

UNIVERSITY OF SWAZILAND

BACHELOR OF SCIENCE

SUPPLEMENTARY EXAMINATION 2016

TITLE OF PAPER : PHYSICAL CHEMISTRY

COURSE NUMBER : C202

TIME : 3 HOURS

INSTRUCTIONS : THERE ARE SIX QUESTIONS

: ANSWER ANY FOUR QUESTIONS

: BEGIN THE ANSWER TO EACH QUESTION ON
A SEPARATE SHEET OF PAPER

: DATA SHEETS ARE PROVIDED WITH THIS
EXAMINATION PAPER

DO NOT OPEN THIS PAPER UNTIL THE INVIGILATOR INSTRUCTS YOU TO DO SO.

Question 1(25 marks)

- a) Define the variable, compressibility factor, z. With the aid of Lennard-Jones potential plot, compressibility and isotherm plots, compare and contrast real and ideal gases.

Your account should make mention of interactions, equations and any necessary theories to help clarify your discussion.

[15]

- b) Write short notes on any One of the following:

- i) Virial equation [10]
- ii) van der waal's equation [10]

Use diagrams, equations or plots to clarify your notes where necessary.

QUESTION 2 [25 marks]

- a) A real gas equation of state for a gas is given by:

$$P = RT(V_m - \beta)^{-1} - (\alpha/T)V_m^{-2} \quad (1)$$

- (i) Derive an expression for $V_{m,c}$, T_c and P_c . [12]
 - (ii) Find an expression for the Boyle's temperature, T_B . [4]
 - (iii) Estimate the temperature at which oxygen behaves as an ideal gas, T_B given the constants: $\alpha=1.748 \text{ L}^2 \text{ atm mol}^{-2} \text{ K}$ and $\beta= 0.0345 \text{ L mol}^{-1}$. [2]
 - (iv) Estimate the radii of real gas molecules using equation (1) for real gases given a critical molar volume of $250 \text{ cm}^3 \text{ mol}^{-1}$ [4]
- b) Using the critical point expressions for $V_{m,c}$, T_c and P_c find an expression or value for compressibility at the critical point, Z_c [3]

Question 3 [25 Marks]

- a) Write short notes on the following

- i) enthalpy change [5]
- ii) Hess's Law [5]

- b) Derive Kirrchoff's equation: [6]

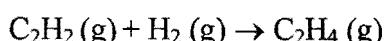
$$\Delta H_r(T_2) = \Delta H_r(T_1) + \Delta_r C_{p,m} \Delta T$$

where $C_{p,m}$ is temperature independent.

- b) Using the data in the table below calculate

- i) $\Delta_r H^\theta$ at 298 K [4]
- ii) $\Delta_r H$ at 346 K [5]

for the hydrogenation reaction:



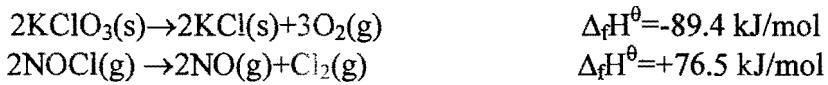
	$\text{C}_2\text{H}_4(\text{g})$	$\text{H}_2(\text{g})$	$\text{C}_2\text{H}_2(\text{g})$
$C_{p,m} \text{ J/mol/K}$	43.56	43.93	28.82
$\Delta_f H^\theta \text{ kJ/mol}$	+52.30	0	+226.8

Question 4 [25 Marks]

- a) Using examples and/or diagrams compare and contrast any one pair of the following terms
- reversible and irreversible expansion [10]
 - path and state functions [10]
- b) 4 moles of pentane occupies 25 L at 315 K.
- Derive an expression for reversible isothermal expansion. [6]
 - Calculate the work done and heat involved when the gas expands isothermally against a constant external pressure of 115 torr until its volume has doubled. [4]
 - Calculate the efficiency of the system in 1 b (ii) above. [5]

Question 5 [25 Marks]

