

D-Galactose

Other names:	(+)-galactose D(+)-galactose D-(+)-galactose Galactose Galactose, D-
Inchi:	InChI=1S/C6H12O6/c7-1-3(9)5(11)6(12)4(10)2-8/h1,3-6,8-12H,2H2/t3-,4+,5+,6-/m0/s1
InchiKey:	GZCGUPFRVQAUUE-KCDKBNATSA-N
Formula:	C6H12O6
SMILES:	O=CC(O)C(O)C(O)C(O)CO
Mol. weight [g/mol]:	180.16
CAS:	59-23-4

Physical Properties

Property code	Value	Unit	Source
gf	-793.74	kJ/mol	Joback Method
hf	-1035.02	kJ/mol	Joback Method
hfus	19.93	kJ/mol	Joback Method
hvap	117.51	kJ/mol	Joback Method
log10ws	1.62		Crippen Method
logp	-3.379		Crippen Method
mcvol	126.320	ml/mol	McGowan Method
pc	6631.37	kPa	Joback Method
ss	205.40	J/molxK	NIST Webbook
tb	844.48	K	Joback Method
tc	1034.02	K	Joback Method
tf	443.48	K	Joback Method
vc	0.460	m3/kmol	Joback Method

Temperature Dependent Properties

Property code	Value	Unit	Temperature [K]	Source
cpg	419.64	J/molxK	1034.02	Joback Method
cpg	390.04	J/molxK	844.48	Joback Method
cpg	395.83	J/molxK	876.07	Joback Method

cpg	401.25	J/mol×K	907.66	Joback Method
cpg	406.32	J/mol×K	939.25	Joback Method
cpg	411.06	J/mol×K	970.84	Joback Method
cpg	415.49	J/mol×K	1002.43	Joback Method
cps	220.54	J/mol×K	296.90	NIST Webbook
cps	203.30	J/mol×K	288.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	210.70	J/mol×K	293.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	213.70	J/mol×K	298.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	217.50	J/mol×K	303.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	222.30	J/mol×K	308.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	227.10	J/mol×K	313.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	229.40	J/mol×K	318.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	235.60	J/mol×K	323.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides

cps	236.80	J/mol×K	328.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	242.80	J/mol×K	333.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	247.00	J/mol×K	338.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	251.20	J/mol×K	343.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	254.70	J/mol×K	348.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	257.50	J/mol×K	353.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	265.10	J/mol×K	358.15	Temperature dependence of the heat capacities in the solid state of 18 mono-, di-, and poly-saccharides
cps	216.30	J/mol×K	300.00	NIST Webbook
dvisc	4.8974616e-08	Paxs	844.48	Joback Method
dvisc	0.0017994	Paxs	443.48	Joback Method
dvisc	0.0000991	Paxs	510.31	Joback Method
dvisc	0.0000107	Paxs	577.15	Joback Method
dvisc	0.0000018	Paxs	643.98	Joback Method
dvisc	0.0000004	Paxs	710.81	Joback Method
dvisc	0.0000001	Paxs	777.65	Joback Method
hfust	43.80	kJ/mol	436.20	NIST Webbook

Sources

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Conductivity in MnSO₄-saccharide-water solutions at 298.15 K: Temperature dependence of the heat capacities in the solid state of 18-crown-6 and polyhydroxy solutes interaction parameters for the Effect of sodium acetate on the polymeric system of poly(2-vinylpyridine)-monosaccharides in aqueous solutions over temperature range 288.15 to 318.15 K: Densities and Viscosities of Polyhydroxy Solutes in Aqueous Yelluriteammonium bromide Saccharide Dimedium Temperature: Electro Interactions in Aqueous Solutions: solvent interactions of measurements and Correlation of drug Water Activity in Ionic Aqueous Solutions Containing and transport properties of some saccharides in Aqueous Drug-ionic liquid systems: Diffusion Coefficients of d-Glucose in (DMSO + H₂O) solutions at different diffusion Coefficients for Six Sugars and 0.1 MPa at Different Pressures (120.0 to 3.2) aqueous solutions: Viscosimetric studies on saccharides in aqueous magnesium chloride solutions and their interaction with 18-crown-6 Aldehyde, Caffeic Acid, d-Galactose, Mannuronic Acid and Correlation of Liquid Ethanol-Water Solutions Ternary Systems of Water + d-Fructose + 1-Butanol, Water + d-Glucose + 1-Butanol, and Water + d-Galactose + 1-Butanol at (288.2, 303.2 and 318.2) K:

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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
hfust:	Enthalpy of fusion 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
ss:	Solid phase molar entropy at standard conditions
tb:	Normal Boiling Point Temperature
tc:	Critical Temperature
tf:	Normal melting (fusion) point
vc:	Critical Volume

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