# CHEMISTRY 9701 THEORY QUESTIONS

# **AS: CHEMICAL ENERGETICS**

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## 3 Bond energies

#### 3(a) Bond energies in diatomic molecules (these are exact values)

#### Homonuclear

Bond	Energy/kJ mol <sup>-1</sup>
H-H	436
D-D	442
N≡N	944
0=0	496
P≡P	485
S=S	425
F-F	158
C1-C1	242
Br–Br	193
I–I	151

#### Heteronuclear

Bond	Energy/kJmol <sup>-1</sup>
H–F	562
H-C1	431
H–Br	366
H–I	299
C≡O	1077

#### 3(b) Bond energies in polyatomic molecules (these are average values)

#### Homonuclear

Bond	Energy/kJ mol <sup>-1</sup>
C-C	350
C=C	610
C≡C	840
C:::C (benzene)	520
N-N	160
N=N	410
0-0	150
Si–Si	222
P_P	200
S-S	264

#### Heteronuclear

Bond	Energy/kJmol⁻¹
C-H	410
C-C1	340
C–Br	280
C-I	240
C-N	305
C=N	610
C≡N	890
C-O	360
C=O	740
C=O in CO <sub>2</sub>	805
N-H	390
N-C1	310
O-H	460
Si-C1	359
Si-H	320
Si–O (in SiO₂(s))	460
Si=O (in SiO₂(g))	640
Р-н	320
P-C1	330
P-O	340
P=O	540
S-H	347
S-C1	250
S-O	360
S=O	500

## **CHEMICAL ENERGETICS (2002-2014)**

(b)	Car	bon disulfide is readily combusted to give CO <sub>2</sub> and SO <sub>2</sub> .	
	(i)	Construct a balanced equation for the complete combustion of ${\rm CS}_2$ .	
	(ii)	Define the term standard enthalpy change of combustion, ∆H <sup>o</sup> c.	
			[3]
(c)		culate the standard enthalpy change of formation of ${\rm CS}_2$ from the following dude a sign in your answer.	ata.
	star	ndard enthalpy change of combustion of CS <sub>2</sub> = -1110 kJ mol <sup>-1</sup>	
	star	ndard enthalpy change of formation of $CO_2 = -395 \mathrm{kJ}\mathrm{mol}^{-1}$	
	star	ndard enthalpy change of formation of SO <sub>2</sub> = -298 kJ mol <sup>-1</sup>	
ıp23			[3]

(c) The standard enthalpy changes of formation of NH<sub>3</sub>(g) and H<sub>2</sub>O(g) are as follows.

$$NH_3(g), \Delta H_f^{\bullet} = -46.0 \text{ kJ mol}^{-1}$$

$$H_2O(g), \Delta H_f^{\bullet} = -242 \text{ kJ mol}^{-1}$$

Use these data and the value of  $\Delta H^{e}_{reaction}$  given below to calculate the standard enthalpy change of formation of NO(g). Include a sign in your answer.

$$4NH_3(g) + 5O_2(g) \rightleftharpoons 4NO(g) + 6H_2O(g)$$

$$\Delta H^{+} = -906 \text{ kJ mol}^{-1}$$

[4]

s/13/qp21

3			ne chemical ncarbonate, Kl							
			ases, the use enthalpy chan				nthalpy cha	ange of rea	ction to be	calculated
	(a)	Stat	e Hess' Law.							
										[2]
			to determine ncarbonate, tw						osition of	potassium
	expe	erim	ent 1					4		
	temp Whe	perat en 0.	of 2.00 moldm ture recorded a 0200 mol of point ith a thermome	as 21.0°C. otassium ca	arbonate	, K <sub>2</sub> CO	3, was add	ded to the	acid and t	
	(b)	(i)	Construct a b	alanced eq	uation fo	or this re				
		(ii)	Calculate the Use relevant same specific	data from	the <i>Data</i>	Bookle	n experim	nent 1, stat	ing your ur	nits.
	(	iii)	Use your ans Give your ans	wer to (ii) to swer in kJ n	o calcula nol <sup>-1</sup> and	te the e	nthalpy ch a sign in	ange per n your answe	nole of K <sub>2</sub> 0 er.	CO <sub>3</sub> .
	\ \		9/							
	(	iv)	Explain why t	he hydroch	loric acid	d must b	e in an ex	cess.		
									•••••	
										[4]

