

MICROWAVE SPECTROSCOPY INFORMATION LETTER

No. XXXII

May 15, 1989

Compiled by:

R. H. Schwemdeman  
Departmen of Chemistry  
Michigan State University

Lise Nygaard  
Chemical Laboratory V  
University of Copenhagen  
H. C. Ørsted Institutet  
5. Universitetsparken  
DK-2100 Copenhagen, Denmark



## MICROWAVE SPECTROSCOPY NEWSLETTER XXXII

## List of Contributors

No.	Contributor	Institute	Page
1	Alonso, J. L.	Universidad de Valladolid	1
2	Amano, T.	Herzberg Institute of Astrophysics	3
3	Baker, J. G.	University of Manchester	4
4	Bauder, A.	Swiss Federal Institute of Technology	5
33	Bauer, A.	Universite de Lille	43
5	Beaudet, R.	University of Southern California	6
33	Bogey, M.	Universite de Lille	43
6	Bohn, R. K.	University of Connecticut	7
44	Botskor, I.	University of Ulm	54
7	Brown, R. D.	Monash University	8
15	Caminati, W.	Universita di Bologna	20
8	Carpenter, J. H.	The University, Newcastle-upon-Tyne	9
15	Cazzoli, G.	Universita di Bologna	20
9	Christen, D.	Universität Tübingen	10
39	Cohen, E. A.	Jet Propulsion Laboratory	49
49	Cook, R. L.	Mississippi State University	61
11	Cox, A. P.	University of Bristol	12
15	Damiani, D.	Universita di Bologna	20
12	Danieli, R.	Ist. di Spettroscopia Molekulare del C.N.R.	13
33	Demuynck, C.	Universite de Lille	43
33	Destombes, J.	Universite de Lille	43
12	Di Lonardo, G.	Ist. di Spettroscopia Molekulare del C.N.R.	13
13	Dreizler, H.	Universität Kiel	14
14	Durig, J. R.	University of South Carolina	19

No.	Contributor	Institute	Page
24	Endo, Y.	Institute of Molecular Science	30
15	Favero, P. G.	Universita' di Bologna	20
16	Galica, J.	Polish Academy of Sciences	21
17	Gerry, M. C. L.	University of British Columbia	22
18	Gillies, C. W.	Rensselaer Polytechnic Institute	23
19	Graybeal, J. W.	Virginia Polytechnic Institute	25
13	Guarnieri, A.	Universität Kiel	14
20	Gutowsky, H. S.	University of Illinois	26
21	Harmony, M. D.	University of Kansas	27
22	Hausamann, D.	German Aerospace Research Est.	28
23	Hayashi, M.	Hirosshima University	29
27	Hillig, K. W.	University of Michigan	35
24	Hirota, E.	Institute of Molecular Science	30
25	Hoeft, J.	Freie Universität Berlin	32
44	Jones, H.	University of Ulm	54
44	Kreiner, W. A.	University of Ulm	54
26	Kroto, H. W.	University of Sussex	33
27	Kuczkowski, R. L.	University of Michigan	35
28	Kukolich, S.	University of Arizona	36
10	Larsen, N. W.	University of Copenhagen	11
29	Lees, R. M.	University of New Brunswick	37
30	Legon, A. C.	University of Exeter	38
31	Li, Y.-S.	Memphis State University	40
32	Lovas, F. J.	National Inst. of Standards and Technology	41
33	Macke, B.	Universite de Lille	43
13	Mäder, H.	Universität Kiel	14

No.	Contributor	Institute	Page
34	Matsumura, C.	National Chemical Laboratory for Industry	44
35	Millen, D. J.	University College London	45
36	Mollendal, H.	University of Oslo	46
25	Nair, K. P. R.	Freie Universität Berlin	32
37	Nandi, R. N.	Saha Institute of Chemical Physics	47
10	Nygaard, L.	University of Copenhagen	11
38	Ogata, T.	Shizuoka University	48
10	Pedersen, T.	University of Copenhagen	11
39	Pickett, H. M.	Jet Propulsion Laboratory	49
39	Poynter, R. L.	Jet Propulsion Laboratory	49
40	Pozdeev, N.	Academy of Sciences of the USSR	50
41	Quade, C. R.	Texas Technological University	51
42	Roberts, J.	North Texas State University	52
43	Roussy, G.	Universite de Nancy	53
44	Rudolph, H. D.	University of Ulm	54
45	Saito, S.	Nagoya University	56
46	Sasada, Y.	Kanto Gakuin University	58
29	Sastry, K. V. L. N.	University of New Brunswick	37
12	Scappini, F.	Ist. di Spettroscopia Molecolare del C.N.R.	13
47	Schwendeman, R. H.	Michigan State University	59
48	Shimizu, T.	University of Tokyo	60
8	Smith, J. G.	The University, Newcastle-upon-Tyne	9
10	Sorenson, G. O.	University of Copenhagen	11
49	Su, C. F.	Mississippi State University	61
32	Suenram, R. D.	National Inst. of Standards and Technology	41
34	Sugie, M.	National Chemical Laboratory for Industry	44

No.	Contributor	Institute	Page
13	Sutter, D. H.	Universitat Kiel	14
50	Takami, M.	Institute of Physical and Chemical Research	62
51	Takano, M.	National Defense Academy (Japan)	63
34	Takeo, H.	National Chemical Laboratory for Industry	44
52	Tanaka, K.	Kyushu University	64
52	Tanaka, T.	Kyushu University	64
53	Tanimoto, M.	Shizuoka University	66
25	Torring, T.	Freie Universitat Berlin	32
54	True, N. S.	University of California, Davis	67
55	Tsunekawa, S.	Toyama University	68
56	Winnewisser, B. P.	Justus-Liebig Universitt Giessen	69
57	Winnewisser, G.	University zu Kln	70
56	Winnewisser, M.	Justus-Liebig Universitt Giessen	69
24	Yamada, C.	Institute of Molecular Science	30
57	Yamada, K. M. T.	University zu Kln	70
58	Yamaguchi, I.	Sophia University	71

1 Name to whom queries should be addressed Prof. José L. Alonso

Mailing address Departamento de Química-Física  
Facultad de Ciencias. Universidad de Valladolid  
Valladolid. 47005. SPAIN

Telephone: (3483)-264000.ex2437 Telefax: (983)257293

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>5</sub> H <sub>8</sub> O <sub>2</sub>	α-Methyl-γ-Butyrolactone	J.C. López J.L. Alonso F.J. Peláez	J.Mol. Spectrosc. <u>131</u> , 9 (1988)
C <sub>3</sub> H <sub>6</sub> O <sub>2</sub> S	Trimethylene sulphone	J.C. López J.L. Alonso	J.Chem.Phys. <u>89</u> , 678 (1988)
C <sub>4</sub> H <sub>6</sub> O <sub>3</sub>	Propylene Carbonate	E.González J.L.Alonso <sup>1</sup> W.Caminati	Spectrum assigned Work in progress.
C <sub>4</sub> H <sub>8</sub> O <sub>2</sub> S	Butadiene Sulphone	J.L.Alonso D.G.Lister J.C. López R.M.Villamañán	Paper in preparation.
C <sub>4</sub> H <sub>6</sub> O <sub>2</sub>	β-Butyrolactone	J.L.Alonso E.González	Paper in preparation.
C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>	3-Methyl oxetane	J.C. López J.L.Alonso <sup>2</sup> H.Wieser	Paper in preparation.
C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>	2-Methyl oxetane	J.C. López J.L.Alonso <sup>2</sup> H.Wieser <sup>2</sup>	Spectrum assigned. Work in progress.
C <sub>7</sub> H <sub>5</sub> OF	o-Fluorobenzaldehyde	R.M.Villamañán	J.Chem.Soc.Far.II
C <sub>7</sub> H <sub>5</sub> OF	m-Fluorobenzaldehyde	J.L.Alonso	(in press)
C <sub>7</sub> H <sub>5</sub> OF	p-Fluorobenzaldehyde	S.R.González R.M.Villamañán J.L.Alonso	J.Mol.Struct. <u>190</u> , 79 (1988)
C <sub>8</sub> H <sub>7</sub> F	o-Fluorostyrene	R.M.Villamañán J.C. López J.L.Alonso	Submitted

$C_8H_7F$	m-Fluorostyrene	R.M.Villamañán J.L.Alonso	Submitted
$C_6H_9NO_2$	1-Nitrocyclohexene	S.R.González R.Mulas J.L.Alonso	J.Mol.Spectrosc. (in press)
$C_5H_4N_2O_2$	p-Nitropyridine	R.Mulas S.R.González J.L.Alonso	J.Mol.Struct. (in press)
$C_4H_7OF$	Butyryl Fluoride	S.R.González J.L.Alonso	Syn-Anti and Syn-Gauche Confor- mers assigned.
$C_7H_4NOF$	m-Fluorophenyl isocyanate	M.E.Charro J.L.Alonso	"cis" Conformer assigned.
$C_4H_6OS$	Tetrahydrothio- phene-3-one	R.M.Villamañán J.C.López J.L.Alonso	Spectrum assigned. Work in progress.

<sup>1</sup>Istituto di Spettroscopia Molecolare CNR and Istituto di Chimica  
Fisica e Spettroscopia. Universita di Bologna (Italy).

<sup>2</sup>University of Calgary. Alberta. Canada.

2 Name to whom queries should be addressed Takayoshi Amano

Mailing address Herzberg Institute of Astrophysics

National Research Council, Ottawa, Ontario, Canada K1A 0R6

Telephone: 613-990-0737 Telefax: 613-952-0974

Electronic mail (BITNET or other): TYA@NRCVMS1

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CNS	NCS radical	T. Amano	Assigned
NS <sub>2</sub>	Nitrogen disulfide	T. Amano	Spectrum recorded
C <sub>2</sub> H <sub>6</sub> (CH <sub>3</sub> CD <sub>3</sub> )	Ethane-d <sub>3</sub>	N. Moazzen-Ahmadi	mm-wave spectrum in torsional excited state

3 Name to whom queries should be addressed Dr John G Baker

Mailing address Schuster Laboratory

University of Manchester

Manchester M13 9PL, United Kingdom

Telephone: (061) 275 4137 Telefax: (061) 273 5867

Electronic mail (BITNET or other): MBCMMJG@UK.AC.UMRCC.CMS

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>2</sub> H <sub>4</sub> O	acetaldehyde	W Liang	far IR torsion spectrum
C <sub>2</sub> H <sub>4</sub> O <sub>2</sub>	methyl formate	H Nadgaran	mmwave torsional states
CF <sub>3</sub> SF <sub>5</sub>	trifluoromethyl sulfur hexafluoride	P N Brier	mmwave supersonic jet high resolution spectrum

4 Name to whom queries should be addressed: Alfred Bauder  
 Mailing address: Laboratorium für Physikalische Chemie  
 ETH - Zentrum  
 CH-8092 Zürich, Switzerland  
 Telephon number: 01 - 256 4341  
 Electronic mail: BAUDER@CZHETH5A (EARN)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{CH}_3\text{F}_3\text{Si}$ ( $\text{CH}_3\text{SiF}_3$ )	Methyl trifluoro-silane	I. Ozier C. Styger	$\Delta K=\pm 3$ transitions started
$\text{CH}_4\text{O}_2$ ( $\text{CH}_3\text{OOH}$ )	Methyl hydroperoxide	A. Bauder in collab. with C. E. Blom (Giessen)	manuscript in preparation
$\text{C}_2\text{H}_3\text{F}_3$ ( $\text{CH}_3\text{CF}_3$ )	1,1,1-Trifluoroethane	I. Ozier B. Vogelsanger	$\Delta K=\pm 3$ transitions assigned
$\text{C}_2\text{H}_5\text{NO}_2$ ( $\text{HOONHOCH}_3$ )	Methoxyformamid	C. Styger	assigned
$\text{C}_3\text{H}_6$	Cyclopropane	T. Brupbacher	centrifugal distort. spectrum measured
$\text{C}_3\text{D}_4$ ( $\text{CD}_2=\text{C=CD}_2$ )	Allene-d <sub>4</sub>	B. Vogelsanger	$\nu_{10}$ and $\nu_{11}$ excited states measured
$\text{C}_4\text{H}_7\text{N}$	Pyrroline	J. Dommen W. Caminati R. Meyer	manuscript in preparation
$\text{C}_6\text{H}_{12}$	Cyclohexane	J. Dommen T. Brupbacher	substitution structure
$\text{C}_7\text{H}_8$	Quadraticane	B. Vogelsanger	manuscript submitted to J. Mol. Spectr.
$\text{C}_2\text{DF}_3\text{O}_2$ ( $\text{CF}_3\text{COOD}$ )	Trifluoroacetic acid-d <sub>1</sub>	L. Martinache M. Wegener	deuterium quadrupole splittings
$\text{C}_2\text{H}_3\text{DO}$ ( $\text{CH}_3\text{CDO}$ )	Acetaldehyde-d <sub>1</sub>	L. Martinache	deuterium quadrupole splittings
$\text{C}_2\text{H}_3\text{DO}_2$ ( $\text{CH}_3\text{COOD}$ )	Acetic acid-d <sub>1</sub>	L. Martinache M. Wegener	deuterium quadrupole splittings
$\text{C}_6\text{H}_4\text{DF}$	Fluorobenzene-d <sub>1</sub>	S. Jans-Bürli	deuterium quadrupole splittings of ortho, meta and para isomer
$\text{C}_6\text{H}_5\text{D}$	Benzene-d <sub>1</sub>	S. Jans-Bürli	deuterium quadrupole splittings remeasur.
$\text{C}_3\text{H}_3\text{F}_3\text{O}_4$ ( $\text{HCOOH}..\text{CF}_3\text{COOH}$ )	Formic acid - trifluoroacetic acid dimer	W. Kresa L. Martinache M. Wegener	D, <sup>13</sup> C, <sup>18</sup> O species assigned
$\text{C}_4\text{H}_5\text{F}_3\text{O}_4$ ( $\text{CH}_3\text{COOH}..\text{CF}_3\text{COOH}$ )	Acetic acid - trifluoroacetic acid dimer	W. Kresa	assigned
$\text{F}_3\text{OP}$ ( $\text{OPF}_3$ )	Phosphoryl trifluoride	C. Styger	$\Delta K=\pm 3$ transitions measured

5 Name to whom queries should be addressed Robert A. Beaudet

Mailing address Department of Chemistry  
University of Southern California  
Los Angeles, CA 90089-0432

Telephone: (213) 743-2997 Telefax: 213-746-5679

Electronic mail (BITNET or other): BEAUDET % RAMOTH@USC-ECL.ARPA

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CHFO <sub>2</sub>	carbon dioxide-hydrogen fluoride cluster	Steven Sharpe, Yongping Zeng	Spectra taken and assigned at 2350 cm <sup>-1</sup>
CHClO <sub>2</sub>	carbon dioxide-hydrogen chloride cluster	Steven Sharpe, Yongping Zeng	Spectra taken and assigned at 2350 cm <sup>-1</sup>
CHBrO <sub>2</sub>	carbon dioxide-hydrogen bromide cluster	Steven Sharpe, Yongping Zeng	Spectra taken at 2350 cm <sup>1</sup>
HFN <sub>2</sub> O	hydrogen fluoride-nitrous oxide cluster	Steven Sharpe, Yongping Zeng	Spectra taken and assigned

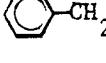
6 Name to whom queries should be addressed

Robert K. Bohn

Mailing address Dept. of ChemistryUniv. of ConnecticutStorrs CT 06268 USATelephone: (203) 486-3044

Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): RBOHN@UCONNVM!

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{ArF}_3\text{P} (\text{PF}_3\text{-Ar})$	Argon-PF <sub>3</sub> Complex	U. of Michigan	Submitted
$\text{F}_3\text{KrP} (\text{PF}_3\text{-Kr})$	Krypton-PF <sub>3</sub> Complex	collaboration.	
$\text{C}_3\text{H}_5\text{NO}_2 (\text{CH}_2=\text{CHCH}_2\text{ONO})$	Allyl Nitrite	K. W. Hwang	4 conformers, 2 are assigned.
$\text{C}_4\text{H}_5\text{ArN}$	 Pyrrole-Ar	U. of Michigan collaboration.	J. Phys. Chem. In press.
$\text{C}_4\text{H}_8\text{N}_2\text{O}$	 Nitrosopyrrolidine	K. W. Hwang	$\tau_0$ structure, $^{14}\text{N}$ quad. coupling. Also, $d_4$ species.
$\text{C}_5\text{H}_{10}\text{N}_2\text{O}$	 Nitrosopiperidine	X. Z. Liu K. W. Hwang	$\tau_0$ structure.
$\text{C}_6\text{H}_{11}\text{NO}$	 1-Formylpiperidine	C. Sahi	Assigned.
$\text{C}_8\text{H}_7\text{N}$	 Benzyl Cyanide	U. of Calif.(Davis)	Ground state, $^{14}\text{N}$ quad. collaboration. coupling assigned.
$\text{C}_8\text{H}_8\text{O}_2 (\text{OHC}-\text{C}_6\text{H}_4-\text{OCH}_3)$	p-Anisaldehyde	U. of Calif.(Davis)	2 conformers assigned. collaboration.
$\text{C}_{10}\text{H}_{12}\text{O}(\text{OHC}-\text{C}_6\text{H}_4-\text{CH}(\text{CH}_3)_2)$	p-Isopropyl Benzaldehyde	R. Bohn	2 conformers assigned.

