

# Combustion Calorimetry and Thermodynamic Functions of Cyanocobalamin

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## Abstract

In a calorimeter with a static bomb and an isothermal shield, the energy of combustion of the cyanocobalamin has been measured at 298.15 K. Physico-chemical methods established the products of combustion of cyanocobalamin in the conditions of calorimetric experiment. The enthalpy of combustion  $\hat{I}^{\circ}cH^{\circ}$  and the thermodynamic parameters  $\hat{I}^{\circ}fH^{\circ}$ ,  $\hat{I}^{\circ}fG^{\circ}$  of the cyanocobalamin at  $T = 298.15$  K and  $p = 0.1$  MPa have been calculated. Thermodynamic parameters  $\hat{I}^{\circ}fH^{\circ}$ ,  $\hat{I}^{\circ}fS^{\circ}$  were determined and used to calculate the enthalpy of formation of cyanocobalamin.

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*Index terms*— vitamin B12; cyanocobalamin; combustion calorimetry; thermodynamic functions.

## 1 Introduction

Cyanocobalamin, also called vitamin B 12 (PubChem CID: 5479203), is a water-soluble vitamin with a key role in the normal functioning of the brain and nervous system, and for the formation of blood. Vitamin B 12 is a cobalt-containing compound synthesized by bacteria and an essential nutrient in mammals, which take it up from diet [1]. The significance of vitamin B 12 adequate nutritional status throughout life span is established and the adverse effects of vitamin B 12 deficiency in human health are currently recognized [2][3][4]. In addition to the well-described reversible hematological and often irreversible neurological changes of severe vitamin B 12 deficiency, epidemiological studies revealed a more common condition, the low vitamin B 12 status particularly in elder and pregnant women [5][6]. Because vitamin B 12 is essential for DNA synthesis and cellular energy production, a low vitamin B 12 status may be a risk factor for altered cellular metabolism and age-related diseases including cognitive decline and cardio-vascular disease [7].

This work is a continuation of systematic studies of vitamins B. Earlier in the articles [8][9][10], we have investigated the thermodynamic properties of vitamins B 2, B 3 and the temperature dependence of the heat capacity of cyanocobalamin. The goals of this work include calorimetric determination of the standard thermodynamic functions of the cyanocobalamin. II.

## 2 Experimental i. Sample

Cyanocobalamin was purchased from Fluka. For phase identification, an X-ray diffraction pattern of the vitamin B 12 sample was recorded on a Shimadzu X-ray diffractometer XRD-6000 (CuK  $\gamma$  radiation, geometry  $\theta$ - $2\theta$ ) in the  $2\theta$  range from  $5^{\circ}$  to  $60^{\circ}$  with scan increment of  $0.02^{\circ}$ . The X-ray data and estimated impurity content (0.1 wt %) in the substance led us to conclude that the cyanocobalamin sample studied was an individual crystalline compound. Cyanocobalamin can crystallize in three modifications [4]: 1) "as-purchased"; 2) "wet"; 3) "dry". According to X-ray diffraction and solid-state NMR spectroscopy, we investigated the "as-purchased" sample cyanocobalamin. This sample is usually obtained by rapid crystallization from water at 343 K.

## 3 ii. Apparatus and measurement procedure

The energy of combustion,  $\hat{I}^{\circ}cU$ , of cyanocobalamin was measured in a calorimeter (V-08) with a static bomb and an isothermal shield. The calorimeter design, the procedure of measuring the energies of combustion and the

## 5 CONCLUSIONS

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43 results of calibration and testing are given elsewhere [11]. It should be noted that while checking the calorimeter  
44 by burning succinic acid, prepared at D.I. Mendeleev Research Institute of Metrology (the value of the standard  
45 enthalpy of combustion of the acid coincided with the certificate value within  $\pm 0.017\%$ ). For complete combustion  
46 of cyanocobalamin we used paraffin as an auxiliary substance.

47 Physico-chemical methods established the products of combustion of cyanocobalamin in the conditions of  
48 calorimetric experiment. Firstly, the solid products of combustion were identified by X-ray diffraction (Shimadzu  
49 X-ray diffractometer XRD-6000). Secondly, the formed liquid droplets were analyzed for phosphorus content using  
50 atomic absorption spectrophotometry (Shimadzu atomic absorption spectrophotometer AA-6300). Thirdly, the  
51 liquid droplets were titrated for total inorganic acids (Mettler Toledo pH meter Five Easy FE-20). Fourthly, the  
52 analysis of the gas phase was carried out by gas chromatography (Shimadzu GC-2010 Plus).