#### experiment 2

The experiment was repeated with 0.0200 mol of potassium hydrogencarbonate,  $KHCO_3$ . All other conditions were the same.

In the second experiment, the temperature fell from 21.0 °C to 17.3 °C.

(c)	(i)	Construct a balanced equation for this reaction.

- (ii) Calculate the quantity of heat absorbed in experiment 2.
- (iii) Use your answer to (ii) to calculate the enthalpy change per mole of KHCO<sub>3</sub>. Give your answer in kJ mol<sup>-1</sup> and include a sign in your answer.

[3]

(d) When KHCO<sub>3</sub> is heated, it decomposes into K<sub>2</sub>CO<sub>3</sub>, CO<sub>2</sub> and H<sub>2</sub>O.

$$2KHCO_3 \rightarrow K_2CO_3 + CO_2 + H_2O$$

Use Hess' Law and your answers to (b)(iii) and (c)(iii) to calculate the enthalpy change for this reaction.

Give your answer in kJ mol<sup>-1</sup> and include a sign in your answer.

[2]

[Total: 11]

w/11/qp21

The unsaturated hydrocarbon, E, is obtained by cracking hexane and is important in the

chemical industry.
The standard enthalpy change of combustion of ${\bf E}$ is $-2059{\rm kJmol^{-1}}$ .
(d) Define the term standard enthalpy change of combustion.
[2]
When 0.47g of <b>E</b> was completely burnt in air, the heat produced raised the temperature of 200g of water by 27.5 °C. Assume no heat losses occurred during this experiment.
(e) (i) Use relevant data from the Data Booklet to calculate the amount of heat released in this experiment.
(ii) Use the data above and your answer to (i) to calculate the relative molecular mass, $M_{\rm r}$ , of ${\bf E}$ .
(f) Deduce the molecular formula of E.
(f) Deduce the molecular formula of 2.  [1]  [2]

3 Alkanes such as methane, CH<sub>4</sub>, undergo few chemical reactions. Methane will, however, react with chlorine but not with iodine.

Relevant standard enthalpy changes of formation for the reaction of methane with chlorine to form chloromethane, CH<sub>3</sub>CI, are given below.

	∆H <sup>o</sup> /kJ mol <sup>-1</sup>
CH <sub>4</sub>	-75
CH <sub>3</sub> C1	-82
HC1	-92

(a) (i) Use the data to calculate  $\Delta H_{\text{reaction}}^{\Phi}$  for the formation of CH<sub>3</sub>C1.

$$\mathrm{CH_4}$$
 +  $\mathrm{C}l_2$   $\rightarrow$   $\mathrm{CH_3C}l$  +  $\mathrm{HC}l$ 

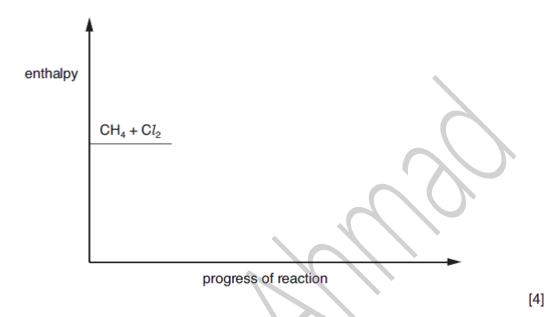
(ii) The corresponding reaction with iodine does not take place.