7 Name to whom queries should be addressed PROFESSOR R.D. BROWN

Mailing address CHEMISTRY DEPARTMENT  
MONASH UNIVERSITY  
WELLINGTON ROAD, CLAYTON VIC 3168, AUSTRALIA

Telephone: 61 3 565 4550 Telefax: (61) (3) 565 4597

Electronic mail (BITNET or other): CHE201N@VAXC.CC.MONASH.EDU.AU

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{C}_2\text{HFO}$ (HFC=C=O)	Fluoroketene	K. Wiedenmann	Vibrational satellites analysed. Manuscript in press, J. Mol. Spec.
$\text{C}_4\text{H}_4\text{N}_2\text{O}_2$	Uracil	A. Pierlot	J. Am. Chem. Soc., 1988, <b>110</b> , 2239.
$\text{C}_5\text{H}_6\text{N}_2\text{O}_2$	Thymine	A. Pierlot	J. Chem. Soc., Chem. Comm., 1988, in Press.
$\text{C}_4\text{H}_5\text{N}_3\text{O}$	Cytosine	A. Pierlot	J. Am. Chem. Soc., 1988, in press.
$\text{C}_5\text{H}_5\text{N}_5$	Adenine	A. Pierlot	Chem. Phys. Letts., 1988, in press.
$\text{C}_3\text{H}_6\text{O}_3$	Glyceraldehyde	S. Mandragona	Analysis in progress.
$\text{C}_2\text{H}_5\text{N}$ ( $\text{H}_2\text{C}=\text{CH-NH}_2$ )	Vinylamine	D. McNaughton	$0^+, 0^-$ mm wave spectrum analysed. Manuscript in preparation.
$\text{C}_2\text{HClO}$	Chloroketene	G. Burns	Vibrational satellites assigned. Manuscript in preparation.
$\text{CH}_3\text{NO}$ ( $\text{HCONH}_2$ )	Formamide	F. Trollope	Vibrational satellite analysis in progress.
$\text{C}_4\text{H}_6\text{N}_2\text{O}_2$	2,5-Piperazinedione (glycine anhydride)	F. Trollope	Analysis in progress.
$\text{C}_2\text{H}_2\text{O}$ ( $\text{H}_2\text{C}=\text{C=O}$ )	Ketene	W. Taylor	mm wave spectrum of $^{13}\text{C}$ and $^{18}\text{O}$ species assigned.

8 Name to whom queries should be addressed: Dr J.H. Carpenter

Mailing address Department of Chemistry,

The University, Newcastle-upon-Tyne,

NE1 7RU, United Kingdom.

Telephone: (091)-232 8511 Telefax: (091)-261 1182

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
<chem>C1F3Si</chem>	Chlorotrifluorosilane	J.H.Carpenter	mm-wave spectrum of ground & excited states
<chem>CDCl3</chem>	Chloroform-d	J.H.Carpenter	mm-wave spectrum of excited states
<chem>CH3Cl3Si</chem>	Methyltrichlorosilane	J.G.Smith	manuscript sent to J. Mol. Spectrosc.
<chem>CH3F3Si</chem>	Methyltrifluorosilane	J.G.Smith	mm-wave spectrum of excited states
<chem>C4H9CCl</chem> <chem>[(CH3)3Cl]</chem>	2-chloro-2-methylpropane ( <u>tert</u> -butyl chloride)	J.G. Smith	mm-wave spectrum of excited states

9 Name to whom queries should be addressed Dines Christen

Mailing address Institut für Physikalische und Theoretische Chemie  
Universität Tübingen  
Auf der Morgenstelle 8, D-7400 Tübingen, B.R.D.

Telephone: (07071) 29 69 24 Telefax:

Electronic mail (BITNET or other): CFOHØØ1@DTUZDV5A

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{CH}_2\text{F}_3\text{N}$	$(\text{CF}_3\text{NH}_2)$ Trifluoromethylamine	P. Zylka	Normal + Deuteriated species
$\text{C}_2\text{H}_6\text{O}_2$	$(\text{CH}_3\text{COCH}_3)$ Dimethylperoxide	P. Zylka	Combination with RFMWDR and FIR
$\text{C}_2\text{F}_3\text{N}_2\text{S}_2$	$(\text{CF}_3\text{C}(\text{N-S})_2)$ 4-(Trifluoromethyl)-1,2,3,5 dithiadiazole	E. Jaudas	Broad band and HR spectra
$\text{F}_6\text{OS}$	$(\text{SF}_5\text{OF})$ Fluoroxysulphurpentafafluoride	E. Jaudas	RFMWDR
$\text{C}_2\text{FH}_6\text{N}$	$((\text{CH}_3)_2\text{NF})$ Dimethylfluoramidine	J. Kadel	Structure with ED
$\text{C}_2\text{F}_3\text{NO}$	$(\text{CF}_3\text{NCO})$ Trifluoromethylisocyanate	D. Christen	LAM analysis with J. Koput, Poznan
$\text{C}_2\text{H}_8\text{O}_2$	$((\text{CH}_2\text{OH})_2)$ Ethylene glycol	D. Christen	MWMWDR, LAM analysis with H. Møllendal, Oslo

10 Name to whom queries should be addressed

Staff member indicated by \*) below

Mailing address      Chemical Laboratory V  
 UNIVERSITY OF COPENHAGEN  
 H. C. Ørsted Institutet  
 5, Universitetsparken  
 DK-2100 Copenhagen, Danmark

Telephone    +45 31 35 31 33      Telefax    45 1 35 06 09

Electronic mail    KM5LN at DKUCCCl1

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>2</sub> H <sub>3</sub> N <sub>3</sub>	1H- and 2H- 1,2,3-triazole		Acta Chem. Scand. <u>A</u> 42 (1988) 500-514.
C <sub>3</sub> H <sub>2</sub> O <sub>4</sub>	1,3-dioxolane- 4,5-dione	*)N.W.Larsen	Vibr. states assigned.
C <sub>5</sub> H <sub>7</sub> NO	furfurylamine and five D species	*)T.Pedersen I.Hedgecock N.W.Larsen L.Nygaard G.O.Sørensen	Nearly finished.
C <sub>7</sub> H <sub>4</sub> F <sub>2</sub> O	3-fluorobenzoyl fluoride		J. Mol. Spectrosc. <u>129</u> (1988) 223-231.
C <sub>7</sub> H <sub>4</sub> F <sub>2</sub> O	4-fluorobenzoyl fluoride		J. Mol. Spectrosc. <u>128</u> (1988) 370-383.

11 Name to whom queries should be addressed Dr. A.P. Cox

Mailing address Department of Physical Chemistry, School of Chemistry,  
University of Bristol, Cantock's Close, BRISTOL, BS8 1TS, U.K.

Telephone: (0272) 303687 Telefax: (0272) 251295

Electronic mail (BITNET or other): COX % BRIS.CSA

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CC <sub>2</sub> F <sub>2</sub> NO	chlorodifluoronitrosomethane	A.C. Fettis	In progress.
CF <sub>3</sub> NO <sub>2</sub>	trifluoroniromethane	P.R.R. Langridge-Smith	In manuscript.
CHF <sub>3</sub> S ) C <sub>2</sub> H <sub>5</sub> S )	trifluoromethylmercaptan	R. Stevens/C.A. Rego	Manuscript prepared for J.Mol.Struct.
CH <sub>5</sub> BO <sub>2</sub> [CH <sub>3</sub> B(OH) <sub>2</sub> ]	methylborondihydroxide	C.A. Rego	In thesis.
C <sub>2</sub> HC <sub>1</sub> F <sub>2</sub> O	chlorodifluoroacetaldehyde	D.W. Knight	Manuscript in preparation.
C <sub>2</sub> H <sub>5</sub> NO	nitrosoethane	J.A. Hardy	Extended study. (see Sussex)
C <sub>3</sub> H <sub>2</sub> N <sub>2</sub> [CH <sub>2</sub> (CN) <sub>2</sub> ]	malononitrile	J. Randell	Structure/quadrupole complete.
C <sub>4</sub> H <sub>9</sub> BF <sub>2</sub>	t-butylborondifluoride	S.D. Hubbard	Barrier/dipole complete.
C <sub>4</sub> H <sub>9</sub> NO	t-butylnitrosomethane	M.J. Corkill	In thesis.
C <sub>4</sub> H <sub>9</sub> NO	t-butylnitrite	M.J. Corkill	Manuscript in preparation.
C <sub>5</sub> H <sub>10</sub> O	pivalaldehyde (t-butyl) formaldehyde	A.D. Couch	Isotopes assigned.
CF <sub>3</sub> NSi	trifluorosilylisocyanide	A.D. Couch/S. Clement	In Chem.Phys.Letters 1989.
CH <sub>3</sub> NSi	trifluorosilylcyanide		
N <sub>2</sub> O <sub>3</sub>	dinitrogentetroioxide	J. Randell	Vibrational satellites and Force field.

12 Name to whom queries should be addressed Dr. Roberto Danieli

Mailing address Istituto di Spettroscopia Molecolare del C.N.R.

Via de' Castagnoli, 1

40126 Bologna, Italy

Telephone: 51-287007 Telefax: 51-229702

Electronic mail (BITNET or other): AB00501 at ICINECA

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CH <sub>3</sub> Br	methyl bromide	F. Scappini	collision induced
CH <sub>3</sub> I	methyl iodide	H. Mäder	transitions
G. Di Lonardo			
SbH <sub>3</sub>	stibine	B.M. Dinelli	RF-IR double
		F. Scappini	resonance
		G. Di Lonardo	
C <sub>4</sub> O <sub>2</sub> Br <sub>2</sub>	1,2-dibromo cyclobuten-3,4-dione	G. Corbelli	
C <sub>2</sub> HF <sub>5</sub> ClBr	halothane	B. Lunelli	MW-spectrum
		F. Scappini	

13 H.Dreizler, A.Guarnieri, H.Mäder, D.Sutter

Mailing address: Institut für Physikalische Chemie  
 -Abteilung Chemische Physik-  
 Christian-Albrechts-Universität zu Kiel  
 Ludewig-Meyn-Str. 8  
 D-2300 Kiel, FR Germany

Electronic Mail: PHC58@DKIUNI0 (EARN)

Telephone Number: 0431/880-2753, -2752, -2751, -2009, -2750

Formula	Name of Compound	Name of Investigator	Present Stage of Progress
$\text{BF}_2\text{OH}$	hydroxodifluoroboron	Vormann	B-hfs complete
$\text{BF}_2\text{OD}$			
$\text{BF}_3$	trifluoroboron	Vormann Stahl	hfs in progress
$\text{BH}_4\text{N}^+$			
$\text{BH}_2\text{NH}_2$	aminoborane	Vormann	N, B-hfs in progress
$\text{FC}^{14}\text{N}$	cyanogen fluoride	Zerbe-Foese	dipole moments of
$\text{FC}^{15}\text{N}$		Hinze	excited states
OCS	carbonyl sulfide	Wolf	relaxation studies
$\text{CD}_2\text{I}_2$	diiodomethane-d <sub>2</sub>	Stahl	assignment in progress
HCN	hydrogen cyanide	Heineking	l-type spectra,
DCN			N-hfs reanalysed
$\text{HNC}^{17}\text{O}$	isocyanic acid	Gerry	N,O-hfs in progress
DNCO		Heineking	N,D-hfs complete
		in coop. with	
		Howard, Vancouver	
$\text{CHClF}_2$	chlorodifluoromethane	Spiehl	mmw spectrum
$\text{CH}_2\text{Br}_2$	dibromomethane	Merke Heineking	Br-hfs in progress
$\text{CH}_2\text{I}_2$	diiodomethane	Stahl in coop. with Vogelsanger, Zürich	assignment in progress
$\text{CH}_2\text{N}^+$			
$\text{CH}_2\text{NH}$	methanimine	Krause	rotational Zeeman effect in progress
$\text{CH}_3\text{NO}_3$	methyl nitrate	Spieckermann	rotational Zeeman effect in progress
$\text{CH}_3\text{Ns}$	methyl azide	Gerry Heineking	incl. <sup>15</sup> N-species, N-hfs ms. submitted
$\text{CH}_4$	methane	Stahl Wolf	hfs, ms. submitted linewidth studies, ms. in prep.

$\text{CH}_3\text{N}:$			
$\text{CH}_3\text{NH}_2$	methylamine	Jäger in coop. with Kreglewski, Poznan	incl. deuterated species hfs- and rot.-vib.- interaction
CNCN	cyanogen isocyanide	Gerry in coop. with Stroh and Winnewisser, Giessen	hfs complete, isotopes
HCCBr DCCBr	bromoacetylene	Harder Heineking	1-type spectra: Br-hfs D-hfs: published Stark effect: ms. in prep.
HCCl DCCl	chloroacetylene	Heineking	1-type spectra: Cl-hfs Stark effect: ms. in prep.
HCCF DCCF	fluoroacetylene	Haekel Heineking	relaxation studies D-hfs: published
HCCI DCCI	iodoacetylene	Andresen Heineking	structure: ms. submitted D-hfs: published Stark effect: ms. in prep.
$\text{C}_2\text{H}_2\text{O}$		Zerbe-Foese Hinze	dipole moments of excited states
$\text{C}_2\text{D}_2\text{O}$		Doose	rotational constants of excited states
$\text{C}_2\text{H}_2\text{F}^{14}\text{N}$ $\text{C}_2\text{H}_2\text{F}^{18}\text{N}$		Zerbe-Foese Hinze	dipole moments of excited states
$\text{C}_2\text{H}_3\text{NO}_2$	nitroethylene	Spieckermann	rotational Zeeman effect in progress
$\text{C}_2\text{H}_4\text{FNO}:$ $\text{CH}_2\text{FCONH}_2$	2-fluoroacetamide	Heineking	N-hfs complete
$\text{C}_2\text{H}_4\text{F}_3\text{N}:$ $\text{CF}_3\text{CH}_2\text{NH}_2$	trifluoroethylamine	Keussen	N-hfs complete N-d <sub>2</sub> -species in progress
$\text{C}_2\text{H}_5\text{I}$	ethyl iodide	Gripp	published
$\text{C}_2\text{H}_5\text{N}:$ $\text{CH}_3\text{CHNH}$	ethanimine	Krause	rotational Zeeman effect in progress
$\text{C}_2\text{H}_6\text{NO}:$ $\text{CH}_3\text{CONH}_2$	acetamide	Heineking	N-hfs complete (groundstate only)
$\text{C}_2\text{H}_6\text{ONO}$	ethyl nitrite	Keussen	N-hfs published inversion splittings
$\text{C}_2\text{H}_6\text{NO}:$ $\text{CH}_2\text{OHCH}_2\text{NH}_2$	aminoethanol	Keussen	N-hfs in progress

$C_2H_6N_2O:$ $(CH_3)_2NNO$	dimethyl nitrosamine	Heineking	N-hfs complete int.rot. in progress
$C_2H_6O$ $C_2D_6O$	dimethylether dimethylether-d <sub>6</sub>	Neustock	mmw spectrum
HCCON	cyanoacetylene	Haekel	relaxation studies
$C_3H_2ClN:$ $CH_2=CClCN$	chloroacrylonitrile	Vormann Heineking	Cl,N-hfs in progress
$C_3H_4:$ $CH_2=C=CD_2$ $CH_2=C=CHD$ $CH_2D-C\equiv CH$	1,1-dideuteroallene monodeuteroallene 3-deuteropropyne	Meyer	rotational Zeeman effect in progress
$C_3H_3DN_2$	1-/4-deuteropyrazole	Böttcher	rotational Zeeman effect in progress
$C_3H_3Br$	cyclopropyl bromide	Heineking in coop. with Li and Lewis-Bevan, Vancouver	Br-hfs complete
$C_3H_3I:$ $CH_2=CI-CH_3$	2-iodopropene	Gripp	int.rot. analysis in progress
$C_3H_5N:$ $C_2H_5NC$	ethyl isocyanide	Krüger	N-hfs reanalysed int.rot. in progress
$CH_2CHCHNH$	syn-/anti-propenimine	Krause	rotational Zeeman effect in progress
$C_3H_7Br:$ $(CH_3)_2CHBr$	isopropyl bromide	Meyer	Br-hfs, dipole moment, excited states in prog.
$C_3H_7Cl:$ $(CH_3)_2CHCl$	isopropyl chloride	Meyer	Cl-hfs, dipole moment, excited states in prog.
$C_3H_7F:$ $(CH_3)_2CHF$	isopropyl fluoride	Meyer	excited states complete
$C_3H_7I:$ $(CH_3)_2CHI$	isopropyl iodide	Gripp	I-hfs nearly complete
$C_3H_7N:$ $C_3H_5NH_2$	cyclopropylamine	Böttcher	rotational Zeeman effect in progress
$C_3H_7NO:$ $(CH_3)_2NCHO$	dimethyl formamide	Heineking	N-hfs, int.rot. in progress

$C_3H_7N:$			
$(CH_3)_2CHNH_2$	isopropylamine	Keussen	incl. deuterated species N-hfs in progress
$C_4H_6N:$			
$C_3H_5CN$	cyclopropyl cyanide	Böttcher	rotational Zeeman effect in progress
$C_4H_5NO:$			
$C_3H_5NCO$	cyclopropyl isocyanate	Heldmann	N-hfs in progress
$C_4H_5NS:$			
$CH_3C_3H_2NS$	4-/5-methylthiazole	Jäger	published
$C_4H_7N:$			
$C_3H_7NC$	n-propyl isocyanide	Krüger	N-hfs, int. rot. in progress
$(CH_3)_2CHNC$	isopropyl isocyanide	Krüger	assignment in progress
$C_4H_5N_2O:$			
$\underline{(CH_2)_4NNO}$	N-nitrosopyrrolidine	Heineking in coop. with Bohn, Storrs, Conn.	N-hfs complete
$C_4H_7Br:$			
$(CH_3)_3CBr$	tert.-butyl bromide	Harder	Br-hfs complete
$C_4H_7N:$			
$\underline{(CH_2)_4NH}$	pyrrolidine	Grabow Ehrlichmann	N,D-hfs complete
$C_4H_7NO:$			
$\underline{(CH_2)_2O(CH_2)_2NH}$	morpholine	Grabow Ehrlichmann	N,D-hfs complete
$C_4H_{10}$	butane	Vormann	int. rot. in progress (torsional groundstate)
$C_4H_5NO:$			
$C_4H_5O(CN)$	2-cyanofuran	Albinus	rotational Zeeman effect in progress
$C_5H_5NO:$			
$C_4H_4NCHO$	2-formyl pyrrole	Heineking	N-hfs complete
$C_5H_5NO$	pyridine-N-oxide	Heineking	N-hfs complete
$C_5H_{10}Si:$			
$(CH_3)_3SiCCH$	trimethylsilylacetylene	Ellendt	assignment in progress
$C_8H_{11}N:$			
$\underline{(CH_2)_5NH}$	piperidine	Grabow Ehrlichmann	N,D-hfs complete
$C_6H_4Cl_2$	1,2-dichlorobenzene 1,3-dichlorobenzene	Merke Keussen in coop. with Onda, Tokio	Cl-hfs in progress
$C_6H_5Cl$	chlorobenzene	Heineking	Cl-hfs complete