- a) Define internal energy change [10]
- b) To Calibrate a calorimeter a 0.120 g naphthalene, $C_{10}H_8(s)$, was burnt at constant volume and it caused the temperature of the calorimeter to rise by 3.05 K. Then 0.10 g of an unknown compound was burned in the same calorimeter, causing a temperature rise of 2.05 K.
- Calculate the heat capacity of the calorimeter [3]
 - Is the unknown compound phenol, $C_6H_5OH(s)$ or ethanol, $CH_3CH_2OH(l)$ whose enthalpies of combustion are $\Delta_fH^\theta = -3054 \text{ kJ mol}^{-1}$ and $-1368 \text{ kJ mol}^{-1}$ respectively. [4]
- c) Calculate the standard enthalpies of formation of:
- $KClO_3(s)$ from the enthalpy of formation of KCl [4]
 - $NOCl(g)$ from the enthalpy of formation of NO [4]
- Given the attached table and the following information:



Useful information:

	Molecular weights/g mol ⁻¹
Benzoic acid	122.12
D-ribose $C_5H_{10}O_5$ (s)	150.13

QUESTION 6 [25 MARKS]

- a) Write short notes on any Two of the following: [10]
- Eutectic temperature and Congruent melting point
 - Zeotrope and Azeotrope
 - Lower consolute and upper consolute temperature

- b) a) Draw a sketch of the phase diagram of carbon dioxide and explain briefly the slopes and curvature of the liquid-solid and the liquid-gas boundaries, respectively. [5]
- c) i) Derive the Clausius-Clapeyron equation for evaporation in the form
- $$\frac{d(\ln p)}{dT}. \quad [5]$$
- ii) The triple point of phenol is at 6.3°C and 37 mm Hg. Predict the boiling point of phenol at 0.091 atm pressure. [5]
-

Useful Relations				General Data			
(RT) _{298.15K} =2.4789 kJ/mol				speed of light	c	2.997 925x10 ⁸ ms ⁻¹	
(RT/F) _{298.15K} =0.025 693 V				charge of proton	e	1.602 19x10 ⁻¹⁹ C	
T/K: 100.15 298.15 500.15 1000.15				Faraday constant	F=Le	9.648 46x10 ⁴ C mol ⁻¹	
T/Cm ⁻¹ : 69.61 207.22 347.62 695.13				Boltzmann constant	k	1.380 66x10 ⁻²³ J K ⁻¹	
1mmHg=133.222 N m ⁻²				Gas constant	R=Lk	8.314 41 J K ⁻¹ mol ⁻¹	
hc/k=1.438 78x10 ⁻² m K						8.205 75x10 ⁻² dm ³ atm K ⁻¹ mol ⁻¹	
1atm	1 cal	1 eV	1cm ⁻¹				
-1.01325x10 ⁵ Nm ⁻²	=4.184 J	=1.602 189x10 ⁻¹⁹ J	=0.124x10 ⁻³ eV	Planck constant	h	6.626 18x10 ⁻³⁴ Js	
-760torr		=96.485 kJ/mol	=1.9864x10 ⁻²³ J		$\hbar = \frac{h}{2\pi}$	1.054 59x10 ⁻³⁴ Js	
-1 bar		= 8065.5 cm ⁻¹					
SI-units:				Avogadro constant	L or N _A	6.022 14x10 ²³ mol ⁻¹	
1 L = 1000 ml = 1000cm ³ = 1 dm ³				Atomis mass unit	u	1.660 54x10 ⁻²⁷ kg	
1 dm = 0.1 m				Electron mass	m _e	9.109 39x10 ⁻³¹ kg	
1 cal (thermochemical) = 4.184 J				Proton mass	m _p	1.672 62x10 ⁻²⁷ kg	
dipole moment: 1 Debye = 3.335 64x10 ⁻³⁰ C m				Neutron mass	m _n	1.674 93x10 ⁻²⁷ kg	
force: 1N=1J m ⁻¹ =1kgms ⁻² =10 ⁵ dyne pressure: 1Pa=1Nm ⁻² =1Jm ⁻³				Vacuum permittivity	$\epsilon_0 = \mu_0^{-1} c^{-2}$	8.854 188x10 ⁻¹² J ⁻¹ C ² m ⁻¹	
1J = 1 Nm				Vacuum permeability	μ_0	4πx10 ⁻⁷ Js ² C ⁻² m ⁻¹	
power: 1W = 1J s ⁻¹			potential: 1V = 1 J C ⁻¹	Bohr magneton	$\mu_B = \frac{e\hbar}{2m_e}$	9.274 02x10 ⁻²⁴ JT ⁻¹	
magnetic flux: 1T=1Vs m ⁻² =1JCsm ⁻²		current: 1A=1Cs ⁻¹		Nuclear magneton	$\mu_N = \frac{e\hbar}{2m_p}$	5.05079x10 ⁻²⁷ JT ⁻¹	
Prefixes:				Gravitational constant	G	6.67259x10 ⁻¹¹ Nm ² kg ⁻²	
p n m m c d k M G				Gravitational acceleration	g	9.80665 ms ⁻²	
pico nano micro milli centi deci kilo mega giga				Bohr radius	a ₀	5.291 77x10 ⁻¹¹ m	
10 ⁻¹² 10 ⁻⁹ 10 ⁻⁶ 10 ⁻³ 10 ⁻² 10 ⁻¹ 10 ³ 10 ⁶ 10 ⁹							