Use bond energy data from the <code>Data Booklet</code> to calculate a 'theoretical value' for  $\Delta H_{\rm reaction}$  for the following equation.

$$CH_4 + I_2 \rightarrow CH_3I + HI$$

(iii) Suggest why this reaction does not in fact occur.

(c) The energy of activation for the formation of CH<sub>3</sub>Cl is 16 kJ mol<sup>-1</sup>. Use this figure and your answer to (a)(i) to complete the reaction pathway diagram below showing the formation of CH<sub>3</sub>Cl from CH<sub>4</sub> and Cl<sub>2</sub>. Show clearly the intermediate organic species and the final products. Indicate on your sketch the relevant enthalpy changes and their values.



w/09/qp21

2 Ketene, C<sub>2</sub>H<sub>2</sub>O, is a member of a class of unsaturated organic compounds that is widely used in pharmaceutical research for the synthesis of organic compounds.

(c) (i)	Define the term standard enthalpy change of formation.

(ii) Use the data below to calculate the standard enthalpy change of formation of ketene.

	ΔH <sup>o</sup> /kJ mol <sup>−1</sup>
standard enthalpy change of formation of CO <sub>2</sub>	-395
standard enthalpy change of combustion of H <sub>2</sub>	-286
standard enthalpy change of combustion of CH <sub>2</sub> =C=O	-1028

[6]

w/08/qp2

(e) Carbon, hydrogen and ethene each burn exothermically in an excess of air.

$$C(s) + O_2(g) \rightarrow CO_2(g)$$

$$\Delta H_{c}^{\Theta} = -393.7 \,\mathrm{kJ} \,\mathrm{mol}^{-1}$$

$$H_2(g) \ + \ {}^{1\!\!/}_2O_2(g) \ \to \ H_2O(I)$$

$$\Delta H_{c}^{\Theta} = -285.9 \,\mathrm{kJ} \,\mathrm{mol}^{-1}$$

$${\rm C_2H_4(g)} \ + \ 3{\rm O_2(g)} \ \rightarrow \ 2{\rm CO_2(g)} \ + \ 2{\rm H_2O(I)} \qquad \Delta H_{\rm \ c}^{\rm \ e} \ = \ -1411.0 \, {\rm kJ \, mol^{-1}}$$

$$\Delta H_{c}^{\circ} = -1411.0 \,\text{kJ mol}^{-1}$$

Use the data to calculate the standard enthalpy change of formation,  $\Delta H_{\rm f}^{\rm e}$ , in kJ mol<sup>-1</sup>, of ethene at 298 K.

$$2C(s) \ + \ 2H_2(g) \ \rightarrow \ C_2H_4(g)$$

[3]

w/07/qp2

The unsaturated hydrocarbon Z is obtained by cracking hexane and is important in the

The sta	andard enthalpy change of combustion of <b>Z</b> is -2059 kJ mol <sup>-1</sup> .
(d) De	efine the term standard enthalpy change of combustion.
	[
	0.47 g of <b>Z</b> were completely burnt in air, the heat produced raised the temperature of water by 27.5 °C.
(e) (i)	Calculate the amount of heat released in this experiment.
(ii)	Use the data above and your answer to (i) to calculate the relative molecular man
	OI Z.
(f) De	educe the molecular formula of <b>Z</b> .
6/qp2	

chemical industry.

(c)	Explain the term standard enthalpy change of formation, $\Delta H_{\rm f}^{ \oplus}$ .	
		[3]
(d)	Calculate the standard enthalpy change of formation of ${\rm CS}_2$ from the following data.	
	standard enthalpy change of formation of SO <sub>2</sub> = -298 kJ mol <sup>-1</sup>	
	standard enthalpy change of formation of CO <sub>2</sub> = -395 kJ mol <sup>-1</sup>	
	standard enthalpy change of combustion of CS <sub>2</sub> = -1110 kJ mol <sup>-1</sup>	
		[3]
v/05/qp2		
(d)	(i) Explain how enthalpy changes, $\Delta H$ values, for covalent bonded molecules can calculated from bond energies.	be
(ii)	Use bond energies from the $\textit{Data Booklet}$ to calculate $\Delta \textit{H}$ for the following dissociation.	ng
X	$2HI(g) \rightarrow H_2(g) + I_2(g)$	