C <sub>6</sub> H <sub>5</sub> BF <sub>2</sub>	phenyldifluoroborane	Vormann	B-hfs published
C <sub>6</sub> H <sub>5</sub> NO:			
C <sub>6</sub> H <sub>4</sub> NCHO	3-/4-formyl pyridine	Stryjewski	N-hfs published
C <sub>6</sub> H <sub>5</sub> NO <sub>2</sub>	nitrobenzene	Heineking	v=0,1,2 N-hfs complete
C <sub>6</sub> H <sub>5</sub> NO <sub>3</sub> :			
C <sub>6</sub> H <sub>4</sub> (OH)NO <sub>2</sub>	o-nitrophenol	Heineking	N-hfs complete
C <sub>6</sub> H <sub>5</sub> D	deuterobenzene	Albinus	rotational Zeeman effect in progress
C <sub>6</sub> H <sub>7</sub> N:			
C <sub>6</sub> H <sub>5</sub> NH <sub>2</sub>	aniline	Oldag	rotational Zeeman effect in progress
C <sub>6</sub> H <sub>7</sub> NO:			
CH <sub>3</sub> C <sub>6</sub> H <sub>4</sub> NO	2-picolin-N-oxide	Heineking Kamura, Tokio	N-hfs, int.rot. in progress
FCIO <sub>3</sub>	perchloryl fluoride	Heldmann	hfs in progress
ClGeH <sub>3</sub>	germyl chloride	Heineking Wlodarczak, Lille	isotopes, Cl-hfs complete in coop. with <sup>73</sup> Ge-hfs in progress
Cl <sub>2</sub> S	sulfur dichloride	Merke	Cl-hfs in progress
Cl <sub>2</sub> SO	thionyl chloride	Merke	Cl-hfs in progress
Cl <sub>2</sub> SO <sub>2</sub>	sulfuryl chloride	Merke	Cl-hfs complete
D <sub>2</sub> O <sub>2</sub>	hydrogen peroxide-d <sub>2</sub>	Doose	rotational constants of excited states
FSiH <sub>3</sub>	silyl fluoride	Guarnieri Neustock	mmw spectrum
HN <sub>3</sub>	hydrazoic acid	Gerry Heineking	incl. <sup>15</sup> N-species N-hfs complete
NH <sub>3</sub>	ammonia	Wolf	collision-induced transitions
SiH <sub>4</sub>	silane	Stahl	hfs, ms. in prep.
N <sub>2</sub> O	nitrous oxide	Nicolaisen	1-type spectrum incl. <sup>15</sup> N-species relaxation studies

14 Name to whom queries should be addressed James R. Durig

Mailing address Department of Chemistry  
University of South Carolina  
Columbia, SC 29208

Telephone number: (803) 777-6612 Telefax: (803) 777-9521Electronic mail (BITNET or other): N670007@UNIVSCVM

FORMULA	NAME OF COMPOUND	NAME OF INVESTIGATOR	PRESENT STAGE OF PROGRESS
$C_2H_7P(CH_3CH_2PH_2)$	ethylphosphine	Groner	excited states in progress
$C_2H_5F_2P(CH_3CH_2PF_2)$	ethyldifluorophosphine	Groner	$d_1, ^{13}C$ partially assigned
$C_2H_5F_2OP(CH_3CH_2POF_2)$	ethylphosphonic difluoride	Johnson	<u>trans</u> assigned
$C_3H_7F_2P((CH_3)_2CHPF_2)$	isopropylidifluorophosphine	Groner	J. Phys. Chem. <u>93(7)</u> , (1989)
$C_2H_6F_2NP((CH_3)_2NPF_2)$	dimethylamino difluorophosphine	Harlan	J. Phys. Chem. <u>93(6)</u> , (1989)
$C_2H_6FPS((CH_3)_2PSF)$	dimethylphosphonothioic fluoride	Chatterjee	manuscript in preparation
$C_3H_9PS((CH_3)_3PS)$	trimethylphosphine sulfide	Chatterjee	Inorg. Chem. <u>28</u> , 298 (1989)
$C_4H_9F((CH_3)_2CHCH_2F)$	2-methyl-1-fluoropropane	Godbey	submitted, J. Chem. Phys.
$C_4H_7F(CH_2=C(CH_3)CH_2F)$	3-fluoro-2-methylpropene	Qiu	submitted, J. Chem. Phys.
$C_4H_7F(trans-CH_3CH=CHCH_2F)$	<u>trans</u> -1-fluoro-2-butene	Groner	<u>gauche</u> assigned, <u>cis</u> in progress
$C_5H_7ClO(CH_2CH_2CH_2CHCClO)$	cyclobutylcarbonyl chloride	Badawi	J. Mol. Struct. <u>190</u> , 475 (1988)
$C_5H_8O(CH_2CH_2CH_2CHCHO)$	cyclobutylcarboxaldehyde	Badawi	submitted, J. Chem. Phys.
$CHF_2N(CF_2=NH)$	difluoromethanimine	Groner	J. Chem. Phys. <u>89</u> , 3983 (1988); $d_1$ in progress
$CBrF_2N(CF_2=NBr)$	N-bromo-difluoromethanimine	Groner	manuscript in preparation
$GeH_3N_3(GeH_3NNN)$	germylazide	Sullivan, Groner	submitted, J. Chem. Phys.
$C_2H_2BrFO(FCH_2CBrO)$	fluoroacetyl bromide	Little	in progress
$C_4H_8O((CH_3)_2CHCHO)$	2-methylpropanal	Guirgis, Stiefvater	submitted, J. Chem. Phys.
$C_4H_{10}Ge(CH_2CH_2CH_2CHGeH_3)$	cyclobutylgermane	Geyer	Chem. Phys. <u>125</u> , 299 (1988)
$C_2H_3FO_2(CH_3OCFO)$	methylfluoroformate	Tolley	$d_1, d_3$ assigned, $^{18}O$ in progress
$C_3H_9P((CH_3)_2CHPH_2)$	isopropylphosphine	Barron	search for 2nd conformer
$C_3H_7F((CH_3)_2CHF)$	isopropylfluoride	Nanaie	manuscript in preparation
$C_5H_9Cl(CH_2(CH_2)_3CHCl)$	chlorocyclopentane	Lee	equatorial conformer assigned
$C_2H_3ClO_2(CH_3OCClO)$	methylchloroformate	Tolley	$^{18}O$ assigned; manuscript in preparation
$C_3H_9P((CH_3)_3P)$	trimethylphosphine	Chatterjee	submitted, Struct. Chem.

15 Name to whom queries should be addressed P.G.FAVEROMailing address Dipartimento di Chimica "G.Ciamician", Universita' di Bologna,  
Via F.Belmi 2, I-40126 Bologna (Italy).Telephone: 0039-51-259500/1 Telefax: 0039-51-259456Electronic mail (BITNET or other): T18101 at ICINECA1 (EARN)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
COS(OCS)	carbonyl sulfide	P.G.Favero <sup>a</sup>	New results on the N <sub>2</sub> <sup>*</sup> -OCS system.
CClN(ClCN)	cyanogen chloride	P.G.Favero <sup>a</sup>	N <sub>2</sub> <sup>*</sup> -ClCN system; work in progress.
COS(OCS)	carbonyl sulfide	G.Cazzoli <sup>a</sup>	Modulated line shape analysis.
CHO <sup>+</sup> (HCO <sup>+</sup> )	formyl ion	G.Cazzoli <sup>a</sup>	Preliminary results on pressure broadening.
CDO <sup>+</sup> (DCO <sup>+</sup> )	formyl ion	G.Cazzoli <sup>a</sup>	HFS and line shape analysis.
BF	boron fluoride	G.Cazzoli <sup>a</sup>	In press, J.Mol.Spectr.
C <sub>6</sub> H <sub>8</sub> O	3-cyclohexenone	D.Damiani <sup>a,b</sup>	In press, J.Mol.Spectr.
C <sub>7</sub> H <sub>10</sub>	3-cyclohexene-1-methylene	D.Damiani <sup>a,b</sup>	Submitted.
C <sub>4</sub> H <sub>8</sub> O	cis-crotyl alcohol	W.Caminati <sup>c</sup>	Submitted, J.Phys.Chem.
C <sub>2</sub> H <sub>7</sub> NS	2-amino-ethane thiol	W.Caminati <sup>c</sup>	Trideuterated species. In press, J.Mol.Str.
C <sub>3</sub> H <sub>6</sub> O <sub>3</sub>	methylglycolate	W.Caminati <sup>d</sup>	Torsional interactions. Submitted.
C <sub>3</sub> H <sub>3</sub> NS	vinylisothiocyanate	W.Caminati	Syn conformer + exc. states.
C <sub>6</sub> H <sub>8</sub> Ge	phenyl germane	W.Caminati	Structure, V <sub>6</sub> barrier.
C <sub>4</sub> H <sub>7</sub> N	3-pyrroline	W.Caminati <sup>d</sup>	Ring puckering, H <sub>N</sub> inversion.
C <sub>2</sub> H <sub>5</sub> NO <sub>2</sub>	N-methoxyformamide <i>(C-N=OCH<sub>3</sub>)</i>	W.Caminati <sup>d</sup>	Manuscript in preparation.

<sup>a</sup> In collaboration with the Istituto di Spettroscopia Molecolare del CNR, Bologna.<sup>b</sup> In collaboration with the Istituto di Chimica Industriale dell' Universita', Messina.<sup>c</sup> In collaboration with the Dipartimento di Chimica Fisica ed Inorganica, Bologna.<sup>d</sup> In collaboration with the Laboratorium fur Physikalische Chemie, ETH, Zurich.

Note: Only a reference participant is listed for each compound, even if most works are with multiple authorship.

16 Name to whom queries should be addressed JERZY GALICA

Mailing address Institute of Molecular Physics of the Polish Academy  
of Sciences  
ul.Smoluchowskiego 17/19, 60-179 Poznań, POLAND

Telephone: 67-40-71 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>5</sub> H <sub>9</sub> N	tertiary butyl cyanide	J.Galica S.Gierszal	linewidth studies in progress

17 Name to whom queries should be addressed Dr. M.C.L. Gerry

Mailing address Department of Chemistry, The University of British Columbia  
2036 Main Mall  
Vancouver, B.C., CANADA, V6T 1Y6

Telephone: (604) 228-2129 Telefax: (604) 228-2847

Electronic mail (BITNET or other): USERMCLG@UBCMTSG

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>2</sub> HBrO (BrHCCO)	bromoketene	N. Westwood W. Lewis-Bevan	Manuscript submitted
C <sub>3</sub> H <sub>5</sub> Br	bromocyclopropane	H. Li (with N. Heineking, H. Dreizler, Kiel)	High order quadrupole coupling analysis
CHNO	isocyanic acid	S. Howard, M. Gerry (with N. Heineking, Kiel)	<sup>17</sup> O, D species analyzed.
C <sub>3</sub> H <sub>3</sub> N(CH <sub>2</sub> CHCN)	acrylonitrile	E. Tien R. Richards	Isotopic substitutions Ms. in preparation
CH <sub>3</sub> NO(CH <sub>2</sub> NOH)	formaldoxime	N. Lee H. Jemson D. MacLennan	Isotopic substitutions. Spectra analysed
CClNS(CNCS)	chlorine thiocyanate	R. Richards	Isotopic substitutions
CH <sub>3</sub> N <sub>3</sub> HN <sub>3</sub>	methyl azide hydrazoic acid	M. Gerry N. Heineking	<sup>14</sup> N hyperfine analysis at U. of Kiel
C <sub>2</sub> N <sub>2</sub> (CNCN)	cyanogen isocyanide	M. Gerry (with M. Winnewisser F. Stroh, Giessen)	<sup>14</sup> N hyperfine analysis at U. of Kiel

18 Name to whom queries should be addressed Charles W. Gillies

Mailing address Department of Chemistry  
Rensselaer Polytechnic Institute  
Troy, New York 12180

Telephone: 518-276-8453 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2HF_3O(CHFCF_2O)$	Trifluoroethylene Oxide	T. Raw	J. Mol. Spectrosc. <u>128</u> , 195 (1988).
$C_2H_4O_3(CH_2CH_2O_00)$	1,2,3-Trioxolane	J. Z. Gillies R. Suenram* F. Lovas*	J. Am. Chem. Soc. <u>110</u> , 7991 (1988).
$CH_3ArNO$	Argon-formamide complex	R. D. Suenram* G. T. Fraser* F. J. Lovas* J. Zozom	J. Chem. Phys. <u>89</u> , 6141 (1988).
$CH_4ArO$	Argon-methanol complex	R. D. Suenram* G. T. Fraser* F. J. Lovas* J. Z. Gillies	Manuscript in preparation.
$C_2H_4O_3$	Ethylene-ozone complex	J. Z. Gillies R. D. Suenram* F. J. Lovas* V. Stahl**	J. Am. Chem. Soc., comm. in press (1989).
$H_2O_4$	Ozone-water complex	J. Z. Gillies R. D. Suenram* F. J. Lovas*	Manuscript in preparation.
$C_2H_2O_3$	Acetylene-ozone complex	J. Z. Gillies F. J. Lovas* K. Matsumura* R. D. Suenram*	Normal, $d_1$ and $d_2$ isotopes assigned
$CH_3F_2O_2P(CH_3OP(O)F_2)$	Methyl Difluorophosphate	L. Rickus H. Justnes J. Zozom R. Suenram* F. Lovas*	Manuscript in preparation.

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>3</sub> H <sub>2</sub> F <sub>4</sub> (CF <sub>2</sub> CHFCHF)	cis-1,2,3,3-Tetra-fluorocyclopropane	R. Beauchamp J. Gillies	Manuscript in preparation

\*National Institute of Standards and Technology

\*\*Institut fur Physikalische Chemie

19 Name to whom queries should be addressed Jack D. Graybeal

Mailing address Department of Chemistry

Davidson Hall, Va. Polytechnic Inst. and State Univ.

Blacksburg, VA 24061-0212

Telephone: (703) 231-5997 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2F_6NO((CF_3)_2NO)$	bis-trifluoromethyl-nitroxide radical	R. M. White	Manuscript in preparation
$C_4H_6N((CH_3)_2CCN)$	2-cyano-2-propyl radical	R. C. Claytor	Manuscript in preparation

20 Name to whom queries should be addressed H. S. Gutowsky

Mailing address University of Illinois

177 Noyes Lab, Box 25

505 S. Mathews Ave., Urbana, IL 61801

Telephone: 217/333-7621 Telefax:

Electronic mail (BITNET or other): Bitnet%"Gutowsky@UIUCSCS"

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR*</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2HNO_2$	OCO-HCN linear dimer	T. D. Klots	Manuscript accepted by JCP
$C_3H_3Si$ [( $CH_3$ ) <sub>2</sub> Si=CH <sub>2</sub> ]	dimethylsilaethylene	H. S. Gutowsky T. Emilsson	Commun. accepted by JACS; detailed analysis in progress
$Ar_2ClD$ & $Ar_3ClD$	$Ar_2$ -DCl trimer & $Ar_3$ -DCl tetramer	T. D. Klots	Manuscript submitted to JCP
$C_3HNO_4$	$H\bar{C}N-(CO_2)_2$ trimer	Jane Chen	Observation and analysis of spectrum in progress
$C_4HNO_6$	$HCN-(CO_2)_3$ tetramer	P. J. Hajduk	Observation and analysis of spectrum in progress
$CF_2$ & $C_2F_4$ [ $CF_2$ & $CF_3CF$ ]	difluorocarbene and perfluoromethyl carbene	Carl Chuang	Observation and analysis of spectra in progress
$B_2ClH_7$ [ $B_2H_6$ -HCl]	bent diborane-HCl dimer	Carl Chuang	Observation and analysis of spectra in progress
$CFH_5N_2$	$NH_3$ -HCN-HF trimer	T. D. Klots	Manuscript in progress
$C_2FH_2NO$	OC-HCN-HF trimer	T. D. Klots	Manuscript in progress
$C_2H_2N_4$ $C_3H_2N_2O$ $C_2H_5N_3$ $C_2H_4N_2O$	$N_2$ -HCN-HCN trimer OC-HCN-HCN trimer $NH_3$ -HCN-HCN trimer $H_2O$ -HCN-HCN trimer	R. Ruoff T. Emilsson	Manuscript in progress

\*Most of this work is a group effort with multiple authorship. Only one or two of several participants are listed for each study.

21 Name to whom queries should be addressed Marlin D. Harmony

Mailing address Department of Chemistry

University of Kansas

Lawrence, KS 66045

Telephone: (913)864-3980 (or 4670) Telefax: (913)842-5612

Electronic mail (BITNET or other): CHEMISTS AT UKANVM

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_3F_2O$	Difluoropropadienone	Tam	Assignment Underway
$C_5H_5N(C_4H_5CN)$	Cyanobicyclobutane	Taylor, Harmony	Manuscript being finalized

22 Name to whom queries should be addressed Dr. Dieter Hausmann

Mailing address Institute of Optoelectronics

German Aerospace Research Est.