THE PERIODIC TABLE OF ELEMENTS

Group	1	2	3	4	5	6	7	8	9	10	11	12	13	14	15	16	17	18
	IA	IIA	IIIB	IVB	VIB	VIB	VIIIB	VIIIB			IB	IIIB	IIIA	IVA	VA	VIA	VIIA	VIIIA
Period 1	1 H 1.008																	2 He 4.003
2	3 Li 6.94	4 Be 9.01																10 Ne 20.18
3	11 Na 22.99	12 Mg 24.31																18 Ar 39.95
4	19 K 39.10	20 Ca 40.08	21 Sc 44.96	22 Ti 47.90	23 V 50.94	24 Cr 52.01	25 Mn 54.9	26 Fe 55.85	27 Co 58.71	28 Ni 58.71	29 Cu 63.54	30 Zn 65.37	31 Ga 69.7	32 Ge 72.59	33 As 74.92	34 Se 78.96	35 Br 79.91	36 Kr 83.80
5	37 Rb 85.47	38 Sr 87.62	39 Y 88.91	40 Zr 91.22	41 Nb 91.22	42 Mo 95.94	43 Tc 98.9	44 Ru 101.1	45 Rh 102.9	46 Pd 106.4	47 Ag 107.9	48 Cd 112.4	49 In 114.8	50 Sn 118.7	51 Sb 121.8	52 Te 127.6	53 I 126.9	54 Xe 131.3
6	55 Cs 132.9	56 Ba 137.3	71 Lu 174.9	72 Hf 178.5	73 Ta 180.9	74 W 183.8	75 Re 186.2	76 Os 190.2	77 Ir 192.2	78 Pt 195.1	79 Au 196.9	80 Hg 200.6	81 Tl 204.4	82 Pb 207.2	83 Bi 208.9	84 Po 210	85 At 210	86 Rn 222
7	87 Fr 223	88 Ra 226.0	103 Lr 257	104 Unq	105 Unp	106 Unh	107 Uns	108 Uno	109 Une									

Lanthanides			57 La 138.9	58 Ce 140.1	59 Pr 140.9	60 Nd 144.2	61 Pm 146.9	62 Sm 150.9	63 Eu 151.3	64 Gd 157.3	65 Tb 158.9	66 Dy 162.5	67 Ho 164.9	68 Er 167.3	69 Tm 168.9	70 Yb 173.0
Actinides			89 Ac 227.0	90 Th 232.0	91 Pa 231.0	92 U 238.0	93 Np 237.1	94 Pu 239.1	95 Am 241.1	96 Cm 247.1	97 Bk 249.1	98 Cf 251.1	99 Es 254.1	100 Fm 257.1	101 Md 258.1	102 No 255

Numbers below the symbol indicates the atomic masses; and the numbers above the symbol indicates the atomic numbers.

SOURCE: International Union of Pure and Applied Chemistry, I mills, ed., Quantities, Units, and symbols in Physical Chemistry, Blackwell Scientific publications, Boston, 1988, pp 86-98.