3	(a)	(i)	What is meant by the standard enthalpy change of formation, $\Delta H^{\circ}_{\ \ t}$ , of a compound? Explain what is meant by the term standard.
		(ii)	Write an equation, with state symbols, for the $\Delta H^{\circ}_{\ \ f}$ of water.
		(iii)	Explain why the $\Delta H^{\circ}_{\dagger}$ for water is identical to the standard enthalpy change of combustion of hydrogen.
			[4]
	(b)		en calcium is placed in water, aqueous calcium hydroxide is formed and hydrogen is en off.
		(i)	Write the equation for the reaction of calcium with water.
		(ii)	When 1.00 g of calcium is placed in 200 g of water, the temperature increases by 12.2 °C when the reaction is completed. The specific heat capacity of water, $c$ , is 4.2 J g <sup>-1</sup> K <sup>-1</sup> .
			Calculate the heat released in the experiment.
<			

(iii) Calculate the standard enthalpy change of reaction in kJ mol-1 for your equation in

[4]
(c) (i) State Hess' Law.
(ii) Use Hess' Law and your result in (b)(iii) to calculate the $\Delta H_f^{\circ}$ of Ca(OH) <sub>2</sub> (aq). You also need the $\Delta H_f^{\circ}$ of water which is $-286\mathrm{kJmol^{-1}}$ .
[4]
(d) Calculate the volume of hydrogen, measured at room temperature and pressure, liberated in the experiment described in (b)(ii).
[2]
[Total : 14]
v/03/qp2

(b)(i).

(d)	Most of the ammonia produced which is not used as fertiliser, is oxidised to nitric acid, $\mbox{HNO}_3$ .
	Construct an equation for the oxidation of ammonia by atmospheric oxygen to form nitric acid.
	[1]

(e) Urea, CO(NH<sub>2</sub>)<sub>2</sub>, is a naturally occurring substance which can be hydrolysed with water to form ammonia according to the following equation.

$$H_2O(I) + CO(NH_2)_2(aq) \rightarrow CO_2(aq) + 2NH_3(aq)$$

The standard enthalpy changes of formation of water, urea, carbon dioxide and ammonia (in aqueous solution) are given below.

compound	ΔH <sub>f</sub> <sup>⊕</sup> /kJ mol <sup>-1</sup>
H <sub>2</sub> O(I)	-287.0
CO(NH <sub>2</sub> ) <sub>2</sub> (aq)	-320.5
CO <sub>2</sub> (aq)	-414.5
NH <sub>3</sub> (aq)	-81.0

Use these data to calculate the standard enthalpy change for the hydrolysis of urea.

[2]

w/02/qp2

3	With the prospect that fossil fuels will become increasingly scarce in the future, many
	compounds are being considered for use in internal combustion engines. One of these is
	DME or dimethyl ether, CH <sub>3</sub> OCH <sub>3</sub> . DME is a gas which can be synthesised from methanol.
	Methanol can be obtained from biomass, such as plant waste from agriculture.

change of compustion, $\Delta H_c^{\bullet}$ , for DIME at 298 K.	
equation	
definition	

(a) Define, with the aid of an equation which includes state symbols, the standard enthalpy

.....[3]

(b) DME may be synthesised from methanol. Relevant enthalpy changes of formation,  $\Delta H_{\rm f}^{\rm e}$ , for this reaction are given in the table below.

compound	ΔH <sup>e</sup> <sub>f</sub> /kJ mol <sup>-1</sup>
CH₃OH(I)	-239
CH <sub>3</sub> OCH <sub>3</sub> (g)	-184
H <sub>2</sub> O(I)	-286