8031 Oberpfaffenhofen, FRG.

Telephone: 8153-28770 Telefax: 8153-28243

Electronic mail (BITNET or other): OE1H at DFVLR OP1

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CHF <sub>3</sub>	Fluoroform	D. Hausmann	FIR spectrum, new lines assigned

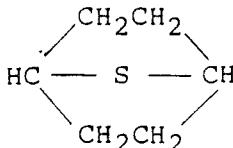
23 Name to whom queries should be addressed Kichiro Hayashi  
 Mailing address Department of Chemistry, Faculty of Science  
Hiroshima University  
Higashi-sendamachi, Naka-ku, Hiroshima 730, Japan  
 Telephone: 082-241-1221, ext. 2388 fax:  
 Electronic mail (BITNET or other):

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_4H_8O(CH_3CH_2OCH=CH_2)$	ethyl vinyl ether	Hayashi, Kakumoto	$r_s$ structure
$C_2H_5F$	ethyl fluoride	Hayashi, Miyazaki	Fermi resonance
$CH_5SiBr$	methyl bromosilane	Hayashi, Ikeda	$r_s$ structure
$C_3H_{10}Si(CH_3CH_2SiH_2CH_3)$	ethyl methyl ether	Hayashi, Ikeda	$r_s$ structure
$C_3H_{10}Si(CH_3CH_2CH_2SiH_3)$	propyl silane	Hayashi, Ikeda	$r_s$ structure
$C_3H_7I$	propyl iodide	Hayashi	$eQq$ , $r_s$ structure
$C_3H_8O$	ethyl methyl ether	Hayashi	excited states

24 Name to whom queries should be addressed Eizi Hirota

Mailing address Institute for Molecular Science  
Myodaiji, Okazaki 444  
Japan

Telephone: 81 (Japan) -564-54-1111, ext. 3320;  
81 (Japan) -564-53-7322 Telefax: 81 (Japan) -564-54-2254  
 (direct)  
 Electronic mail (BITNET or other):

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_5H_8$	1,4-pentadiene <chem>[CH2=CHCH2CH=CH2]</chem>	E. Hirota	One rotamer assigned.
$C_6H_{10}S$	7-thiabicyclo [2.2.1]heptane	K. Irie	Work in progress.
			
HNO	nitroxyl	K. Takagi S. Saito	Manuscript in preparation.
$C_1OS[C_1SO]$	$C_1SO$ radical	S. Saito	Work almost completed.
$CD_3O$	methoxy radical-d <sub>3</sub>	Y. Endo	Manuscript in preparation.
$CH_3CO$	acetyl radical	Y. Endo	Work in progress.
$F_3Si[SiF_3]$	silicon trifluoride	M. Tanimoto	Assigned.
$C_2D[CCH]$	ethynyl radical	Y. Endo	Manuscript in preparation.
$C_4H_6D_2$	cyclobutane-1,2-d <sub>2</sub>	E. Hirota	cis, work in progress.
$C_4H_4D_4$	cyclobutane-1,1,3,3-d <sub>4</sub>	E. Hirota	Work almost completed.

NaO	sodium monoxide	C. Yamada	Paper accepted in J.Chem.Phys.
LiO	lithium monoxide	C. Yamada	Paper accepted in J.Chem.Phys.
KO	potassium monoxide	C. Yamada	$2\Pi, 2\Sigma^+$ assigned.
RbO	rubidium monoxide	C. Yamada	Assigned.
AlO	aluminum monoxide	C. Yamada	Assigned.
Cl <sub>2</sub> Si SiCl <sub>2</sub>	silicon dichloride (dichlorosilylene)	M. Fujitake	Work almost completed.
ClFSi SiFCl	silicon chloride fluoride (chlorofluorosilylene)	M. Fujitake	Assigned.
CCl <sub>2</sub>	dichlorocarbene	M. Fujitake	Assigned.

25 Name to whom queries should be addressed Prof. Dr. J. Hoeft

Mailing address Institut für Molekulphysik, Freie Universität Berlin,  
Arnimallee 14, D-1000 Berlin 33, Fed. Rep. Germany

Telephone: (030) 838 3590 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
Y <sub>0</sub>	Yttrium monoxide	Törring, Fehlhammer	rot. spectrum, Hfs assigned
InCl	Indium monochloride	Hoeft, K.P.R. Nair	Chem.Phys.Lett. (in press)
InBr	Indium monobromide	Hoeft, K.P.R. Nair	manuscript in preparation

26 Name to whom all queries should be addressed H.W.Kroto

Mailing address School of Chemistry and Molecular Sciences  
University of Sussex  
Brighton BN1 9QJ UK

Telephone number: 0273-606755

<u>Formula</u>	<u>Name of Compound</u>	<u>Name of Investigator</u>	<u>Present Stage of Progress</u>
C <sub>2</sub> H <sub>5</sub> NO CH <sub>3</sub> CH <sub>2</sub> NO	Nitrosoethane	M Maier D Silverton	See also* A P Cox
C <sub>5</sub> H <sub>10</sub> Si (CH <sub>3</sub> ) <sub>3</sub> SiC≡CH	Trimethylsilyl -ethyne	A Alexander D R M Walton	ms in prep
C <sub>6</sub> H <sub>9</sub> NSi (CH <sub>3</sub> ) <sub>3</sub> SiC≡CCN	Trimethylsilyl -cyanoethyne	"	"
C <sub>8</sub> H <sub>9</sub> NSi (CH <sub>3</sub> ) <sub>3</sub> Si(C≡C) <sub>2</sub> CN	Trimethylsilylcyanato -cyanoethyne	"	"*
CF <sub>2</sub> HP CF <sub>2</sub> =PH	2,2-Difluorophospha -ethene	J F Nixon N P C Simmons	ms in prep
BrCH <sub>2</sub> P CH <sub>2</sub> =PBr	1-Bromophospha -ethene	J F Nixon O Ohashi	ms in prep*
CFH <sub>2</sub> P CH <sub>2</sub> =PF	1-Fluorophospha -ethene	J F Nixon O Ohashi D R M Walton	submitted
C <sub>2</sub> F <sub>3</sub> P CF <sub>3</sub> C≡P	3,3,3-Trifluoro-1-phosphapropyne	J F Nixon N P C Simmons	ms in prep*
C <sub>2</sub> NP N≡C-C≡P	C-Cyanophospha -ethyne	O Ohashi J F Nixon	ms in prep*
BFS FB=S	Fluorosulphido -boron	T Cooper	ms in prep*
BBrS BrB=S	Bromosulphido -boron	T Cooper	ms in prep*
BClSe ClB=S	Chloroselenido -boron	M King	in prog*
C <sub>3</sub> HNO HC≡CNCO	Isocyanatoethyne	T Cooper D R M Walton	ms in prep
C <sub>4</sub> H <sub>3</sub> NO CH <sub>3</sub> C≡CNCO	Isocyanatopropyne	S Aziz M Durrant D R M Walton	ceased
AlC <sub>3</sub> H <sub>12</sub> N (CH <sub>3</sub> ) <sub>3</sub> NaH <sub>3</sub>	Trimethylaminoalane	C Kirby D Smith	Broadband

<u>Formula</u>		<u>Name of Compound</u>	<u>Name of Investigator</u>	<u>Present Stage of Progress</u>
C <sub>5</sub> H <sub>3</sub> N	CH <sub>2</sub> =CHC≡CCN	1-Cyano-but-3-en-1-yne	K Phillips D R M Walton D McNaughton	submitted
C <sub>5</sub> H <sub>3</sub> N	HC≡CCH=CHCN	1-Cyano but-1-en-3-yne	J August D McNaughton	submitted
C <sub>4</sub> H <sub>3</sub> N	HC≡CCH <sub>2</sub> NC	1-isocyano prop-2-yne	N Romeril M F Lappert D McNaughton	ms complete
C <sub>3</sub> H <sub>3</sub> N	HC≡CCH=NH	C-ethynyl methanimine	D McNaughton O Osman	submitted
C <sub>5</sub> H <sub>9</sub> N	CH <sub>3</sub> CH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> CN	Valeronitrile	J August K Georgiou A C Legon	in prog.

\* some preliminary data in Chem.Soc.Revs., 11, 435 (1982)

27 Name to whom queries should be addressed Robert L. Kuczkowski, Kurt W. Hillig, II

Mailing address Department of Chemistry

The University of Michigan

Ann Arbor, MI 48109-1055

Telephone: RLK: (313) 764-7540 / KWH: (313) 747-2867 Telefax: (313) 747-4865

Electronic Mail:

RLK: USERGBRK@UMICHUB.BITNET (BITNET) or Robert\_Kuczkowski@UB.CC.UMICH.EDU (INTERNET)  
 KWH: USERK9F3@UMICHUM.BITNET (BITNET) or Kurt\_Hillig@UM.CC.UMICH.EDU (INTERNET)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>4</sub> H <sub>4</sub> O <sub>3</sub>	cyclobutadiene ozonide	Lorenzak	In press, JPC
C <sub>4</sub> H <sub>8</sub> O	1-butene oxide	Badawi	2 Isotopes
C <sub>7</sub> H <sub>12</sub> O	8-oxabicyclo[3.2.1]octane	Badawi, Lorenzak, H Wieser (Calgary)	Assigned
C <sub>4</sub> H <sub>3</sub> F <sub>3</sub>	2,3,3-trifluorocyclobutene	Lorenzak, Sibley, Craig (Oberlin)	Assigned
C <sub>3</sub> H <sub>9</sub> BF <sub>3</sub> N	trimethylamine-BF <sub>3</sub>	Riggs, LaBarge, Dreizler (Kiel)	eQq analysis
C <sub>5</sub> H <sub>10</sub> O	pivaldehyde	Cox (Bristol)	<sup>13</sup> C Isotopes
C <sub>4</sub> H <sub>4</sub> O·Ar	argon-furan complex	Oh	Dipole and CD analysis
C <sub>4</sub> H <sub>5</sub> N·Ar	argon-pyrrole complex	Oh	Dipole and CD analysis
Ar·F <sub>3</sub> P	argon-PF <sub>3</sub> complex	LaBarge, Bohn, Taleb-Bendiab	Dipole, CD, etc. In press, JCP
F <sub>3</sub> KrP	krypton-PF <sub>3</sub> complex	LaBarge, Bohn, Taleb-Bendiab	Dipole, eQq, CD, etc. In press, JCP
F <sub>3</sub> NeP	neon-PF <sub>3</sub> complex	Hillig	Assigned
H <sub>2</sub> F <sub>3</sub> OP	water-PF <sub>3</sub> complex	LaBarge, Hillig	Assigned
C <sub>3</sub> H <sub>9</sub> NO <sub>2</sub> S	trimethylamine-SO <sub>2</sub>	Oh	Isotopes
C <sub>4</sub> H <sub>4</sub> NO <sub>2</sub> S	pyridine-SO <sub>2</sub> complex	Oh	Assigned
C <sub>2</sub> H <sub>4</sub> O <sub>2</sub> S	ethylene-SO <sub>2</sub> complex	Andrews, LaBarge	Isotopes
C <sub>6</sub> H <sub>6</sub> O <sub>2</sub> S	benzene-SO <sub>2</sub> complex	Taleb-Bendiab, LaBarge	Assigned

28 Name to whom queries should be addressed Dr. Stephen Kukolich

Mailing address University of Arizona

Department of Chemistry

Tucson, AZ 85721

Telephone: (602)621-2969 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CH <sub>2</sub> O <sub>2</sub>	Formic Acid (IR)	Bumgarner, Choe	MS Accepted
DFN <sub>2</sub> O	N <sub>2</sub> O - DF Complexes	D. Pauley	MS Accepted
CO	Co - Intercomb. Bands	Choe, Lee, LeFlock	MS Accepted

29 Name to whom queries should be addressed \_\_\_\_\_ R.M. Lees or K.V.L.N. Sastry  
 Mailing address \_\_\_\_\_ Department of Physics  
 \_\_\_\_\_ University of New Brunswick  
 \_\_\_\_\_ Fredericton, N.B., Canada E3B 5A3  
 Telephone: (506)-453-4723 Telefax: \_\_\_\_\_  
 Electronic mail (BITNET or other): LEES@UNB

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CH <sub>4</sub> O (CD <sub>3</sub> OH)	Methanol	R.M. Lees M. Mollabashi (with I. Mukhopadhyay, UBC and J. Johns, Herzberg Inst.)	Fourier transform FIR and laser assignments
CD <sub>3</sub> OH, <sup>13</sup> CH <sub>3</sub> OH, <sup>13</sup> CD <sub>3</sub> OH	"	K.V.L.N. Sastry J. VanderLinde R.M. Lees	Dipole moment measurements
CH <sub>3</sub> <sup>18</sup> OH	"	R.M. Lees (with W. Lewis-Bevan, Southern Illinois U.)	FTIR and FTFIR and FIR laser assignments
CH <sub>5</sub> N (CH <sub>3</sub> NH <sub>2</sub> )	Methylamine	R.M. Lees K.V.L.N. Sastry (with W. Lewis-Bevan)	Spectrum assigned FIR laser assignments

30 Name to whom queries should be addressed A.C. Legon

Mailing address Department of Chemistry, University of Exeter,  
Stocker Road, Exeter EX4 4QD, United Kingdom.

Telephone: 0392 - 263488 Telefax: 0392 - 263108

Electronic mail (BITNET or other): LEGON.AC@UK.AC.EXETER (onJANET)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
ClH <sub>3</sub> O	H <sub>2</sub> O•••HCl	L.C. Willoughby	Manuscript in preparation
BrH <sub>3</sub> O	H <sub>2</sub> O•••HBr	A.P. Suckley	Preliminary publication: <i>Chem. Phys. Letters</i> , <b>150</b> , 153, (1988).
ClH <sub>4</sub> N	H <sub>3</sub> N•••HCl	N.W. Howard	Paper published: <i>J. Chem. Phys.</i> , <b>88</b> , 4694, (1988)
ClH <sub>4</sub> P	H <sub>3</sub> P•••HCl	L.C. Willoughby	Manuscript in preparation.
H <sub>4</sub> IP	H <sub>3</sub> P•••HI	N.W. Howard	Spectrum assigned.
H <sub>3</sub> IS	H <sub>2</sub> S•••HI	A.P. Suckley	Spectrum assigned.
BrHN <sub>2</sub>	N <sub>2</sub> •••HBr	N.W. Howard	Paper published: <i>J. Chem. Phys.</i> <b>90</b> , 672, (1989).
ClHO <sub>2</sub> S	SO <sub>2</sub> •••HCl	A.J. Travis	Further work.
CHArF <sub>3</sub>	Ar•••HCF <sub>3</sub>	E.J. Goodwin	Spectrum assigned.
C <sub>2</sub> H <sub>5</sub> N	CH <sub>4</sub> •••HCN	A.L. Wallwork	Assigned. Isotopic work in progress
C <sub>2</sub> H <sub>2</sub> O <sub>2</sub> S	(SO <sub>2</sub> ,HCCH)	N.W. Howard	Spectrum observed.
C <sub>2</sub> O <sub>3</sub>	OCO•••CO	A.P. Suckley	IR and MW spectra assign. Isotopic work in progress.
C <sub>2</sub> H <sub>3</sub> F <sub>3</sub>	CH <sub>2</sub> CF <sub>2</sub> •••HF	Z. Kisiel	Spectrum assigned.
C <sub>2</sub> H <sub>4</sub> F <sub>2</sub>	CH <sub>2</sub> CHF•••HF	Z. Kisiel	Search in progress.
C <sub>2</sub> H <sub>3</sub> ArF	CH <sub>2</sub> CHF•••Ar	Z. Kisiel	Spectrum assigned.
C <sub>3</sub> H <sub>3</sub> O	(H <sub>2</sub> CO,HCCH)	N.W. Howard	Paper published: <i>J. Chem. Phys.</i> , <b>88</b> , 6793, (1988).

[Continued  
over page]

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>3</sub> H <sub>5</sub> Cl	H <sub>2</sub> C=C=CH <sub>2</sub> •••HCl	A.J. Travis	Spectrum assigned.
C <sub>3</sub> H <sub>10</sub> ClN	(CH <sub>3</sub> ) <sub>3</sub> N•••HCl	C.A. Rego	Preliminary publication: <i>Chem. Comm.</i> , 1496, (1988).
C <sub>3</sub> H <sub>10</sub> BrN	(CH <sub>3</sub> ) <sub>3</sub> N•••HBr	C.A. Rego/ A.L. Wallwork	Assigned. <sup>14</sup> N/ Br nuclear quadrupole hfs analysed.
C <sub>6</sub> H <sub>6</sub> ClN	CH <sub>3</sub> NH <sub>2</sub> •••HCl	C.A. Rego	Assigned. Isotopic work in progress.
C <sub>4</sub> H <sub>10</sub> N <sub>2</sub>	(CH <sub>3</sub> ) <sub>3</sub> N•••HCN	C.A. Rego	Paper published: <i>J. Chem. Phys.</i> , 89, 696, (1988).
C <sub>4</sub> H <sub>9</sub> NSi	(CH <sub>3</sub> ) <sub>3</sub> SiCN	C.S. Cooper	<sup>14</sup> N - nuclear quadrupole coupling.
C <sub>5</sub> H <sub>11</sub> N	(CH <sub>3</sub> ) <sub>3</sub> N•••HCCH	C.A. Rego	Paper published: <i>J. Mol. Structure</i> , 189, 137, (1988)
C <sub>5</sub> H <sub>10</sub> FN	(CH <sub>3</sub> ) <sub>3</sub> CCN•••HF	N. Gerry	Spectrum assigned. Manuscript in preparation.
C <sub>5</sub> H <sub>9</sub> N	(CH <sub>3</sub> ) <sub>3</sub> CNC	A.L. Wallwork	Heavy atom r <sub>g</sub> - geometry; paper in press: <i>J. Mol. Structure</i> .
C <sub>7</sub> H <sub>13</sub> N	1 - azabicyclo[2.2.2] octane	M.Ellis	<sup>14</sup> N-nuclear quadrupole coupling.
C <sub>10</sub> H <sub>15</sub> Cl	1 - chloroadamantane	M. Ellis	Cl - nuclear quadrupole coupling.