Standard molar enthalpies of formation at 298.15 K

Temperature dependence of heat capacities, $C_{p,m} = a+bT+cT^{-2}$

	M_r	$\Delta H_f^\theta / \text{kJ/mol}$		M_r	$\Delta H_f^\theta / \text{kJ/mol}$		$a/\text{J K}^{-1}\text{mol}^{-1}$	$b/10^3 \text{J K}^2\text{mol}^{-1}$	$c/10^5 \text{J Kmol}^{-1}$
H ₂ O(g)	18.015	-241.8	O ₃ (g)	47.998	+142.7	Gases (298–2000K)			
H ₂ O(l)	18.015	-285.8	NO(g)	30.006	+90.2	He, Ne, Ar, Kr, Xe	20.78	0	0
H ₂ O ₂ (l)	34.015	-187.8	NO ₂ (g)	46.006	+33.2	H ₂	27.28	3.26	0.50
NH ₃ (g)	17.031	-46.1	N ₂ O ₄ (g)	92.012	+9.2	O ₂	29.96	4.18	-1.67
N ₂ H ₄ (l)	32.045	+50.6	SO ₂ (g)	64.063	-296.8	N ₂	28.58	3.77	-0.50
N ₃ H(l)	43.028	+264.1	H ₂ S(g)	34.080	-20.6	Cl ₂	37.03	0.67	-2.85
N ₃ H(g)	43.028	+294.1	SF ₆ (g)	146.054	-1209	CO ₂	44.23	8.79	-8.62
HNO ₃ (l)	63.013	-174.1	HF(g)	20.006	-271.1	H ₂ O	30.54	10.29	0
NH ₂ OH(s)	33.030	-114.2	HCl(g)	36.461	-92.3	NH ₃	29.75	25.10	-1.55
NH ₄ Cl(s)	53.492	-314.4	HCl(aq)	36.461	-167.2	CH ₄	23.64	47.86	-1.92
HgCl ₂ (s)	271.50	-224.3	HBr(g)	80.917	+36.4	C(S)	16.86	4.77	-8.54
H ₂ SO ₄ (l)	98.078	-814.0	Hl(g)	127.912	+26.5				
H ₂ SO ₄ (aq)	98.078	-909.3	CO ₂ (g)	44.010	-393.5				
NaCl(s)	58.443	-411.0	CO(g)	28.011	-110.5				
NaOH(s)	39.997	-428.7	Al ₂ O ₃ (s)	101.945	-1675.7				
KCl(s)	74.555	-435.9	SiO ₂ (s)	60.085	-910.9	Standard molar enthalpies of formation and combustion at 298.15 K.			
KBr(s)	119.011	-392.2	FeS(s)	87.91	-100.0				
KI(s)	166.006	-327.6	FeS ₂ (s)	119.975	-178.2				
DIATOMICS Eg. N ₂ , O ₂ , H ₂	0	AgCl(s)	143.323	-127.1					
						M_r	$\Delta H_f^\theta / \text{kJ/mol}$	$\Delta H_c^\theta / \text{kJ/mol}$	
						CH ₄ (g)	16.043	-74.81	
						C ₂ H ₂ (g)	26.038	+226.8	-1300
						C ₂ H ₄ (g)	28.054	+52.30	-1411
						C ₂ H ₆ (g)	30.070	-84.64	-1580
						C ₃ H ₆ cyclopropane(g)	42.081	53.35	-2091
						C ₃ H ₆ propene(g)	42.081	20.5	-2058
He	3.5	0.021	4.22	0.084		C ₄ H ₁₀ n-butane(g)	58.124	-128.11	-2877
Ar	83.81	1.188	87.29	6.506		C ₅ H ₁₂ n-pentane(g)	72.151	-146.4	-3536
H ₂	13.96	0.117	20.38	0.9183		C ₆ H ₁₂ cyclohexane(l)	84.163	-158.2	-3920
N ₂	63.15	0.719	77.35	5.586		C ₆ H ₁₄ n-hexane(l)	86.178	-198.7	-4163
O ₂	54.36	0.444	90.18	6.820		C ₆ H ₆ benzene(l)	78.115	+48.99	-3268
Cl ₂	172.12	6.406	239.05	20.410		C ₈ H ₁₈ n-octane(l)	114.233	-249.8	-5471
Br ₂	265.80	10.573	332.35	29.45		C ₁₀ H ₈ naphthalene(l)	128.175	+78.53	-5157
I ₂	386.75	15.52	458.39	41.80		CH ₃ OH(l)	32.042	-239.0	-728.1
Hg	234.29	2.292	629.73	59.296		CH ₃ CHO(g)	44.054	-166.0	-1193
Ag	1234	11.30	2436	250.63		CH ₃ CH ₂ OH(l)	46.070	-277.0	-1368
Na	370.95	2.601	1156	98.01		CH ₃ COOH(l)	60.053	-484.2	-874.5
CO ₂	217.0	8.33	194.64	25.23		CH ₃ COOC ₂ H ₅ (l)	88.107	-486.6	-2231
H ₂ O	273.15	6.008	373.15	40.656	(44.016 at 298.15 K)	C ₆ H ₅ OH(s)	94.114	-165.0	-3054
NH ₃	195.40	5.652	239.73	23.351		C ₆ H ₅ NH ₂ (l)	93.129	-31.1	-3393
H ₂ S	187.61	2.377	212.80	18.673		NH ₂ CO.NH, urea(s)	60.056	-333.0	-632.2
CH ₄	90.68	0.941	111.66	8.18		CH ₂ (NH ₂)CO ₂ H, glycine(s)	76.068	-537.2	-964.4
C ₂ H ₆	89.85	2.86	184.55	14.7		C ₆ H ₁₂ O ₆ , α -D-glucose(s)	180.159	-1274	-2802
C ₆ H ₆	278.65	10.59	353.25	30.8		C ₆ H ₂₂ O ₆ , β -D-glucose(s)	180.159	-1268	-2808
CH ₃ OH	175.25	3.159	337.22	35.27	(37.99 at 298.15K)	C ₁₂ H ₂₂ O ₁₁ , sucrose(s)	342.303	-2222	-5645
						CH ₃ CH(OH)COOH	90.079	-694.0	-1344
						lactic acid(s)			