Use these values to calculate  $\Delta H_{\rm reaction}^{\bullet}$  for the synthesis of DME, using the following equation. Include a sign in your answer.

$$2CH_3OH(I) \rightarrow CH_3OCH_3(g) + H_2O(I)$$

$$\Delta H_{\text{reaction}}^{\Theta} = \dots \text{kJ mol}^{-1}$$
[3]

s/12/qp23

2			anol, CH <sub>3</sub> OH, are co be used in car engir	_	sible replacements for fossil
	(a) Define, with the aid of an equation which includes state symbols, the standar change of combustion, ΔH <sup>o</sup> <sub>c</sub> , for methanol at 298 K.				
		equation			
		definition			
					121
		•••••			[3]
					ogen. Relevant ΔH <sup>e</sup> c values
	TOT	this reaction are give	en in the table below		
			compound	ΔH <sup>o</sup> <sub>c</sub> /kJ mol <sup>-1</sup>	
			CO(g)	-283	
			$H_2(g)$	-286	
			CH <sub>3</sub> OH(g)	-726	
	(b)		calculate ΔH <sup>e</sup> <sub>reaction</sub> sign in your answer		methanol, using the following
			CO(a) + 2H (a	ı) → CH-OH(a)	
			CO(g) + 211 <sub>2</sub> (g	) → CH₃OH(g)	
					∆H <sup>o</sup> reaction =kJ mol <sup>-1</sup>
					reaction

s/12/qp22

[3]

	Methanol, $\mathrm{CH_3OH}$ , is considered to be a possible alternative to fossil fuels, particularly for use in vehicles.					
	Nethanol can be produced from fossil fuels and from agricultural waste. It can also be ynthesised from carbon dioxide and hydrogen.					
(a)	Define, with the aid of an equation which includes state symbols, the standard enthalpy change of formation of carbon dioxide.					
	equ	ation				
	def	inition				
					[3]	
(b)	Rel	evant ∆H° va	lues for the reaction	that synthesises meth	nanol are given in the table.	
			compound	$\Delta H_7^{\circ}/\text{kJ} \text{mol}^{-1}$		
			CO <sub>2</sub> (g)	-394		
			CH <sub>3</sub> OH(g)	-201	1	
			H₂O(g)	-242		
	(i)	Use these va	alues to calculate $\Delta t$	7 this synthes	is of methanol.	
		Include a sig	n in your answer.			
			CO <sub>-</sub> (a) + 3H <sub>-</sub> (a) s		(a)	
			2(3)		9/	
		1				
				ΔH <sup>e</sup> <sub>reac</sub>	<sub>otion</sub> =kJ mol <sup>-1</sup>	
	(ii)	Suggest on answer.	e possible environ	mental advantage of	this reaction. Explain your	
	[5]					

3

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3	1	ı u	$\nu$	_

2 Halogenoalkanes have been widely used as aerosol propellants, refrigerants and solvents for many years.

Fluoroethane, CH<sub>3</sub>CH<sub>2</sub>F, has been used as a refrigerant. It may be made by reacting ethene with hydrogen fluoride.

You are to calculate a value for the C-F bond energy in fluoroethane.

(a) Use relevant bond energies from the Data Booklet, and the equation below to calculate a value for the bond energy of the C-F bond.

$$\mathrm{CH_2} = \mathrm{CH_2}(\mathrm{g}) \qquad + \qquad \mathrm{HF}(\mathrm{g}) \qquad \longrightarrow \qquad \mathrm{CH_3}\mathrm{CH_2F}(\mathrm{g}) \qquad \Delta H^{\oplus} = -\ 73\ \mathrm{kJ}\ \mathrm{mol}^{-1}$$

(b) Another halogenoalkane which was used as a refrigerant, and also as an aerosol propellant, is dichlorodifluoromethane, CC1<sub>2</sub>F<sub>2</sub>.

