediction Method
logp	3.106		Crippen Method
mcvol	197.790	ml/mol	McGowan Method
pc	2826.33	kPa	Joback Method
rinpol	2185.00		NIST Webbook
tb	823.30	K	Joback Method
tc	1052.13	K	Joback Method
tf	367.01	K	Aqueous Solubility Prediction Method
tf	367.12	K	Aqueous Solubility Prediction Method
vc	0.741	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	611.91	J/mol×K	1052.13	Joback Method
cpg	557.11	J/mol×K	823.30	Joback Method
cpg	568.42	J/mol×K	861.44	Joback Method
cpg	578.78	J/mol×K	899.58	Joback Method
cpg	588.24	J/mol×K	937.72	Joback Method
cpg	596.87	J/mol×K	975.85	Joback Method
cpg	604.74	J/mol×K	1013.99	Joback Method

dvisc	0.0000255	Paxs	823.30	Joback Method
dvisc	0.0008679	Paxs	481.12	Joback Method
dvisc	0.0003532	Paxs	538.15	Joback Method
dvisc	0.0001707	Paxs	595.18	Joback Method
dvisc	0.0000937	Paxs	652.21	Joback Method
dvisc	0.0000567	Paxs	709.24	Joback Method
dvisc	0.0000369	Paxs	766.27	Joback Method
hfust	28.23	kJ/mol	367.40	NIST Webbook
hfust	37.30	kJ/mol	368.00	NIST Webbook
hfust	25.04	kJ/mol	369.00	NIST Webbook
hsubt	110.10 ± 0.50	kJ/mol	353.00	NIST Webbook

Sources

Aqueous Solubility Prediction Method: <http://onschallenge.wikispaces.com/file/view/AqueousDataset002.xlsx/351826032/AqueousDa>

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

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

Thermodynamic analysis of the solubility of ketoprofen in some <https://www.doi.org/10.1016/j.fluid.2010.03.031>

Solubility and Speciation of Ketoprofen <https://www.doi.org/10.1021/acs.jced.5b00812>

and Aspirin in Supercritical CO₂ by <https://www.doi.org/10.1021/je9002866>

Partial Thermodynamics of Ketoprofen in Ethanol + Water <http://onschallenge.wikispaces.com/file/view/AqueousDataset003.xlsx/351830174/AqueousDa>

Observed Solubility Prediction Method: <http://pubs.acs.org/doi/abs/10.1021/ci9903071>

Crippen Method: <https://www.doi.org/10.1021/je034163p>

The Solubility of Benzocaine, Lidocaine, and Procaine in Liquid and <http://webbook.nist.gov/cgi/cbook.cgi?ID=C22071154&Units=SI>

Supercritical Carbon Dioxide:

Legend

- cpg:** Ideal gas heat capacity
- dvisc:** Dynamic viscosity
- gf:** Standard Gibbs free energy of formation
- hf:** Enthalpy of formation at standard conditions
- hfus:** Enthalpy of fusion at standard conditions
- hfust:** Enthalpy of fusion at a given temperature
- 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
rinpola:	Non-polar retention indices
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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