↓ Sublimation: ^avarious pressures: ^bat 1atm

Source: American Institute of Physics handbook, McGraw-Hill.

Heat capacities at 25°C

	$C_{v,m}$ JK ⁻¹ mol ⁻¹	$C_{p,m}$ JK ⁻¹ mol ⁻¹
He, Ne, Ar, Kr, Xe	12.47	20.78
H ₂	20.50	28.81
O ₂	21.01	29.33
N ₂	20.83	29.14
CO ₂	28.83	37.14
NH ₃	27.17	35.48
CH ₄	27.43	35.74
N ₂ O ₄		77.28
NO ₂		37.20

F.P Depression, B.P. Elevation

Solvent	F.P °C	K _f °C kg mol ⁻¹	B.P (°C, 101kNm ⁻²)	K _b °C kg mol ⁻¹
Water	0	1.86	100.0	0.52
Benzene	5.51	5.10	80.1	2.60
Acetic Acid	16.6	3.90	118.1	3.10
Cyclohexane	6.5	20.2	81.4	2.79
Camphor	177.7	40.0	205	-
Nitrobenzene	5.7	6.9	210.9	5.24
Ethanol	-177		78.5	1.22
Chloroform	-64		61.3	3.63

Third Law entropies at 25°C, Sm^θ/J K⁻¹ mol⁻¹

Solids	Liquids	Gases
Ag	42.68	Hg
C(gr)	5.77	Br ₂
C(d)	2.44	
Cu	33.4	O ₂
Zn	41.6	Cl ₂
I ₂	116.7	
S(Rh)	31.9	H ₂ O
		70.0
		CO ₂
AgCl	96.2	HNO ₃
AgBr	104.6	CH ₃ OH
CuSO ₄ ·5H ₂ O	305.4	C ₆ H ₆
HgCl ₂	144	CH ₃ COOH
Sucrose	360.2	C ₆ H ₁₂
		298.2