State two reasons why compounds such as  $\mathrm{CH_3CH_2F}$  and  $\mathrm{CC}\mathit{l_2F_2}$  have been used as aerosol propellants and refrigerants.



s/11/qp22

1	Hydrazine, $N_2H_4$ , can be used as a rocket fuel and is stored as a liquid. It reacts exothermically with oxygen to give only gaseous products.					
	The enthalpy change of a reaction such as that between hydrazine and oxygen may be calculated by using standard enthalpy changes of formation.					
	(a)	Def	ine the term <i>stand</i>	dard enthalpy cha	nge of formation,	ΔH <sub>f</sub> <sup>⇔</sup> .
						[3]
	(b)	Hyd	Irazine reacts with	oxygen accordin	g to the following	equation.
			N	$_{2}H_{4}(I) + O_{2}(g) -$	$\rightarrow$ N <sub>2</sub> (g) + 2H <sub>2</sub> O	(g)
		(i)	Use the data in reaction.	the table to ca	alculate the stan	dard enthalpy change of this
				compound	ΔH <sub>f</sub> <sup>+</sup> /kJ mol <sup>-1</sup>	
				N <sub>2</sub> H <sub>4</sub> (I)	50.6	
				H <sub>2</sub> O(g)	-241.8	
				$\setminus \lambda$	Ý	
					ΔH <sup>+</sup> =	kJ mol <sup>-1</sup>
		(ii)	Although the al	nove reaction is	highly exothern	nic, hydrazine does not burn
		(11)	spontaneously in	oxygen.	riigiliy exotileili	no, nyarazine does not bum
			Suggest a reaso	n for this.		
s/10/	qp23	3				

	(ii)	Use the <code>Data Booklet</code> to calculate the enthalpy change that occurs when one mole of gaseous magnesium ions, ${\rm Mg^{2+}}$ , is formed from one mole of gaseous magnesium atoms.
		Include a sign in your answer.
		enthalpy change = kJ mol <sup>-1</sup>
		[3]
/09/qp2		
(d)	Use	equation for the complete combustion of ethyne is given below.  appropriate bond energy data from the Data Booklet to calculate a value for the halpy change of combustion of ethyne.
		$C_2H_2(g) + \frac{5}{2}O_2(g) \rightarrow 2CO_2(g) + H_2O(g)$
		[3]
(e)	The	value for the standard enthalpy change of combustion of ethyne is -1300 kJ mol <sup>-1</sup> .
	(i)	Define the term standard enthalpy change of combustion.
	(ii)	Explain why your answer to (d) does not have the same value as the standard enthalpy change of combustion.
		[3]
/06/qp2		

2 Ethanol, C<sub>2</sub>H<sub>5</sub>OH, is a most important industrial chemical and is used as a solvent, a fuel and an intermediate in large scale organic synthesis.

Ethanol is prepared industrially by the reaction of ethene and steam in the presence of a catalyst.

$$C_2H_4(g) + H_2O(g) \rightarrow C_2H_5OH(g)$$

The standard enthalpy change of the reaction can be determined by using the standard enthalpy changes of combustion,  $\Delta H_0^{\circ}$ , at 298 K.

$$\Delta H_{c}^{\circ}/\text{kJ mol}^{-1}$$
 $C_{2}H_{4}(g)$  -1411
 $C_{2}H_{5}OH(I)$  -1367

(a) Calculate the standard enthalpy change for the following reaction.

$$C_2H_4(g) + H_2O(I) \rightarrow C_2H_5OH(I)$$

		[2	]
(b)	(i)	Define the term standard enthalpy change of combustion.	
	(II)	Explain why the state symbols for water and ethanol given in the equation in (a have been changed from those quoted in the industrial process.	)
X			٠
	(iii)	Write the equation for the complete combustion of ethanol.	
		[4	j

s/02/qp2