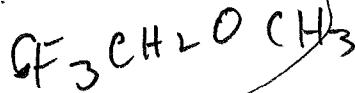
31 Name to whom queries should be addressed Dr. Ying S. Li

Mailing address Memphis State University  
Department of Chemistry  
Memphis, TN 38152

Telephone: 901-678-4427 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>9</sub> H <sub>14</sub> O	Bicyclo[3,3,1]nonan-9-one	San Li	Submitted
C <sub>3</sub> H <sub>5</sub> Cl <sub>3</sub> O	2,2,2-Trichloroethylmethyl ether	G. Lui J. Lu	Assigned
C <sub>8</sub> H <sub>14</sub> O	4-Ethylcyclohexanone	Arthur Lee	Assigned
C <sub>7</sub> H <sub>10</sub> O	3-Methyl-2-Cyclohexen-1-one	San Li	Assigned
CH <sub>3</sub> F <sub>2</sub> P	Difluoromethylphosphine	W. Wang R. A. Beaudet Y. S. Li	Submitted
C <sub>2</sub> H <sub>7</sub> SiI	Dimethylsilyliodide	P. Gironer J. R. Durig Y. S. Li	Quadrupole
C <sub>3</sub> H <sub>5</sub> F <sub>3</sub> O	2,2,2-Trifluoroethylmethyl ether	J. R. Durig Y. S. Li	Assigned



32 Name to whom queries should be addressed F. J. Lovas and R. D. Suenram

Mailing address National Institute of Standards and Technology  
Molecular Spectroscopy Division, Rm. B268/Bldg. 221  
Gaithersburg, MD 20899

Telephone: (301) 975-2385 Telefax: 301-869-7761  
or (301) 975-2165

Electronic mail (BITNET or other): LOVAS@NBS (BITNET)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
Ar-H <sub>2</sub> O	Argon water complex	R.D. Suenram F.J. Lovas	Four isotopic species assigned.
ArO <sub>2</sub> S (Ar-SO <sub>2</sub> )	Argon sulfur dioxide complex	L. Coudert K. Matsumura	New rotational measurements and tunneling analysis, Ar <sup>34</sup> SO <sub>2</sub> assigned.
CH <sub>2</sub> O <sub>2</sub> (CO-H <sub>2</sub> O)	Water carbon monoxide complex	D. Yaron <sup>a</sup> K.I. Peterson <sup>b</sup> D. Zolandz <sup>a</sup> W. Klemperer <sup>a</sup> R.D. Suenram F.J. Lovas	Normal, deuterated, <sup>17</sup> O, and <sup>18</sup> O isotopic species assigned. Tunneling motion and structure analysis, manuscript in prep.
CH <sub>2</sub> O <sub>2</sub> S (H <sub>2</sub> S-CO <sub>2</sub> )	Hydrogen sulfide carbon dioxide complex	J.K. Rice <sup>c</sup> F.J. Lovas R.D. Suenram L.H. Coudert K. Matsumura W. Stahld <sup>d</sup> D.J. Pauley <sup>e</sup> S. Kukolich <sup>e</sup>	Rotational spectrum for 3 isotopic species assigned. Tunneling motions and structure analyzed. Manuscript in prep.
C <sub>2</sub> H <sub>2</sub> O <sub>3</sub> (O <sub>3</sub> -HCCH)	Ozone-acetylene complex	J.Z. Gillies <sup>f</sup> C.W. Gillies <sup>f</sup> F.J. Lovas K. Matsumura R.D. Suenram	Rotational spectrum for 3 isotopic species assigned. Structure and dipole determined. Manuscript in prep.

<sup>a</sup>Department of Chemistry, Harvard University, Cambridge, MA.

<sup>b</sup>Department of Chemistry, University of Rhode Island, Kingston, RI.

<sup>c</sup>Naval Research Laboratory, Washington, D.C.

<sup>d</sup>Institute for Physical Chemistry, University of Kiel, Germany.

<sup>e</sup>Department of Chemistry, University of Arizona, Tucson, AZ.

<sup>f</sup>Department of Chemistry, Rensselaer Polytechnic Inst., Troy, NY.

<sup>g</sup>Department of Chemistry, Wesleyan University, Middletown, CT.

*No.  
Dr. Lovas and Suenram  
shuttle meteo - Red Line  
Shady Grove*

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2H_4O_2$ $((H_2CO)_2)$	Formaldehyde dimer	F.J. Lovas R.D. Suenram L.H. Coudert T.A. Blake <sup>g</sup> K.G. Grant <sup>g</sup> S.E. Novick <sup>g</sup>	Rotation-tunneling spectrum assigned for normal and perdeuterated species. Manuscript in prep.
$C_2H_4O_3$ $(O_3-H_2CCH_2)$	Ozone ethylene complex	J.Z. Gillies <sup>f</sup> C.W. Gillies <sup>f</sup> R.D. Suenram F.J. Lovas W. Stahl <sup>d</sup>	Rotational spectrum for 3 isotopic species assigned. Manuscript submitted to J. Am. Chem. Soc.
$C_2Si$ $(SiC_2)$	Silicon dicarbide	R.D. Suenram F.J. Lovas K. Matsumura	Produced by laser ablation, $J=1-0$ transition measured for $^{28}Si$ , $^{29}Si$ and $^{30}Si$ isotopes. Dipole determined.
$ClFH_2$ $(HF-HCl \text{ and } HCl-HF)$	Hydrogen fluoride Hydrogen chloride complex	G.T. Fraser A. Pine	$K=1$ spectrum of HF-HCl assigned, $K=0,1$ spectrum observed for HCl-HF. Manuscript submitted to JCP.
$H_2O_3S$ $(H_2O-SO_2)$	Water-sulfur dioxide complex	K. Matsumura F.J. Lovas R.D. Suenram	Tunneling rotation spectrum assigned for 3 isotopic species. Manuscript submitted JCP.
$H_2O_4$ $(O_3-H_2O)$	Ozone-water complex	J.Z. Gillies <sup>f</sup> C.W. Gillies <sup>f</sup> R.D. Suenram F.J. Lovas	Rotational spectrum for 3 isotopic species assigned. Manuscript in prep.
$H_4O_2$ $(H_2O)_2$	Water dimer	G.T. Fraser R.D. Suenram L.H. Coudert	$A_1$ state and c-type spectrum assigned. Manuscript submitted to JCP.
$(D_2O)_2$	Perdeutero-water dimer	R.D. Suenram G.T. Fraser F.J. Lovas	$A_2$ and $B_2$ states assigned and $K=1$ transitions for all states observed.

33 Name to whom queries should be addressed B. MACKE

Mailing address LABORATOIRE DE SPECTROSCOPIE HERTZIENNE  
S.F.R. DE PHYSIQUE - BATIMENT P.5  
UNIVERSITE DE LILLE 1  
59655 - VILLENEUVE D'ASCQ CEDEX FRANCE

Telephone: 20.43.47.84 Telefax: 20.43.49.95

Electronic mail (BITNET or other):

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
H <sub>2</sub> O	water vapor	( A. Bauer ( M. Godon ( J. Carlier	Lineshape
MgCl	magnesiumchloride	(	mm wave spectrum
C <sub>2</sub> H( <sup>13</sup> CCH, <sup>13</sup> CCH)	ethynyl radical	( M. Bogey ( H. Bolvin ( C. Demuynck	" " "
N(I)	nitrogen atom	( C. Demuynck	" " "
C <sub>2</sub> HNS(HCSCN)	thioformyl cyanide	( J.L. Destombes	" " "
C <sub>5</sub> H <sub>6</sub> ArH <sub>3</sub> <sup>+</sup>	cyclopentadiene	(	H/D forms mm, submm wave spectra
C <sub>4</sub> H <sub>6</sub> O	{  2,5 dihydrofuran 2,3 dihydrofuran	) ( in collaboration with Valladolid	mm wave spectrum ring puchering
C <sub>2</sub> D <sub>3</sub> N(CD <sub>3</sub> CN)	methyl cyanide, d <sub>3</sub>	) J. Burie	excited states
C <sub>2</sub> H <sub>2</sub> O	cyclopropenone	) J. Demaison	mm wave spectrum
GeH <sub>3</sub> Cl	germyl chloride	) G. Wlodarczak	" "
C <sub>4</sub> H <sub>5</sub> N(CH <sub>2</sub> CHCH <sub>2</sub> CN)	allyl cyanide	)	" "
C <sub>3</sub> H <sub>6</sub> (CH <sub>2</sub> CHCH <sub>3</sub> )	propene	)	internal rotation
C <sub>2</sub> H <sub>3</sub> N(CH <sub>3</sub> CN)	methyl cyanide	( F. Rohart	low temperature relaxation
C <sub>3</sub> HN(HC <sub>3</sub> N)	cyanoacetylene	( D. Derozier	
CH <sup>15</sup> N(HC <sup>15</sup> N)	hydrogen cyanide	B. Segard	optical bistability

34 Name to whom queries should be addressed C. Matsumura, H. Takeo, M. Sugie

Mailing address National Chemical Laboratory for Industry

1-1 Higashii, Tsukuba, Ibaraki 305

Japan

Telephone: (0298) 54-4521 Telefax: (0298) 54-4488

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{C}_2\text{H}_2\text{N}_2$ ( $\text{HN}=\text{CHCN}$ )	C-cyanomethanimine	H. Takeo	reaction mechanism
$\text{C}_3\text{H}_2\text{NCl}$ ( $\text{H}-\text{C}\equiv\text{C}-\text{CH}=\text{N}-\text{Cl}$ )	N-chloropropargylimine	M. Sugie	spectrum assigned
$\text{C}_3\text{H}_6\text{O}$ ( $\text{CH}_3\text{CH}=\text{CHOH}$ )	2-methylvinylalcohol	H. Takeo	spectrum assigned
$\text{C}_3\text{H}_5\text{N}$		M. Sugie	manuscript in preparation
$\text{C}_3\text{H}_6\text{NCl}$		T. Egawa	joint analysis with electron diffraction

35 Name to whom queries should be addressed Professor D.J. Millen

Mailing address Department of Chemistry  
University College London  
20 Gordon Street, LONDON WC1H 0AJ

Telephone: 01-387 7050 Telefax: 01-380 7463

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{CH}_2\text{FN}$ {HCN...HF}	hydrogen-bonded complex of hydrogen cyanide and hydrogen fluoride	L C Willoughby	Further work on vibrational satellites and isotopic species
$\text{C}_2\text{H}_2\text{N}_2$ {HCN...HCN}	hydrogen-bonded dimer of hydrogen cyanide	K Georgiou	Further work in progress
$\text{C}_3\text{H}_7\text{FO}$ ( $\text{C}_3\text{H}_7\text{O}$ ...HF)	hydrogen-bonded complex of oxetane and hydrogen fluoride	R A Collins	Further work on satellites
$\text{C}_2\text{H}_7\text{FO}$ { $(\text{CH}_3)_2\text{O}$ ...HF}	hydrogen-bonded dimer of dimethyl ether and hydrogen fluoride	H M North J Kisiel	Spectrum partially assigned
$\text{C}_4\text{H}_4\text{FN}$ { $\text{CH}_3\text{CCCN}$ ...HF}	hydrogen-bonded complex of methylcyanoacetylene and hydrogen fluoride	K Georgiou H M North	Work in progress
$\text{C}_5\text{H}_{10}\text{FN}$ { $(\text{CH}_3)_3\text{CCN}$ ...HF}	hydrogen-bonded complex of $\gamma$ -butyl cyanide and hydrogen fluoride	S L A Adebayo L C Willoughby	Work in progress

Prof. R.H. Donweszen  
 Department of Chemistry  
 Michigan State University  
 East Lansing, MI 48824-1322  
 U.S.A.

36 Name to whom queries should be addressed Harald Møllendal

Mailing address Department of Chemistry, The University of Oslo

P. O. Box 1033, Blindern

N-0315 Oslo 3, Norway

Telephone: 02/455674 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

FORMULA	NAME OF COMPOUND	NAME OF INVESTIGATOR	PRESENT STAGE OF PROGRESS
C <sub>3</sub> H <sub>8</sub> OS (CH <sub>3</sub> SCH <sub>2</sub> CH <sub>2</sub> OH)	2-(methylthio)ethanol	H. Møllendal	Acta Chem. Scand. <u>v3</u> (1969) 46
C <sub>3</sub> H <sub>5</sub> NO	2-azetidinone	H. Møllendal	In press, Acta Chem. Scand.
C <sub>4</sub> H <sub>10</sub> O <sub>2</sub> (CH <sub>3</sub> OCH <sub>2</sub> CH <sub>2</sub> CH <sub>2</sub> OH)	3-methoxypropanol	H. Møllendal	In press, Acta Chem. Scand.
C <sub>3</sub> H <sub>8</sub> OS (CH <sub>3</sub> CH(OH)CH <sub>2</sub> SH)	1-mercaptopropanol	H. Møllendal	Writing up
C <sub>5</sub> H <sub>8</sub> O (H <sub>2</sub> C=CH-CH(CH)-CH=CH <sub>2</sub> )	1,4-pentadien-3-ol	H. Møllendal	Assigned
C <sub>4</sub> H <sub>6</sub> O <sub>2</sub>	Cyclopropyl carboxylic acid	H. Møllendal	Partial assignment

37 Name to whom queries should be addressed	<u>Dr. R. N. Nandi</u>		
Mailing address	<u>Microwave Spectroscopy Laboratory</u>		
	<u>Saha Institute of Nuclear Physics</u>		
	<u>92, Acharya Prafulla Chandra Road, Calcutta 700 009, INDIA</u>		
Telephone:	<u>35-4281 - 85</u>		
	<u>Telefax:</u>		
Electronic mail (BITNET or other):			
<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_7H_4BrN$	3-bromobenzonitrile	S. Maiti A. Datta A. I. Jaman	Inconsistent Spectra, work abandoned.
$C_4H_5NO$	Allyl Isocyanate	S. Maiti A. Datta A. I. Jaman R. N. Nandi	Work in progress.
$C_7H_3F_2N$	2,3-difluorobenzonitrile		Work started.

38 Name to whom queries should be addressed Teruhiko OGATA

Mailing address Faculty of Liberal Arts  
Shizuoka University

Ohya, Shizuoka, 422 JAPAN

Telephone: (0542) 37-1111 ex. 8222 Telefax: (0542) 37-9184

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2H_2ClF_3(CF_2Cl-CH_2F)$	1-Chloro-1,1,2-trifluoroethane	M. Yoshikawa	Manuscript in Preparation
$C_4H_6$ ( $H_2C=C=CHCH_3$ )	Methylallene	K. Matsui	$^{13}C$ -2 isotope species in progress
$C_4H_5F$ ( $HC\equiv C-CHFCH_3$ )	3-Fluoro-1-butyne	N. Yamada	Manuscript in preparation
$C_3H_2ClF$ ( $H_2C=C=CFCI$ )	1-Chloro-1-fluoroallene	T. Fukukawa	Manuscript in preparation

39 Name to whom queries should be addressed H. M. Pickett, E. A. Cohen, R. L. Poynter

Mailing address Jet Propulsion Lab.

Telephone: (818) 354-4701      Telefax:                           
                          
                        

Electronic mail (BITNET or other): HPICKETT@CALTECH, EAC@CALTECH

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
COF <sub>2</sub> (OCF <sub>2</sub> )	Carbonyl fluoride	Cohen, L. Brown	Submm, I. R. (ν <sub>1</sub> , 2ν <sub>2</sub> )
BrHO (HOBr)	Hypobromous acid	G. McRae, Cohen in collaboration with Nat. Chem. Lab for Industry (Japan)	Rotational spectrum manuscript prepared. High resolution I.R. in preparation
Cl <sub>2</sub> O <sub>2</sub> (ClOOCl)	Chlorine peroxide (ClO dimer)	M. Birk, R. Friedl, Cohen	Manuscripts in prep- aration
N <sub>2</sub> O <sub>5</sub>	Nitrogen pentoxide	Cohen (with M. Fujitake and E. Hirota, IMS)	In progress
NH <sub>3</sub> (H <sub>3</sub> N)	Ammonia	Poynter, J. Margolis (with H. Sasada, Keio University)	ν <sub>4</sub> , 2ν <sub>2</sub> manuscript in preparation

40 Name to whom queries should be addressed N.Pozdeev

Mailing address Department of Physics, Bashkirian Scientific Centre  
Ural Branch of Academy of Sciences of the USSR  
Tuckaeva 50, 450000, Ufa, USSR

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2H_3ClO(CH_3CClO)$	acetylchloride		centrif. dist., Starck effect
$C_4H_6ClO_2P((CH_3)_2CCOOPCl)$	4,5 dimethyl- 2 chloride- 1,3 dioxophospholen		assigned
$C_4H_8O$	tetrahydrofuran		rot.-vibr. in excited pseudo- rotational states
$C_4H_8S$	tetrahydrothiophene		

41 Name to whom queries should be addressed C. Richard Quade  
Mailing address Department of Physics  
Texas Tech University  
Lubbock, Texas 79409  
Telephone: 806-742-3767 Telefax: \_\_\_\_\_  
Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{CH}_2\text{D}_2\text{O}(\text{CH}_2\text{DOD})$	Methyl Alcohol	Chun Fu Su Dept. of Phys. Miss. State, MS 39762	a-dipole, 0-1, and one Q-type, gauche to gauche series for each molecule assigned.
$\text{CD}_3\text{O} (\text{CD}_2\text{OD})$			

42 Name to whom queries should be addressed Dr. Jim Roberts

Mailing address Department of Physics  
University of North Texas  
Denton, TX 76203

Telephone: 817-565-3281 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$^{12}\text{CH}_3\ ^{13}\text{C}^{15}\text{N}$	METHYL CYANIDE	M. Al-Share, H. S. Tam and J. A. Roberts	Rotational Components for $1 \leq J \leq 5$ measured and assigned for the $v_8$ with $v=2, 3$ . Large frequency departures $\approx 6(J+1)$ MHz observed for some components.