53 iii.

## 54 4 Results and Discussion

55 The experimental data on burning of cyanocobalamin are presented in Table 1. As a result, the energies and  
56 enthalpies of combustion of riboflavin at  $T = 298.15\text{K}$  and standard pressure were determined. The values are  
57 for the reaction:  $\text{C}_{63}\text{H}_{88}\text{N}_{14}\text{O}_7(\text{cr}) + 79.75\text{O}_2(\text{g}) + 63\text{O}_2(\text{g}) + 42.875\text{H}_2\text{O}(\text{l}) + 0.75\text{CoO}(\text{cr}) +$   
58  $+ 0.125\text{Co}_2\text{P}_2\text{O}_7(\text{cr}) + 0.75\text{H}_3\text{PO}_4(\text{sl-n}, 700^\circ\text{C}) + 7\text{N}_2(\text{g})$

59 In brackets are given the physical states of reagents: (cr), crystalline; (g), gaseous; (l), liquid; (sl-n), solution.  
60 It should be noted that we have used a significant amount of physico-chemical methods (see section 2.2) in the  
61 study of combustion products of cyanocobalamin which is an organometallic compound. The data on the enthalpy  
62 of combustion of the crystalline cyanocobalamin was used to estimate enthalpy of combustion and formation at  
63  $T = 298.15\text{K}$  and  $p = 0.1\text{MPa}$  (Table 2). Due to the fact that the standard enthalpy of formation of dicobalt  
64 diphosphate absent in the literature, we calculated the standard enthalpy and entropy of formation of  $\text{Co}_2\text{P}_2\text{O}_7$   
65 at  $298.15\text{K}$  (Table 2). In works [12,13], the absolute entropy and the standard Gibbs function of formation  
66 of dicobalt diphosphate were determined.

67 The Gibbs function of formation  $\hat{I}^\circ$  f  $G^\circ$  of the cyanocobalamin was evaluated from the  $\hat{I}^\circ$  f  $H^\circ$  and  $\hat{I}^\circ$  f  $S^\circ$   
68 [10] values (Table 2). The values conform to the following process:  $63\text{C}(\text{gr}) + 44\text{H}_2(\text{g}) + 7\text{N}_2(\text{g}) + 7\text{O}_2(\text{g}) +$   
69  $\text{Co}(\text{cr}) + \text{P}(\text{cr}) \rightarrow \text{C}_{63}\text{H}_{88}\text{N}_{14}\text{O}_7(\text{cr})$

70 where in the brackets are indicated the physical states of reagents: (gr), graphite; (g), gaseous; (cr), crystalline.  
71 IV.

## 72 5 Conclusions

73 The general aim of these investigations was to report the results of the thermodynamic study of the  
74 cyanocobalamin. The standard enthalpy of formation is determined by using combustion calorimetry. Much  
75 of the work is devoted to the study of the mechanism of combustion of cyanocobalamin and determination of  
76 thermodynamic functions of the combustion products.

77 V.

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Value	Experiment					
	1	2	3	4	5	6
m sam (g) a	0.15345	0.1600	0.1466	0.1580	0.1595	0.1607
m par (g) a	0.6920	0.6989	0.7024	0.7021	0.7050	0.7059
m thread (g) a	0.0025	0.0022	0.00235	0.00215	0.0021	0.0019
W (J?K -1 ) b	14805	14805	14805	14805	14805	14805
?t (K) c	2.443105	2.475455	2.463935	2.483210	2.493200	2.497915
?? ? U ? (J) d	36170.4	36649.1	36478.6	36763.9	36911.8	36981.6
?? ? U par (J) e	32347.1	32668.2	32831.8	32818.7	32954.3	32997.3
?? ? U thread (J) e	42.0	36.5	39.3	36.0	35.6	32.1
?? ? U HNO3 (J) f ?? ? U ? (J) g	5.9 9.8	10.5	8.2 16.4	8.8 -	9.4	11.7
		16.4			26.2	26.2
?? ? U (J?g -1 ) h	24666.0	24689.4	24663.7	24686.1	24694.0	24683.9

?? ? U = 24681 ± 10 J?g -1 = 33452 ± 14 kJ?mol -1 , the mean energy of combustion of cyanocobalamin, ?? ?  
 °U = 33435 ± 14 kJ?mol -1 , the energy of combustion of cyanocobalamin at standard pressure.

[Note: a b W, the energy equivalent of the calorimeter. c ?t, the temperature increase in the experience, adjusted for heat transfer.d h ? ?]

Figure 1: Table 1 :

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Figure 2: Table 2 :



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