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FRAFLUOROHYDRAZINE



ARLINGTON HALL

Cates Tag

 $(F_2 N - NF_2)$

A NEW HIGH ENERGY LIQUID

OXIDIZER

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ROHM & HAAS COMPANY

REDSTONE ARSENAL RESEARCH DIVISION HUNTSVILL, ALABAMÁ

REPORT NO. 3.22

TETRAFLUOROHYDRAZINE (A new, high energy, liquid oxidizer)

Вy

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ABSTRACT

Tetrafluorohydrazine gives the following theoretical equilibrium impulses at 1000 psi \rightarrow 1 atmos. with the following fuels:

 $H_2 = 357 \text{ sec.}; N_2H_4 = 333; NH_3 = 325; UMDH = 316; Hidyne = 314; JP=4 = 297.$

The physical and chemical properties of tetrafluorohydrazine indicate that it would be much easier to handle as a liquid oxidizer than are oxygen and fluorine.

The preparation, chemical and physical properties of tetrafluorohydrazine are briefly reviewed.

TETRAFLUOROHYDRAZINE (A now, high energy, liquid oxidizer)

The new heat of formation of tetrafluorohydrazine $(\Delta H_{125}^{0}[N_{2}F_{4}] = -2.0 \pm 2.5 \text{ kcal./mole}^{1}$ which has been determined at the National Bureau of Standards indicates that tetrafluorohydrazine is a much more powerful oxidizer than previously believed. The calculated specific impulses (Table I) would indicate that tetrafluorohydrazine lies between fluorine and oxygen in its oxidative power. However, its physical and chemical properties are much superior to those of fluorine and oxygen (Table II).

It is immediately apparent from these tables that tetrafiuorohydrazine exceeds liquid oxygen with all fuels listed (except hydrogen) and approaches the performance of fluorine as a high energy oxidizer. Because of its properties outlined in Table II it is much more readily handled than are fluorine and oxygen. Because of its low freezing point it would be easier to work with than is N_2O_4 .

Preparation

The preparation of tetrafluorohydrazine was first reported by this Division². Subsequent reports in this series have covered the development of satisfactory methods of preparation of tetrafluorohydrazine.

Tetrafluorohydrazine is prepared by the reaction at elevated temperatures of nitrogen trifluoride with various fluorine acceptors such as copper, arsenic, bismuth, antimony, glass^{2,3} and carbon⁴. The

- ² Rohm & Haas Company Quarterly Progress Report on Physical Chemistry, No. P-57-4, March, 1957.
- ³ C. B. Colbarn, A. Kennedy JACS 80 5004 (1958).
- ⁴ Stauffer Chemical Company Report No. SCC-26-QPR-6 December, 1958.

¹ G. T. Armstrong, S. Marantz and C. F. Coyle NBS Report No. 6584 October, 1959.

- 2 -

Table I

Fuel	Oxidizer					
	N ₂ F	4	$\begin{array}{c} N_2 F_4^2 \\ N_2 O_4 \end{array}$	O ₂	F2	N ₂ O ₄
Hydrogen	Eq. ³	357		391 ⁵	410 ⁵	
, ,	Froz. ⁴	37.6		3885	3985	
Hydrazine	Eq.	333		3135	363 ⁵	2925
	Froz.	311		301 ⁵	3345	2835
Ammonia	Eq.	325		2945	357 ⁵	
	Froz.	306		285 ⁵	330 ⁵	
UDMH	Eq.	316		310 ⁵		2855
	Froz.	286		295 ⁵		2745
Hidyne ⁶	Eq.	314		3065		2825
	Froz.	285		291 ⁵		271 ⁵
JP-4	Eq.	297	314	3005,7		2765,7
(CH1-93)	Froz.	277	285	286577		2635; 1

Specific Impulse¹ of Various Fuel Oxidizer Mixtures

¹ Calculated using ideal assumptions, expansion from 1000 psi to 1 atmos. See Rohm & Haas Company Quarterly Progress Report on Physical Chemistry, No. P-59-16; September 1959 for method of calculation.

² For carbon containing fuels it is interesting to note that a mixture of N_2F_4 and N_2O_4 will give a higher specific impulse than either constituent alone in much the same fashion as do mixtures of F_2 and O_2 .

³ Equilibrium flow.

⁴ Frozen flow.