43 Name to whom quieries should be addressed Georges ROUSSY

Mailing address LABORATOIRE DE SPECTROSCOPIE ET DES TECHNIQUES MICROONDÉS

UNIVERSITÉ NANCY I - U.A. C.N.R.S. 1105

B.P. 239 - 54506 VANDOEUVRE LES NANCY CEDEX (FRANCE)

Telephone : 83.91.20.48 Telefax :

Electronic mail (BITNET or other) :

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>7</sub> H <sub>12</sub> S	1-thia spiro (2-5) octane	H. BOULEBNANE E. ALAMI R. VILLAMANAN*	To be published in J. Mol. Spectry
C <sub>8</sub> H <sub>14</sub> O <sub>2</sub>	1-4 dioxa Spiro (4-5) decane	E. ALAMI G. ROUSSY	Work in progress
C <sub>8</sub> H <sub>14</sub> S <sub>2</sub>	1-4 dithia spiro (4-5) decane	E. ALAMI G. ROUSSY	Work in progress
C <sub>4</sub> H <sub>5</sub> O <sub>2</sub>	2-methyl-1,3 dioxolane	E. ALAMI G. ROUSSY	Work started
C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>	2,2-dimethyl-1,3 dioxolane	E. ALAMI G. ROUSSY	Work started

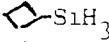
\* Permanent address : Departamento de Química-Física  
UNIVERSIDAD de VALLADOLID  
SPAIN

44 Name to whom queries should be addressed: H.D.Rudolph

Mailing address Institute of Physical Chemistry  
University of Ulm  
P.O.B. 4066, Oberer Eselsberg  
7900 Ulm, W-Germany

Telephone number 0731-176-1302

<u>Formula</u>	<u>Name of Compound</u>	<u>Name of Investigator</u>	<u>Present state of Progress</u>
AqH	Silver Hydride	H. Birk, H. Jones	Diode laser
BaH	Barium Monohydride	C. Magg, H. Birk H. Jones	Diode laser. Chem. Phys. Lett. <b>149</b> , 321 (1988)
CF <sub>2</sub>	Difluoro carbene	H.Birk and H.Jones	Diode laser ν <sub>1</sub> in free expansion jet
CF <sub>3</sub> J	Trifluoriodo-methane	W.Höhe, L.Jörissen W.A.Kreiner (with H.Bürger, Wuppertal)	ν <sub>1</sub> fundamental
HCCl	Chlorocarbene	H.Birk and H.Jones	Diode laser ν <sub>3</sub>
HCCF	Fluoroacetylene	J.Lindenmaier D.Papousek, H.Jones	Diode laser ν <sub>3</sub> paper ready
C <sub>2</sub> H <sub>7</sub> N (H <sub>3</sub> CCH <sub>2</sub> NH <sub>2</sub> )	Ethylamine	E.Fischer, I.Botskor, H.D.Rudolph	ms in preparation
C <sub>3</sub> H <sub>7</sub> Cl (H <sub>3</sub> CCCH <sub>2</sub> Cl)	Propargyl-chlorid	L.Braun I.Botskor	ms in preparation
C <sub>3</sub> H <sub>7</sub> N (H <sub>2</sub> CCHCH <sub>2</sub> NH <sub>2</sub> )	Allylamine	K.H.Wiedenmann, I.Botskor, H.D.Rudolph	structure evaluation work completed
C <sub>3</sub> H <sub>8</sub> Ge $\Delta$ -GeH <sub>3</sub>	Cyclopropyl-germane	K.Epple, H.D. Rudolph	spectra of 31 isotopic species assigned, structure
C <sub>3</sub> H <sub>8</sub> Si $\Delta$ -SiH <sub>3</sub>	Cyclopropyl-silane	J.Mennicke, B.Mir, H.D.Rudolph	spectra of several isotopic species, structure

$C_4H_{10}Si$	Cyclobutyl-silane 	A.Ruck, B.Mir, H.D. Rudolph	heavy atom structure, additional isotopic species of both conformeres
CsH	Cesium Hydride	U. Magg and H. Jones	Diode laser Chem. Phys. Lett. 148, 6 (1988)
GaH	Gallium Hydride	R-D. Urban, U. Magg, and H. Jones	Diode laser, Chem. Phys. Lett. 154, 135 (1988) deuteride in progress
IF	Iodine Fluoride	K.P.R. Nair, U. Magg, H. Birk and H. Jones	Diode laser sub. Z. Naturforsch
InH	Indium Hydride	A. H. Bahnmaier, R-D.Urban, and H.Jones	Diode laser accepted Chem. Phys. Lett. deuteride in progress
PbH	Lead Hydride	U. Magg and H. Jones	Diode laser submitted Chem. Phys. Lett
PH <sub>3</sub>	Phosphine	D. Papousek, H. Birk, U. Magg, and H. Jones	Diode laser A <sub>1</sub> -A <sub>2</sub> splittings J.Mol.Spectr in press
RbH	Rubidium Hydride	U. Magg, H. Birk, and H. Jones	Diode laser Chem. Phys. Lett. 151, 503 (1988)
SrH	Strontium Hydride	U. Magg, H. Birk, and H. Jones	Diode laser Chem. Phys Lett. 151, 263-266 (1988)
TlH	Thallium Hydride	R-D. Urban, U. Magg A.H.Bahnmaier,H. Jones	Diode laser sub. Chem. Phys. Lett. pot. function Deuteride in progress
ZnH	Zink Hydride	U. Magg, R.-D. Urban, H. Jones	Diode laser, ms in preparation

45 Name to whom queries should be addressed Shuji SAITO

Mailing address Department of Astrophysics

Nagoya University

Chikusa, Nagoya 464-01, Japan

Telephone: 052-781-5111, ex. 6672 telefax: 052-782-1029

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>3</sub> H	C <sub>3</sub> H radical	S. Yamamoto	Astrophys. J. Submitted.
C <sub>3</sub> D		S. Yamamoto	<sup>34</sup> S, <sup>13</sup> C species.
C <sub>2</sub> S(CCS)	CCS radical	S. Saito K. Kawaguchi*	Manuscript in preparation.
		M. Tanimoto** S. Saito S. Yamamoto K. Kawaguchi*	v <sub>1</sub> , v <sub>2</sub> , v <sub>3</sub> excited states. Manuscript in preparation.
C <sub>3</sub> H	Cyclic C <sub>3</sub> H radical	S. Yamamoto	D, <sup>13</sup> C species. Manuscript in preparation.
CP	CP radical	S. Saito	Astrophys. J. in press.
C <sub>2</sub> H <sub>2</sub> N C <sub>2</sub> D <sub>2</sub> N	CH <sub>2</sub> CN radical	S. Saito	Astrophys. J. <u>334</u> , L113 (1988). H, D species. Manuscript in preparation.
CH <sub>2</sub> N	CH <sub>2</sub> N radical	S. Yamamoto	Manuscript in preparation.
HS <sub>2</sub>	HS <sub>2</sub> radical	S. Yamamoto	Manuscript in preparation.
C <sub>5</sub> H	C <sub>5</sub> H radical	S. Takano	v <sub>9</sub> excited state. Manuscript in preparation.
C <sub>3</sub> H	C <sub>3</sub> H radical	M. Kanada	<sup>13</sup> C species, v <sub>4</sub> excited states. Manuscript in preparation.
C <sub>3</sub> N	C <sub>3</sub> N radical	H. Mikami	v <sub>5</sub> excited state. Astron. Astrophys. Lett. submitted.

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
CHN <sub>2</sub>	HNCN radical	S. Yamamoto	Spectrum assigned.
H <sub>2</sub> NS	H <sub>2</sub> NS radical	S. Yamamoto	Spectrum assigned.
D <sub>2</sub> N	NH <sub>2</sub> -d <sub>2</sub> radical	M. Kanada	Manuscript in preparation.
CaS	Calcium sulfide	S. Takano	Manuscript in preparation.
MgS	Magnesium sulfide	S. Takano	Manuscript in preparation.
C <sub>2</sub> <sup>DN</sup>	HCCN-d radical	F. X. Brown S. Saito	Spectrum assigned.
H <sub>2</sub> NO	H <sub>2</sub> NO radical	H. Mikami	Spectrum assigned.

\*Nobeyama Radio Observatory. Shizuoka University.

46 Name to whom queries should be addressed Yoshiaki SASADA

Mailing address College of Engineering,

Kanto Gakuin University,

Kanazawa-ku, Yokohama, 236, JAPAN

Telephone: 045 - 781 - 2001 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_4H_3SCl \left( \begin{array}{c} Cl \\   \\ S \end{array} \right)$	3-chlorothiophene	Y. Sasada	Work in progress
$C_4H_3SI \left( \begin{array}{c} I \\   \\ S \end{array} \right)$	2-iodothiophene	Y. Sasada Y. Niide I. Ohkoshi	Two excited states manuscript in preparation
$C_4H_3SI \left( \begin{array}{c} I \\   \\ S \end{array} \right)$	3-iodothiophene	Y. Sasada Y. Niide I. Ohkoshi	Work in Progress

47 Name to whom queries should be addressed: Richard H. Schwendeman

Mailing address Department of Chemistry

Michigan State University

East Lansing, MI 48824-1322

Telephone: 517-353-9412 Telefax: 517-353-1793

Electronic mail (BITNET or other) BITNET%"SCHWEN@MSUCEM"

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{CH}_2\text{O}(\text{CD}_2\text{O})$	Formaldehyde-d <sub>2</sub>	S.-C. Hsu	IR-MW sideband laser $\nu_2$ band, in press J Mol. Spectrosc.
$\text{CH}_3\text{I}(\text{CD}_3\text{I})$	Methyl iodide-d <sub>3</sub>	H.-G. Cho	IR-MW sideband laser IR-RF double resonance $\nu_2$ band, in press J Mol. Spectrosc.
$\text{H}_3\text{N}(\text{NH}_3)$	Ammonia	Y. Matsuo S. K. Lee	IR-MW sideband laser
$\text{CH}_3\text{F}({}^{13}\text{CH}_3\text{F})$	Methyl fluoride	Y. Matsuo	IR-MW sideband laser IR-IR double resonance
$\text{CH}_3\text{F}$	Methyl fluoride	H.-G. Cho Y. Matsuo	IR-MW sideband laser $\nu_3 + \nu_6 - \nu_6$ band
$\text{CH}_3\text{Br}(\text{CD}_3\text{Br})$	Methyl bromide-d <sub>3</sub>	H. G. Cho	IR-MW sideband laser IR-RF double resonance $\nu_2$ band
$\text{CH}_3\text{I}$	Methyl iodide	W. Fawzy	IR-RF double resonance Lineshape analysis
$\text{CH}_4\text{O}(\text{CH}_3\text{OH})$	Methanol	D. Peterson	IR-MW double resonance $\nu_8$ band
$\text{CF}_3\text{I}$	Trifluoromethyl iodide	D. Peterson	IR-MW double resonance $\nu_1$ band

48 Name to whom queries should be addressed Dr.T.Shimizu

Mailing address Department of Physics Faculty of Science University  
of Tokyo, Hongo 7-3-1, Bunkyo-Ku Tokyo 113, Japan

Telephone: 03-812-2111 Telefax: 03-814-9717

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
H <sub>3</sub> N(NH <sub>3</sub> )	Ammonia	H.Odashima	Pressure-broadening parameter determined
COS(OCS)	Carbonyl Sulfide	H.Odashima	Pressure-broadening parameter determined

49 Name to whom queries should be addressed C. F. Su or R. L. Cook

Mailing address Department of Physics  
Mississippi State University  
Mississippi State, MS 39762

Telephone 601-325-2806 Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2H_6O_2$	Meso-Bisoxirane (gauche form)		Assignments of a- and c-type transitions have been completed. (paper in preparation)
$CH_2H_6O_2$	d1-Bisoxirane (gauche form)		in progress
$C_3H_5SCl$	3-chloropropylene sulfide		Assignments of a- and b-type transitions have been completed. Calculation of coupling constant of Cl has been completed. (paper in preparation)

50 Name to whom queries should be addressed M. Takami

Mailing address RIKEN, The Inst. Phys. Chem. Res., Wako, Saitama 351-01,  
Japan

Telephone: 0484(62)1111 ext.3611 Telefax: 0484(62)1449

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
H <sub>4</sub> Sn(SnH <sub>4</sub> )	Stannane	L.G.Jörissen Y.Ohshima	v <sub>1</sub> /v <sub>3</sub> MW spectrum J.Chem.Phys. in press Extended analysis in progress
CF <sub>4</sub>	Carbontetrafluoride	M.Takami	v <sub>3</sub> MW spectrum extended analysis in progress

51 Name to whom queries should be addressed Prof. Mitsuru Takano

Mailing address Department of Mathematics and Physics

National Defense Academy

1-10-20, Hashirimizu, Yokosuka 239, Japan

Telephone: 0468-41-3810 ext 2212 Telefax: 0468-43-6236

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$C_2H_4BrCl$ ( $ClCH_2CH_2Br$ )	1-Bromo-2-chloroethane	M. Takano I. Ohkoshi Y. Niide	In progress
$C_2H_4FI$ ( $FCH_2CH_2I$ )	1-Iodo-2-fluoroethane	Y. Niide I. Ohkoshi	In progress
$CH_2BrCl$	Bromochloromethane	Y. Niide I. Ohkoshi	Manuscript submitted <i>J. Mol. Spectrosc.</i>
$C_2H_2OBrCl$ ( $BrCH_2COCl$ )	Bromoacetyl Chloride	Y. Niide I. Ohkoshi	In progress

52 Name to whom queries should be addressed Takehiko Tanaka, Keiichi Tanaka  
 Mailing address Department of Chemistry, Faculty of Science,  
Kyushu University 33, Hakozaki, Higashiku,  
Fukuoka, 812 JAPAN  
 Telephone number 092 - 641 - 1101 - Ext 4250

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>3</sub> H <sub>4</sub>	Allene	K. Tanaka	Vibrationally induced transitions. Manuscript in preparation.
C <sub>7</sub> H <sub>6</sub> O <sub>2</sub>	Tropolone	H. Onaka	Ground state spectrum assigned.
C <sub>4</sub> HF (HC≡C-C≡CF)	Fluorodiacetylene	T. Okabayashi	Submitted
C <sub>3</sub> FN (FC≡CC≡N)	Fluorocyanooacetylene	T. Okabayashi	Excited vibrational states. Manuscript in preparation
CHNO (HCNO)	Fulminic acid	R. Takashi	LMDR (Laser-microwave double resonance) in progress
POF <sub>3</sub>	Phosphoryl fluoride	K. Someya	LMDR manuscript in preparation
SiHF <sub>3</sub> (SiHF <sub>3</sub> , SiDF <sub>3</sub> )	Trifluorosilane	K. Harada	LMDR in progress
CH <sub>3</sub> F (CH <sub>3</sub> F, CD <sub>3</sub> F)	Methyl fluoride	K. Harada	LMDR manuscript in preparation
CH <sub>3</sub> I	Methyl iodide	K. Harada	LMDR in progress
C <sub>2</sub> HF (HC≡CF)	Fluoroacetylene	Y. Nakahara	LMDR in progress
C <sub>2</sub> H <sub>3</sub> N (CH <sub>3</sub> CN)	Methyl Cyanide	T. Oyama	LMDR in progress
C <sub>2</sub> H <sub>3</sub> N (CH <sub>3</sub> NC)	Methyl Isocyanide	T. Oyama	LMDR manuscript in preparation

C <sub>2</sub> HN (DC≡CC≡N)	Cyanoacetylene	K. Tanaka	LMDR in progress
CFN (FC≡N)	Cyanogen fluoride	S. Matsuba	LMDR in progress

53 Name to whom queries should be addressed Mitsutoshi Tanimoto

Mailing address Department of Chemistry, Faculty of Science

Shizuoka University

836 Ohya, Shizuoka 442, JAPAN

Telephone: 0542-37-1111 (5600) Telefax: 0542-37-9895

Electronic mail (BITNET or other): \_\_\_\_\_

FORMULA	NAME OF COMPOUND	NAME OF INVESTIGATOR	PRESENT STAGE OF PROGRESS
O <sub>2</sub> S (SO <sub>2</sub> )	Sulfur Dioxide	Y. Morino M. Tanimoto with S. Saito	r-structure potential function
Cl <sub>2</sub> Si(SiCl <sub>2</sub> )	Dichlorosilylene	M. Tanimoto with C. Matsumura H. Takeo, and M. Fujitake	manuscript in preparation
ClSi	Silicon monochloride	M. Tanimoto with S. Saito	v=1 spectrum observed
F <sub>3</sub> Si (SiF <sub>3</sub> )	Silicon trifluoride	M. Tanimoto with S. Saito	work in progress

54 Name to whom queries should be addressed: Nancy S. True

Mailing address Department of Chemistry

The University of California

Davis, CA 95616

Telephone: 916-752-0874 Telefax: 916-752-6363

Electronic mail (BITNET or other)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>6</sub> H <sub>5</sub> -CH <sub>2</sub> X	Benzyl Halides	S. A. Sorenson N. S. True	M.S. in preparation
C(O)F-O-C(H)-CH <sub>2</sub>	Vinyl Fluoroformate	N. S. True	M.S. in preparation
C(O)Cl-O-C(H)-CH <sub>2</sub>	Vinyl Chloroformate	N. S. True	Spectrum Assigned, <sup>35</sup> Cl & <sup>37</sup> Cl isotopes
p-COH-C <sub>6</sub> H <sub>4</sub> -OCH <sub>3</sub>	p-Anisaldehyde	S. A. Sorenson N. S. True	M.S. in preparation