Standard molar Gibbs free energy and molar entropy of formation at 298.15 K

	M _r	ΔG _f ^θ /KJ/mol	S ^θ /J K ⁻¹ mol ⁻¹		M _r	ΔG _f ^θ /KJ/mol	S ^θ /J K ⁻¹ mol ⁻¹
H ₂ O(g)	18.015	-228.57	188.83	O ₃ (g)	47.998	163.2	238.93
H ₂ O(l)	18.015	-120.35	109.6	NO(g)	30.006	86.55	210.76
H ₂ O ₂ (l)	34.015	-120.35	109.6	NO ₂ (g)	46.006	51.31	240.06
NH ₃ (g)	17.031	-16.45	192.45	N ₂ O ₄ (g)	92.012	97.89	304.29
N ₂ H ₄ (l)	32.045	149.43	121.21	SO ₂ (g)	64.063	-300.19	248.22
N ₃ H(l)	43.028	327.3	140.6	H ₂ S(g)	34.080	-33.56	205.79
N ₃ H(g)	43.028	328.1	238.97	SF ₆ (g)	146.054	-1105.3	291.82
HNO ₃ (l)	63.013	-80.71	155.60	HF(g)	20.006	-273.2	173.78
NH ₂ OH(s)	33.030			HCl(g)	36.461	-95.30	186.91
NH ₄ Cl(s)	53.492	-202.87	94.6	HCl(aq)	36.461	-131.23	56.5
HgCl ₂ (s)	271.50	-178.6	146.0	HBr(g)	80.917	-53.45	198.70
H ₂ SO ₄ (l))	98.078	-690.00	156.90	HI(g)	127.912	1.70	206.59
H ₂ SO ₄ (aq)	98.078	-744.53	20.1	CO ₂ (g)	44.010	-394.36	213.74
NaCl(s)	58.443	-384.14	72.13	CO(g)	28.011	-137.17	197.67
NaOH(s)	39.997	-379.49	64.46	Al ₂ O ₃ (□,s)	101.945	-1582.3	50.92
KCl(s)	74.555	-409.14	82.59	SiO ₂	60.09	-856.64	41.84
KBr(s)	119.011	-380.66	95.90	FeS(s)	87.91	-100.4	60.29
KI(s)	166.006	-324.89	106.32	FeS ₂ (s)	119.975	-166.9	52.93
				AgCl(s)	143.323	-109.79	96.2
He(g)	4.003	0	126.15	Hg(g)	200.59	31.82	174.96
Ar(g)	39.95	0	154.84	Hg(l)	200.59	0	76.02
H ₂ (g)	2.016	0	130.684	Ag(g)	107.87	245.65	173.00
N ₂ (g)	28.013	0	191.61	Ag(s)	107.87	0	42.55
O ₂ (g)	31.999	0	205.138	Na(g)	370.95	76.76	153.71
O ₃ (g)	47.998	163.2	238.93	Na(s)	22.99	0	51.21
Cl ₂ (g)	70.91	0	223.07				
Br ₂ (g)	159.82	3.110	245.46				
Br ₂ (l)	159.82	0	152.23				
I ₂ (g)	253.81	19.33	260.69				
I ₂ (s)	253.81	0	116.135				

	M _r	ΔG _f ^θ /KJ/mol	S ^θ /J K ⁻¹ mol ⁻¹
organic compounds			
CH ₄ (g) methane	16.043	-50.72	186.26
C ₂ H ₂ (g) ethyne	26.038	209.20	200.94
C ₂ H ₄ (g) ethene	28.05	68.15	219.56
C ₂ H ₆ (g) ethane	30.070	-32.82	229.60
C ₃ H ₆ cyclopropane(g)	42.081	104.45	237.55
C ₃ H ₆ propene(g)	42.081	62.78	267.05
C ₄ H ₁₀ n-butane (g)	58.124	-17.03	310.23
C ₅ H ₁₂ n-pentane(g)	72.151	-8.20	348.40
C ₆ H ₁₂ cyclohexane (l)	84.163	26.8	
C ₆ H ₁₄ n-hexane (l)	86.178		204.3
C ₆ H ₆ benzene (l)	78.115	124.3	173.3
C ₆ H ₆ benzene (g)	78.115	129.72	269.31
C ₈ H ₁₈ n-octane (l)	114.233	6.4	361.1
C ₁₀ H ₈ naphthalene (l)	128.175		
CH ₃ OH (g)	32.042	-161.96	239.81
CH ₃ OH (l)	32.042	-166.27	126.8
CH ₃ CHO (g)	44.054	-128.86	250.3
CH ₃ CH ₂ OH (l)	46.07	-174.78	160.7
CH ₃ COOH (l)	60.053	-389.9	159.8
CH ₃ COOC ₂ H ₅ (l)	88.107	-332.7	259.4
C ₆ H ₅ OH (s)	94.114	-50.9	146.0
C ₆ H ₅ NH ₂ (l)	93.129		
CH ₂ (NH ₂)CO ₂ H, glycine (s)	75.068	-373.4	103.5
C ₆ H ₁₂ O ₆ , α-D-glucose (s)	180.159		
C ₆ H ₂₂ O ₆ , β-D-glucose (s)	180.159	-910	212
C ₁₂ H ₂₂ O ₁₁ , sucrose (s)	342.303	-1543	360.2
CH ₃ CH(OH)COOH	90.079		
lactic acid (s)			

Source: American Institute of Physics handbook, McGraw-Hill.