⁵ These values taken from Rocketdyne Publication 505-X (Revised 1 January, 1959) "Rocket Engine Propellants".

⁶ 60 wt. % UDMH 40% diethylene triamine.

⁷ These numbers are for RP-1 which has essentially the same composition as JP-4.

-3-

Table II

Physical Properties of N_2F_4 , F_2 , O_2 , N_2O_4

	N ₂ F ₄	O ₂ ³	F ₂ ³	N2043
Boiling Point ⁰ C.	-73 ¹	-183	-187	21.3
Melting Point [°] C.	-163 (approx.)	-218.4	-223	-9.3
Critical Temperature ^o C.	36 ¹			
Critical Pressure Atmos.	771			
Density gr/cc.	1.5^2 at -100° C.	1.14	-183 1.1	08 ⁻¹⁸⁷ 1, 491 ⁰

¹ C. B. Colburn, A. Kennedy JACS <u>80</u> 5004 (1958).

² Private communication Charles Mader to W. D. Niederhauser, 15 November, 1957.

³ These values are handbook values.

reaction can be represented as follows:

 $2 NF_3 + M \rightarrow N_2 F_A + 2 MF$

The loss of hydrogen by difluoroamine¹on contact with various solids has also been reported to give tetrafluorohydrazine². Electrical discharge through nitrogen trifluoride has also been reported as yielding tetrafluorohydrazine and difluorodiazine³.

By far the most efficient method for the preparation of tetrafluorohydrazine is the reaction of nitrogen trifluoride with carbon⁴. This process has been perfected⁵ and scaled up to a small pilot plant operation⁶. The yield of tetrafluorohydrazine from the reaction of nitrogen trifluoride with carbon is 77% at 440°C. with a residence time

³ J. W. Frazer, I. Inorg. Z Nuclear Check 11 165-167 (1959).

⁴ Stauffer Chemical Company Report No. SCC-26-QPR-6 December, 1958.

⁵ Rohm & Haas Company Quarterly Progress Report on Physical Chemistry, No. P-59-9, June, 1959.

⁶ Ibid., Quarterly Progress Report on Solid Propellant Processing, No. P-59-8 June, 1959.

¹ A. Kennedy, C. B. Colburn, JACS <u>81</u> 2906 (1959). ² E. A. Lawton & J. Q. Weber JACS <u>81</u> 4755 (1959).

of approximately six seconds. The yield is between 62-71% in the reaction between nit: ogen trifluoride and copper at 375°C. and with a residence time of thirteen minutes. These reactions take place in flow systems and hence the process can be made continuous.

Summary of Chemical and Physical Properties of Tetrafluorohydrazine

Heat of formation	= -2, 0 + 2.5 kcal./mole
Boiling point	- 73°C.
Freezing point	-162°C. (approximately)
Critical temperature	36 [°] C.
Critical pressure	77 atmospheres (estimated)
Vapor pressure equation	
log P(mm)	= -692/T + 6.33
Density	1.5 gm./cc.
Heat of vaporization	3170 cal. /mole

Mass Cracking Pattern of N₂F₄

(Obtained on a Consolidated Model 620 Mass Spectrometer)

m/e	Ion	Pattern
52	NF ₂ ⁺	90.6%
33	NF ⁺	100.0%
28	N_2^+	7.7%
19	· F ⁺	4.7%
14	Nt	8.5%

Sensitivity div./m 19.0

Instrument sensitivity 100 div. /m for m/e 43 for n butane.

Infrared Spectrum of Tetrafluorohydrazine

(See Fig. 1 Rohm & Haas Company QPR Physical Chemistry Report, P-57-10, page 1).

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-4-

- 5 ...

N. M. R. Spectrum of Tetrafluorohydrazine

The F^{19} nuclear magnetic resonance spectrum of tetrafluorohydrazine consisted of a single broad unresolved band at a field of approximately 75 p. p. m. lower than that of the F^{19} nuclei of trifluoroacetic acid. Resolution of the expected triplet was not observed.

Tetrafluorohydrazine does not react with glass, stainless steel, or copper at room temperature.

Summary

Tetrafludrohydrazine is a powerful new oxidizing agent which lies between liquid oxygen and liquid fluorine in its performance behaviour with conventional fuels. Its physical and chemical properties indicate that it would be much easier to work with in liquid motors than fluorine and oxygen. Tetrafluorohydrazine can be prepared with relative ease.

Acknowledgement

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