55 Name to whom queries should be addressed Shozo Tsunekawa

Mailing address Department of Physics

Toyama University

Toyama 930 Japan

Telephone: (0764)41-1271 (ext 318) Telefax: \_\_\_\_\_

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
$\text{CH}_5\text{N}$ ( $\text{CH}_3\text{NH}_2$ , $\text{CH}_3\text{NHD}$ )	Methylamine	K. Takagi S. Tsunekawa M. Iziri	Manuscript in Preparation
$\text{CH}_4\text{SH}$ ( $\text{CH}_3\text{SD}$ , $\text{CD}_3\text{SH}$ )	Methyl Mercaptan	T. Hayashi S. Tsunekawa K. Nakagawa (Zyousai Univ.)	Work in progress ( $\text{CD}_3\text{SH}$ species)
$\text{H}_4\text{N}_2$ ( $\text{ND}_2\text{ND}_2$ )	Hydrazine	S. Tsunakawa	Work in progress
$\text{C}_2\text{H}_5\text{O}$ ( $\text{CH}_3\text{CONH}_2$ )	Acetamide	K. Nakagawa S. Tsunekawa	Excited State
$\text{CH}_4\text{O}$ ( $^{13}\text{CH}_3\text{OH}$ , $\text{CH}_3\text{OH}$ )	Methyl alcohol	K. Takagi M. Hayashi Y. Hoshino M. Ohishi	Work in progress

56 Name to whom queries should be addressed: Manfred Winnewisser

Mailing address: Physikalisch-Chemisches Institut  
Justus-Liebig-Universität Giessen  
Heinrich-Buff-Ring 58  
D-6300 Giessen, Fed. Rep. Germany

Telephone number: (0641) 702-5790 (BITNET: MWI at DGIHRZ01)

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STATE OF PROGRESS</u>
C <sub>2</sub> H <sub>3</sub> NO (CH <sub>3</sub> CNO)	Acetonitrile oxide	(J. Galica, Poznan)  B. P. Winnewisser, M. Winnewisser	Exc. states MMW work in progress
C <sub>3</sub> OS (OCCCS)	Tricarbon oxide sulfide	M. Winnewisser	Comb. states, MMW work in progress
CD <sub>2</sub> N <sub>2</sub> (ND <sub>2</sub> NC)	d <sub>2</sub> -Isocyanamide	F. Stroh, M. Winnewisser	MMW work in progress
CH <sub>2</sub> N <sub>2</sub> (HNCNH)	Carbodiimide	M. Birk, M. Winnewisser, (E. A. Cohen, JPL)	0 <sup>+</sup> , 0 <sup>-</sup> states ass. in MMW, FIR MS in prep.
C <sub>2</sub> H <sub>2</sub> N <sub>2</sub> O (NCCONH <sub>2</sub> )	Cyanoformamide	(J. J. Christiansen, Copenhagen)	Ground state and exc. states: MS in prep.
CH <sub>4</sub> O <sub>2</sub> (CH <sub>3</sub> OOH)	Methyl hydroperoxide	C. E. Blom (M. Tyblewski, A. Bauder, Zürich)	Manus. in preparation
CH <sub>4</sub> O <sub>2</sub> (CH <sub>3</sub> OOD)	Methyl deuteriohydroperoxide	C. E. Blom	"
C <sub>3</sub> H <sub>4</sub> O (CH <sub>2</sub> CHCHO)	Acrolein s-cis, s-trans conformers ass.	C. E. Blom	Exc. states of
CHNO (HCNO)	Fulminic acid	G. Wagner, J. Preusser, M. Winnewisser B. P. Winnewisser	<sup>13</sup> C, <sup>15</sup> N species meas. in progress, 1-type doublets assigned.
C <sub>2</sub> N <sub>2</sub> (CNCN)	Isocyanogen	F. Stroh, M. Winnewisser, (M.C.L.Gerry, Vancouver)	MW and MMW lines assigned, <sup>14</sup> N hpf resolved with FT in Kiel.

57 Name to whom queries should be addressed K.M.T. Yamada and G. Winnewisser

Mailing address I.Physikalisches Institut, Universität zu Köln  
D-5000 Köln 41, West Germany

---

Telephone: 0221-470-3567 Telefax: TELEX 8 882 291 UNIK D

Electronic mail (BITNET or other): ABA04 at DKØRRZKØ.BITNET

---

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
NO	Nitric Oxide	A. Saleck	$^{18}\text{O}$ , $^{15}\text{N}$ measured in mmw.
$\text{H}_2\text{S}_2$	Hydrogen disulfied	P. Mittler	$^{34}\text{S}$ , D, and Ex.St. measured in mmw.
$\text{H}_2\text{S}_3$	Hydrogen trisulfied	D. Mauer	Papers published.
$\text{H}_2\text{S}_4$	Hydrogen tetrasulfied	A. Saleck	Measurement in progress in mmw.
$\text{H}_2\text{C}_2\text{O}$	Ketene	T. Wang	Ex.St. in mmw. Work in progress.

(continued)

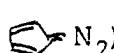
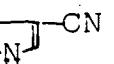
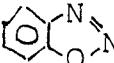
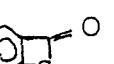
$C_7H_4ClN$	2-Chlorobenzonitrile	M.Onda	hfs in progress
$C_7H_5NO$	Phenyl cyanate	M.Onda	ms in prep.
$C_7H_6O_2$	Benzoic acid	M.Onda	Isotopic species in progress
$C_7H_6O_3$	Salicylic acid	M.Onda	in progress
$C_7H_6OS$	Thiobenzoic acid	M.Onda	Spectr. assigned
$C_7H_7Cl$	Benzyl chloride	K.Suga	in progress
$C_7H_8O$	Anisole	M.Onda	$CD_3$ , excited state ms in prep.
$C_7H_8O$	Benzyl alcohol	K.Suga	in progress
$C_8H_7N$	Benzyl cyanide	K.Suga	Spectr. assigned
$C_8H_8O$	Acetophenone	M.Onda	ms in prep.
$C_{12}H_9F$	4-Fluorobiphenyl	M.Onda	Spectr. assigned

58 Name to whom queries should be addressed Ichiro Yamaguchi

Mailing address Department of Chemistry, Sophia University,  
7 Kioicho, Chiyodaku, Tokyo 102, Japan

Telephone: 03-238-3358 Telefax: 03-238-3885

Electronic mail (BITNET or other): \_\_\_\_\_

<u>FORMULA</u>	<u>NAME OF COMPOUND</u>	<u>NAME OF INVESTIGATOR</u>	<u>PRESENT STAGE OF PROGRESS</u>
C <sub>3</sub> H <sub>6</sub> OS (CH <sub>3</sub> CH <sub>2</sub> COSH)	Thiopropionic acid	I.Yamaguchi	in progress
C <sub>2</sub> H <sub>4</sub> OS (HSCH <sub>2</sub> CHO)	Mercaptoacetaldehyde	H.Hatcho I.Yamaguchi	in progress
C <sub>2</sub> H <sub>4</sub> O <sub>2</sub> S (HSCH <sub>2</sub> COOH)	Mercaptoacetic acid	I.Yamaguchi	in progress
C <sub>4</sub> H <sub>6</sub> O <sub>2</sub> (CH <sub>3</sub> CH=CHCOOH)	2-Butenoic acid	I.Yamaguchi	J.Mol.Struct. 190, 377(1988)
C <sub>4</sub> H <sub>7</sub> NO (CH <sub>3</sub> CH=CHCH=NOH)	Crotonaldehyde oxime	M.Hamano	ms in prep.
C <sub>4</sub> H <sub>7</sub> NO (CH <sub>2</sub> =C(CH <sub>3</sub> )CH=NOH)	syn-2-Methylacrylaldehyde oxime	Y.Yanagawa	ms in prep.
C <sub>4</sub> H <sub>9</sub> NO (CH <sub>3</sub> (CH <sub>2</sub> ) <sub>2</sub> CH=NOH)	Butyraldehyde oxime	O.Ohashi	in progress
C <sub>5</sub> H <sub>9</sub> NO (CH <sub>2</sub> (CH <sub>2</sub> ) <sub>3</sub> C=NOH)	Cyclopentanone oxime	A.Murakami	ms in prep.
C <sub>6</sub> H <sub>11</sub> NO (CH <sub>2</sub> (CH <sub>2</sub> ) <sub>4</sub> C=NOH)	Cyclohexanone oxime	O.Ohashi	ms in prep.
C <sub>4</sub> H <sub>7</sub> NO (HNCH <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> C=O)	2-Pyrrolidone	O.Ohashi	ms in prep.
C <sub>5</sub> H <sub>9</sub> NO (HNCH <sub>2</sub> (CH <sub>2</sub> ) <sub>3</sub> C=O)	2-Piperidone	O.Ohashi	assigned
C <sub>6</sub> H <sub>9</sub> NO (CH <sub>2</sub> (CH <sub>2</sub> ) <sub>2</sub> CHCHC=NOH)	2-Cyclohexen-1-one oxime	Y.Sato	assigned
C <sub>4</sub> H <sub>9</sub> NO ((CH <sub>3</sub> ) <sub>2</sub> CHCH=NOH)	Isobutyraldehyde oxime	A.Murakami	assigned
C <sub>3</sub> H <sub>7</sub> ND (CH <sub>3</sub> CH <sub>2</sub> CH=NOH)	E-sp-Propionaldehyde oxime	F.Kato	assigned
CHNS (HSCN)	Thiocyanic acid	T.Sakaizumi	in progress
C <sub>2</sub> H <sub>3</sub> NO (CH <sub>3</sub> OCN)	Methyl cyanate	H.Mure T.Sakaizumi	<sup>14</sup> N, <sup>15</sup> N assigned
C <sub>3</sub> H <sub>5</sub> NO (CH <sub>3</sub> CH <sub>2</sub> OCN)	Ethyl cyanate	H.Mure T.Sakaizumi	ms in prep.
C <sub>5</sub> H <sub>4</sub> N <sub>2</sub> (  N <sub>2</sub> )	Diazocyclopentadiene	S.Fukuda T.Sakaizumi	ms in prep.
C <sub>5</sub> H <sub>4</sub> N <sub>2</sub> (  N H )	Pyrrole-2-carbonitrile	T.Sakaizumi	ms in prep.
C <sub>6</sub> H <sub>4</sub> N <sub>2</sub> S (  O N S )	1,2,3-Benzothiadiazole	T.Sakaizumi	ms in prep.
C <sub>7</sub> H <sub>4</sub> OS (  O S )	Benzothiet-2-one	S.Fukuda T.Sakaizumi	assigned

## FORMULA INDEX

AgH - silver hydride - 44	CF <sub>2</sub> - difluorocarbene - 20,44
AlO - aluminum monoxide - 24	CF <sub>2</sub> O - carbonyl fluoride - 39
ArF <sub>3</sub> P - argon-PF <sub>3</sub> complex - 6,27	CF <sub>3</sub> I - trifluoriodomethane - 44,47
ArH <sub>2</sub> O - argon water complex - 32	CF <sub>3</sub> NO <sub>2</sub> - trifluoronitromethane - 11
ArH <sub>3</sub> <sup>+</sup> - Ar-H <sub>3</sub> <sup>+</sup> complex - 33	CF <sub>3</sub> NSi - trifluorosilylisocyanide - 11
ArO <sub>2</sub> S - argon sulfur dioxide complex - 32	CF <sub>3</sub> NSi - trifluorosilylcyanide - 11
Ar <sub>2</sub> ClH - Ar <sub>2</sub> -HCl trimer - 20	CF <sub>4</sub> - carbon tetrafluoride - 50
Ar <sub>3</sub> ClH - Ar <sub>3</sub> -HCl tetramer - 20	CF <sub>8</sub> S - trifluoromethyl sulfur pentafluoride - 3
BBrS - bromosulfidoboron - 26	CHArF <sub>3</sub> - Ar-HCF <sub>3</sub> complex - 30
BClSe - chloroselenidoboron - 26	CHBrO <sub>2</sub> - carbon dioxide-hydrogen bromide cluster - 5
BF - boron fluoride - 15	CHCl - chlorocarbene - 44
BFS - fluorosulfidoboron - 26	CHClF <sub>2</sub> - chlorodifluoromethane - 13
BF <sub>2</sub> HO - hydroxodifluoroboron - 13	CHClO <sub>2</sub> - carbon dioxide-hydrogen chloride cluster - 5
BF <sub>3</sub> - trifluoroboron - 13	CHCl <sub>3</sub> - chloroform - 8
BH <sub>4</sub> N - aminoborane - 13	CHFO <sub>2</sub> - carbon dioxide-hydrogen fluoride cluster - 5
B <sub>2</sub> ClH <sub>7</sub> - bent diborane-HCl dimer - 20	CHF <sub>2</sub> N - difluoromethanimine - 14
BaH - barium monohydride - 44	CHF <sub>2</sub> P - 2,2-difluorophosphaethene - 26
BrHN <sub>2</sub> - N <sub>2</sub> -HBr complex - 30	CHF <sub>3</sub> - fluoroform - 22
BrHO - hypobromous acid - 39	CHF <sub>3</sub> S - trifluoromethylmercaptan - 11
BrH <sub>3</sub> O - H <sub>2</sub> O-HBr complex - 30	CHN - hydrogen cyanide - 13,33
BrIn - indium monobromide - 25	CHNO - isocyanic acid - 13,17
CBrF <sub>2</sub> N - N-bromodifluoromethanimine - 14	CHNO - fulminic acid - 52,56
CClF <sub>2</sub> NO - chlorodifluoronitrosomethane - 11	CHNS - thiocyanic acid - 58
CClN - cyanogen chloride - 15	CHN <sub>2</sub> - HNCN radical - 45
CClNS - chlorine thiocyanate - 17	CHO <sup>+</sup> - formyl ion - 15
CCl <sub>2</sub> - dichlorocarbene - 24	
CFN - cyanogen fluoride - 13,52	

$\text{CH}_2\text{BrCl}$	- bromochloromethane	- 51	$\text{CH}_3\text{O}$	- methoxy radical	- 24
$\text{CH}_2\text{BrP}$	- 1-bromophosphaethene	- 26	$\text{CH}_4$	- methane	- 13
$\text{CH}_2\text{Br}_2$	- dibromomethane	- 13	$\text{CH}_4\text{ArO}$	- argon-methanol complex	- 18
$\text{CH}_2\text{FN}$	- HCN-HF complex	- 35	$\text{CH}_4\text{O}$	- methanol	- 29,41,47,55
$\text{CH}_2\text{FP}$	- 1-fluorophosphaethene	- 26	$\text{CH}_4\text{O}_2$	- methyl hydroperoxide	- 4,56
$\text{CH}_2\text{F}_3\text{N}$	- trifluoromethylamine	- 9	$\text{CH}_4\text{S}$	- methyl mercaptan	- 55
$\text{CH}_2\text{I}_2$	- diiodomethane	- 13	$\text{CH}_5\text{BO}_2$	- methylborondihydroxide	- 11
$\text{CH}_2\text{N}$	- $\text{CH}_2\text{N}$ radical	- 45	$\text{CH}_5\text{BrSi}$	- methylbromosilane	- 23
$\text{CH}_2\text{N}_2$	- isocyanamide	- 56	$\text{CH}_5\text{FN}_2$	- $\text{NH}_3$ -HCN-HF trimer	- 20
$\text{CH}_2\text{N}_2$	- carbodiimide	- 56	$\text{CH}_5\text{N}$	- methylamine	- 13,29,55
$\text{CH}_2\text{O}$	- formaldehyde	- 47	$\text{CH}_6\text{ClN}$	- $\text{CH}_3\text{NH}_2$ -HCl complex	- 30
$\text{CH}_2\text{O}_2$	- formic acid	- 28	CNS	- NCS radical	- 2
$\text{CH}_2\text{O}_2$	- water carbon monoxide complex	- 32	CO	- carbon monoxide	- 28
$\text{CH}_2\text{O}_2\text{S}$	- hydrogen sulfide carbon dioxide complex	- 32	COS	- carbonyl sulfide	- 13,15,48
$\text{CH}_3\text{ArNO}$	- argon-formamide complex	- 18	CP	- CP radical	- 45
$\text{CH}_3\text{Br}$	- methyl bromide	- 12,47	$\text{C}_2\text{F}_3\text{NO}$	- trifluoromethylisocyanate	- 9
$\text{CH}_3\text{Cl}_3\text{Si}$	- methyltrichlorosilane	- 8	$\text{C}_2\text{F}_3\text{N}_2\text{S}_2$	- 4-(trifluoromethyl)-1,2,3,5-dithiadiazole	- 9
$\text{CH}_3\text{F}$	- methyl fluoride	- 47,52	$\text{C}_2\text{F}_3\text{P}$	- 3,3,3-trifluoro-1-phosphapropyne	- 26
$\text{CH}_3\text{F}_2\text{O}_2\text{P}$	- methyldifluorophosphate	- 18	$\text{C}_2\text{F}_4$	- perfluoromethyl carbene	- 20
$\text{CH}_3\text{F}_2\text{P}$	- difluoromethylphosphine	- 31	$\text{C}_2\text{F}_6\text{NO}$	- bis-trifluoromethylnitroxide radical	- 19
$\text{CH}_3\text{F}_3\text{Si}$	- methyl trifluorosilane	- 4,8	$\text{C}_2\text{H}$	- ethynyl radical	- 24,33
$\text{CH}_3\text{I}$	- methyl iodide	- 12,47,52	$\text{C}_2\text{HBr}$	- bromoacetylene	- 13
$\text{CH}_3\text{N}$	- methanimine	- 13	$\text{C}_2\text{HBrClF}_3$	- halothane	- 12
$\text{CH}_3\text{NO}$	- formamide	- 7	$\text{C}_2\text{HBrO}$	- bromoketene	- 17
$\text{CH}_3\text{NO}$	- formaldoxime	- 17	$\text{C}_2\text{HCl}$	- chloroacetylene	- 13
$\text{CH}_3\text{NO}_3$	- methyl nitrate	- 13	$\text{C}_2\text{HClF}_2\text{O}$	- chlorodifluoroacetaldehyde	- 11
$\text{CH}_3\text{N}_3$	- methyl azide	- 13,17			

