

Maltose

Other names:	.alpha.-trehalose 4-(«alpha»-D-Glucopyranosido)-«alpha»-glucopyranose 4-(«alpha»-D-Glucosido)-D-glucose 4-(Â«alphaÂ»-D-Glucopyranosido)-Â«alphaÂ»-glucopyranose 4-(Â«alphaÂ»-D-Glucosido)-D-glucose 4-O-.alpha.-D-glucopyranosyl-D-glucose 4-O-«alpha»-d-Glucopyranosyl-d-glucose 4-O-Â«alphaÂ»-d-Glucopyranosyl-d-glucose Advantose 100 Cextromaltose D-(+)-Maltose D-Glucose, 4-O-«alpha»-D-glucopyranosyl- D-Glucose, 4-O-Â«alphaÂ»-D-glucopyranosyl- D-Maltose D-trehalose Finetose Finetose F Malt sugar Maltobiose Maltodiose Maltos Martos-10 Sunmalt Sunmalt S alpha,alpha-trehalose «alpha»-Malt sugar Â«alphaÂ»-Malt sugar
Inchi:	InChI=1S/C12H22O11/c13-1-3-5(15)6(16)9(19)12(22-3)23-10-4(2-14)21-11(20)8(18)7(10)
InchiKey:	GUBGYTABKSRVRQ-GXGLMDGZSA-N
Formula:	C12H22O11
SMILES:	OCC1OC(OC2C(CO)OC(O)C(O)C2O)C(O)C(O)C1O
Mol. weight [g/mol]:	342.30
CAS:	69-79-4

Physical Properties

Property code	Value	Unit	Source
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gf	-1334.42	kJ/mol	Joback Method
hf	-1959.15	kJ/mol	Joback Method
hfus	68.92	kJ/mol	Joback Method
hvap	185.55	kJ/mol	Joback Method
log10ws	0.36		Estimated Solubility Method
logp	-5.397		Crippen Method
mccvol	222.790	ml/mol	McGowan Method
pc	4311.22	kPa	Joback Method
tb	1289.46	K	Joback Method
tc	1783.82	K	Joback Method
tf	767.77	K	Joback Method
vc	0.777	m ³ /kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	889.71	J/mol×K	1289.46	Joback Method
cpg	874.12	J/mol×K	1371.85	Joback Method
cpg	848.56	J/mol×K	1454.25	Joback Method
cpg	812.80	J/mol×K	1536.64	Joback Method
cpg	766.60	J/mol×K	1619.03	Joback Method
cpg	709.75	J/mol×K	1701.43	Joback Method
cpg	641.99	J/mol×K	1783.82	Joback Method
cps	434.70	J/mol×K	300.00	NIST Webbook
cps	424.90	J/mol×K	293.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	436.10	J/mol×K	298.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	442.60	J/mol×K	303.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides

cps	452.00	J/mol×K	308.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	460.40	J/mol×K	313.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	470.90	J/mol×K	318.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	485.30	J/mol×K	323.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	394.80	J/mol×K	288.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	503.80	J/mol×K	333.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	515.00	J/mol×K	338.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	528.10	J/mol×K	343.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	542.20	J/mol×K	348.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides

Partial Molar Volumes and Isentropic Compressibilities of Some Saccharides in Aqueous Solutions of: Phosphate-Based Ionic Liquid Salts:

<https://www.doi.org/10.1021/je400977r>

Evaluation of the effect of ionic-liquids as soluting-out agents on the solubility of organic compounds in aqueous solutions: Polyhydroxy Solutes in Aqueous Solutions: Molecular Interactions of Saccharides and Their Derivatives with Thiamine: Influence of Pyridoxine HCl-Vitamin B6 on Aqueous Solutions: Isentropic Compressibility and Refractive Indices of Saccharides and Their Derivatives in Aqueous Solutions:

<https://www.doi.org/10.1021/acs.jced.5b00845>

http://pubs.acs.org/doi/suppl/10.1021/ci034243x/suppl_file/ci034243xsi20040112_053635.txt

<https://www.doi.org/10.1016/j.fluid.2017.12.006>

<https://www.doi.org/10.1021/acs.jced.5b00940>

<https://www.doi.org/10.1021/acs.jced.7b00937>

<https://www.doi.org/10.1021/acs.jced.8b00681>

Legend

cpg:	Ideal gas heat capacity
cps:	Solid phase 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
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
vc:	Critical Volume

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