$C_2HClO$ - chloroketene - 7	$C_2H_3F_3$ - $CH_2CF_2$ -HF complex - 30
$C_2HF$ - fluoroacetylene - 13,44,52	$C_2H_3N$ - methyl cyanide - 33,42,52
$C_2HFO$ - fluoroketene - 7	$C_2H_3N$ - methyl isocyanide - 52
$C_2HF_3O$ - trifluoroethylene oxide- 18	$C_2H_3NO$ - acetonitrile oxide - 56
$C_2HF_3O_2$ - trifluoroacetic acid - 4	$C_2H_3NO$ - methyl cyanate - 58
$C_2HI$ - iodoacetylene - 13	$C_2H_3NO_2$ - nitroethylene - 13
$C_2HN$ - HCCN radical - 45	$C_2H_3N_3$ - 1H-1,2,3-triazole - 10
$C_2HNO_2$ - OCO-HCN linear dimer - 20	$C_2H_3N_3$ - 2H-1,2,3-triazole - 10
$C_2HNS$ - thioformyl cyanide - 33	$C_2H_3O$ - acetyl radical - 24
$C_2H_2BrClO$ - bromoacetyl chloride - 51	$C_2H_4BrCl$ - 1-bromo-2-chloroethane - 51
$C_2H_2BrFO$ - fluoroacetyl bromide - 14	$C_2H_4FI$ - 2-iodo-1-fluoroethane - 51
$C_2H_2ClF_3$ - 1-chloro-1,1,2-trifluoroethane - 38	$C_2H_4FNO$ - 2-fluoroacetamide - 13
$C_2H_2FN$ - fluoroacetonitrile - 13	$C_2H_4F_2$ - $CH_2CHF$ -HF complex - 30
$C_2H_2FNO$ - OC-HCN-HF trimer - 20	$C_2H_4F_3N$ - trifluoroethylamine - 13
$C_2H_2N$ - $CH_2CN$ radical - 45	$C_2H_4N_2O$ - $H_2O$ -HCN-HCN trimer - 20
$C_2H_2N_2$ - C-cyanomethanimine - 34	$C_2H_4O$ - acetaldehyde - 3,4
$C_2H_2N_2$ - hydrogen cyanide dimer - 35	$C_2H_4OS$ - mercaptoacetaldehyde - 58
$C_2H_2N_2O$ - cyanoformamide - 56	$C_2H_4O_2$ - methyl formate - 3
$C_2H_2N_4$ - $N_2$ -HCN-HCN trimer - 20	$C_2H_4O_2$ - acetic acid - 4
$C_2H_2O$ - ketene - 7,13,57	$C_2H_4O_2$ - formaldehyde dimer - 32
$C_2H_2O_2S$ - $SO_2$ -HCCH complex - 30	$C_2H_4O_2S$ - ethylene- $SO_2$ complex - 27
$C_2H_2O_3$ - acetylene-ozone complex - 18,32	$C_2H_4O_2S$ - mercaptoacetic acid - 58
$C_2H_3ArF$ - $CH_2CHF$ -Ar complex - 30	$C_2H_4O_3$ - 1,2,3-trioxolane - 18
$C_2H_3ClO$ - acetyl chloride - 40	$C_2H_4O_3$ - ethylene-ozone complex - 18,32
$C_2H_3ClO_2$ - methylchloroformate - 14	$C_2H_5F$ - ethyl fluoride - 23
$C_2H_3FO_2$ - methylfluoroformate - 14	$C_2H_5F_2OP$ - ethylphosphonic difluoride - 14
$C_2H_3F_3$ - 1,1,1-trifluoroethane - 4	$C_2H_5F_2P$ - ethyldifluorophosphine - 14

C <sub>2</sub> H <sub>5</sub> I	- ethyl iodide	- 13	C <sub>2</sub> S	- CCS radical	- 45
C <sub>2</sub> H <sub>5</sub> N	- vinylamine	- 7	C <sub>2</sub> Si	- silicon dicarbide	- 32
C <sub>2</sub> H <sub>5</sub> N	- ethanimine	- 13	C <sub>3</sub> FN	- fluorocyanooacetylene	- 52
C <sub>2</sub> H <sub>5</sub> N	- CH <sub>4</sub> -HCN complex	- 30	C <sub>3</sub> F <sub>2</sub> O	- difluoropropadienone	- 21
C <sub>2</sub> H <sub>5</sub> NO	- nitrosoethane	- 11,26	C <sub>3</sub> H	- C <sub>3</sub> H radical	- 45
C <sub>2</sub> H <sub>5</sub> NO	- acetamide	- 13,55	C <sub>3</sub> H	- cyclic C <sub>3</sub> H radical	- 45
C <sub>2</sub> H <sub>5</sub> NO <sub>2</sub>	- ethyl nitrite	- 13	C <sub>3</sub> HN	- cyanoacetylene	- 13,33,52
C <sub>2</sub> H <sub>5</sub> NO <sub>2</sub>	- N-methoxyformamide	- 4,15	C <sub>3</sub> HNO	- isocyanatoethyne	- 26
C <sub>2</sub> H <sub>5</sub> N <sub>3</sub>	- NH <sub>3</sub> -HCN-HCN trimer	- 20	C <sub>3</sub> HNO <sub>4</sub>	- HCN-(CO <sub>2</sub> ) <sub>2</sub> trimer	- 20
C <sub>2</sub> H <sub>6</sub>	- ethane	- 2	C <sub>3</sub> H <sub>2</sub> ClF	- 1-chloro-1-fluoroallene	- 38
C <sub>2</sub> H <sub>6</sub> FN	- dimethylfluoramine	- 9	C <sub>3</sub> H <sub>2</sub> ClN	- chloroacrylonitrile	- 13
C <sub>2</sub> H <sub>6</sub> FPS	- dimethylphosphonothioic fluoride	- 14	C <sub>3</sub> H <sub>2</sub> ClN	- N-chloropropargylimine	- 34
C <sub>2</sub> H <sub>6</sub> F <sub>2</sub> NP	- dimethylaminodifluorophosphine	- 14	C <sub>3</sub> H <sub>2</sub> F <sub>4</sub>	- cis-1,1,2,3- tetrafluorocyclopropane	- 18
C <sub>2</sub> H <sub>6</sub> N <sub>2</sub> O	- dimethyl nitrosamine	- 13	C <sub>3</sub> H <sub>2</sub> N <sub>2</sub>	- malononitrile	- 11
C <sub>2</sub> H <sub>6</sub> O	- dimethylether	- 13	C <sub>3</sub> H <sub>2</sub> N <sub>2</sub> O	- OC-HCN-HCN trimer	- 20
C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>	- dimethylperoxide	- 9	C <sub>3</sub> H <sub>2</sub> O	- cyclopropenone	- 33
C <sub>2</sub> H <sub>6</sub> O <sub>2</sub>	- ethylene glycol	- 9	C <sub>3</sub> H <sub>2</sub> O <sub>4</sub>	- 1,3-dioxolane-4,5-dione	- 10
C <sub>2</sub> H <sub>7</sub> FO	- dimethylether-HF complex	- 35	C <sub>3</sub> H <sub>3</sub> Cl	- propargyl chloride	- 44
C <sub>2</sub> H <sub>7</sub> ISi	- dimethylsilyliodide	- 31	C <sub>3</sub> H <sub>3</sub> ClO <sub>2</sub>	- vinyl chloroformate	- 54
C <sub>2</sub> H <sub>7</sub> N	- ethylamine	- 44	C <sub>3</sub> H <sub>3</sub> FO <sub>2</sub>	- vinyl fluoroformate	- 54
C <sub>2</sub> H <sub>7</sub> NO	- aminoethanol	- 13	C <sub>3</sub> H <sub>3</sub> F <sub>3</sub> O <sub>4</sub>	- formic acid-trifluoroacetic acid dimer	- 4
C <sub>2</sub> H <sub>7</sub> NS	- 2-aminoethane thiol	- 15	C <sub>3</sub> H <sub>3</sub> N	- acrylonitrile	- 17
C <sub>2</sub> H <sub>7</sub> P	- ethylphosphine	- 14	C <sub>3</sub> H <sub>3</sub> N	- C-ethynylmethanimine	- 26
C <sub>2</sub> NP	- C-cyanophosphaethyne	- 26	C <sub>3</sub> H <sub>3</sub> NS	- vinylisothiocyanate	- 15
C <sub>2</sub> N <sub>2</sub>	- isocyanogen	- 13,17,56	C <sub>3</sub> H <sub>4</sub>	- allene	- 4,13,52
C <sub>2</sub> O <sub>3</sub>	- OCO-CO complex	- 30	C <sub>3</sub> H <sub>4</sub>	- propyne	- 13

C <sub>3</sub> H <sub>4</sub> N <sub>2</sub>	- pyrazole - 13	C <sub>3</sub> H <sub>7</sub> F <sub>2</sub> P	- isopropyldifluorophosphine - 14
C <sub>3</sub> H <sub>4</sub> O	- acrolein - 56	C <sub>3</sub> H <sub>7</sub> I	- isopropyl iodide - 13
C <sub>3</sub> H <sub>4</sub> O	- H <sub>2</sub> CO-HCCH complex - 30	C <sub>3</sub> H <sub>7</sub> I	- propyl iodide - 23
C <sub>3</sub> H <sub>5</sub> Br	- cyclopropyl bromide - 13,17	C <sub>3</sub> H <sub>7</sub> N	- propenimine - 13
C <sub>3</sub> H <sub>5</sub> Cl	- H <sub>2</sub> C=C-CH <sub>2</sub> -HCl complex - 30	C <sub>3</sub> H <sub>7</sub> N	- cyclopropylamine - 13
C <sub>3</sub> H <sub>5</sub> ClS	- 3-chloropropylene sulfide - 49	C <sub>3</sub> H <sub>7</sub> N	- allylamine - 44
C <sub>3</sub> H <sub>5</sub> Cl <sub>3</sub> O	- 2,2,2-trichloroethylmethyl ether - 31	C <sub>3</sub> H <sub>7</sub> NO	- dimethylformamide - 13
C <sub>3</sub> H <sub>5</sub> F <sub>3</sub> O	- 2,2,2-trifluoroethylmethyl ether - 31	C <sub>3</sub> H <sub>7</sub> NO	- E-sp-propionaldoxime - 58
C <sub>3</sub> H <sub>5</sub> I	- 2-iodopropene - 13	C <sub>3</sub> H <sub>8</sub> Ge	- cyclopropylgermane - 44
C <sub>3</sub> H <sub>5</sub> N	- ethyl isocyanide - 13	C <sub>3</sub> H <sub>8</sub> O	- ethylmethylether - 23
C <sub>3</sub> H <sub>5</sub> N	- 2-methyl-3H-azirine - 34	C <sub>3</sub> H <sub>8</sub> OS	- 2-(methylthio)ethanol - 36
C <sub>3</sub> H <sub>5</sub> NO	- 2-azetidinone - 36	C <sub>3</sub> H <sub>8</sub> OS	- 1-mercaptopropanol - 36
C <sub>3</sub> H <sub>5</sub> NO	- ethyl cyanate - 58	C <sub>3</sub> H <sub>8</sub> Si	- dimethylsilaethylene - 20
C <sub>3</sub> H <sub>5</sub> NO <sub>2</sub>	- allyl nitrite - 6	C <sub>3</sub> H <sub>8</sub> Si	- cyclopropylsilane - 44
C <sub>3</sub> H <sub>6</sub>	- cyclopropane - 4	C <sub>3</sub> H <sub>9</sub> BF <sub>3</sub> N	- trimethylamine-BF <sub>3</sub> - 27
C <sub>3</sub> H <sub>6</sub>	- propene - 33	C <sub>3</sub> H <sub>9</sub> N	- isopropylamine - 13
C <sub>3</sub> H <sub>6</sub> ClN	- N-chloroazetidine - 34	C <sub>3</sub> H <sub>9</sub> NO <sub>2</sub> S	- trimethylamine-SO <sub>2</sub> - 27
C <sub>3</sub> H <sub>6</sub> O	- 2-methylvinylalcohol - 34	C <sub>3</sub> H <sub>9</sub> P	- isopropylphosphine - 14
C <sub>3</sub> H <sub>6</sub> OS	- thiopropionic acid - 58	C <sub>3</sub> H <sub>9</sub> P	- trimethylphosphine - 14
C <sub>3</sub> H <sub>6</sub> O <sub>2</sub> S	- trimethylene sulfone - 1	C <sub>3</sub> H <sub>9</sub> PS	- trimethylphosphine sulfide - 14
C <sub>3</sub> H <sub>6</sub> O <sub>3</sub>	- glyceraldehyde - 7	C <sub>3</sub> H <sub>10</sub> BrN	- (CH <sub>3</sub> ) <sub>3</sub> N-HBr complex - 30
C <sub>3</sub> H <sub>6</sub> O <sub>3</sub>	- methylglycolate - 15	C <sub>3</sub> H <sub>10</sub> ClN	- (CH <sub>3</sub> ) <sub>3</sub> N-HCl complex - 30
C <sub>3</sub> H <sub>7</sub> Br	- isopropyl bromide - 13	C <sub>3</sub> H <sub>10</sub> Si	- ethylmethysilane - 23
C <sub>3</sub> H <sub>7</sub> Cl	- isopropyl chloride - 13	C <sub>3</sub> H <sub>10</sub> Si	- propylsilane - 23
C <sub>3</sub> H <sub>7</sub> F	- isopropyl fluoride - 13,14	C <sub>3</sub> H <sub>12</sub> AlN	- trimethylaminoalane - 26
C <sub>3</sub> H <sub>7</sub> FO	- oxetane-HF complex - 35	C <sub>3</sub> N	- C <sub>3</sub> N radical - 45
		C <sub>3</sub> OS	- tricarbon oxide sulfide - 56

$C_4Br_2O_2$	- 1,2-dibromocyclobuten-3,4-dione	$C_4H_6N$	- 2-cyano-2-propyl radical	- 19	
	- 12				
$C_4HF$	- fluorodiacetylene	- 52	$C_4H_6N_2O_2$	- 2,5-piperazinedione	- 7
$C_4HNO_6$	- HCN-(CO <sub>2</sub> ) <sub>3</sub> tetramer	- 20	$C_4H_6O$	- 2,5-dihydrofuran	- 33
$C_4H_3ClS$	- 3-chlorothiophene	- 46	$C_4H_6O$	- 2,3-dihydrofuran	- 33
$C_4H_3F_3$	- 2,3,3-trifluorocyclobutene	- 27	$C_4H_6OS$	- tetrahydrothiophene-3-one	- 1
$C_4H_3IS$	- 2-iodothiophene	- 46	$C_4H_6O_2$	- meso-bisoxirane	- 49
$C_4H_3IS$	- 3-iodothiophene	- 46	$C_4H_6O_2$	- dl-bisoxirane	- 49
$C_4H_3N$	- 1-isocyano-prop-2-yne	- 26	$C_4H_6O_2$	- $\beta$ -butyrolactone	- 1
$C_4H_3NO$	- isocyanatopropyne	- 26	$C_4H_6O_2$	- cyclopropylcarboxylic acid	- 36
$C_4H_4ArO$	- argon-furan complex	- 27	$C_4H_6O_2$	- 2-butenoic acid	- 58
$C_4H_4FN$	- methylcyanoacetylene-HF complex	- 35	$C_4H_6O_2S$	- butadiene sulfone	- 1
$C_4H_4N_2O_2$	- uracil	- 7	$C_4H_6O_3$	- propylene carbonate	- 1
$C_4H_4O_3$	- cyclobutadiene ozonide	- 27	$C_4H_7F$	- 3-fluoro-2-methylpropene	- 14
$C_4H_5ArN$	- pyrrole-Ar	- 6,27	$C_4H_7F$	- trans-1-fluoro-2-butene	- 14
$C_4H_5F$	- 3-fluoro-1-butyne	- 38	$C_4H_7FO$	- butyryl fluoride	- 1
$C_4H_5F_3O_4$	- acetic acid-trifluoroacetic acid dimer	- 4	$C_4H_7N$	- pyrroline	- 4,15
$C_4H_5N$	- cyclopropyl cyanide	- 13	$C_4H_7N$	- n-propyl isocyanide	- 13
$C_4H_5N$	- allyl cyanide	- 33	$C_4H_7N$	- isopropyl isocyanide	- 13
$C_4H_5NO$	- cyclopropyl isocyanate	- 13	$C_4H_7NO$	- crotonaldoxime	- 58
$C_4H_5NO$	- allyl isocyanate	- 37	$C_4H_7NO$	- syn-2-methylacrylaldoxime	- 58
$C_4H_5NS$	- 4-methylthiazole	- 13	$C_4H_7NO$	- 2-pyrrolidone	- 58
$C_4H_5NS$	- 5-methylthiazole	- 13	$C_4H_8$	- cyclobutane	- 24
$C_4H_5N_3O$	- cytosine	- 7	$C_4H_8N_2O$	- N-nitrosopyrrolidine	- 6,13
$C_4H_6$	- 1,2-butadiene	- 38	$C_4H_8O$	- 2-methylpropanal	- 14
$C_4H_6ClO_2P$	- 4,5 dimethyl-2-chloro-1,3-dioxaphospholene	- 40	$C_4H_8O$	- cis-crotyl alcohol	- 15
			$C_4H_8O$	- ethyl vinyl ether	- 23
			$C_4H_8O$	- 1-butene oxide	- 27

C <sub>4</sub> H <sub>8</sub> O	- tetrahydrofuran	- 40	C <sub>5</sub> H <sub>4</sub> N <sub>2</sub> O <sub>2</sub>	- p-nitropyridine	- 1
C <sub>4</sub> H <sub>8</sub> O	- 3-methyloxetane	- 1	C <sub>5</sub> H <sub>5</sub> N	- cyanobicyclobutane	- 21
C <sub>4</sub> H <sub>8</sub> O	- 2-methyloxetane	- 1	C <sub>5</sub> H <sub>5</sub> NO	- 2-formyl pyrrole	- 13
C <sub>4</sub> H <sub>8</sub> O <sub>2</sub>	- 2-methyl-1,3-dioxolane	- 43	C <sub>5</sub> H <sub>5</sub> NO	- pyridine-N-oxide	- 13
C <sub>4</sub> H <sub>8</sub> S	- tetrahydrothiophene	- 40	C <sub>5</sub> H <sub>5</sub> NO <sub>2</sub> S	- pyridine-SO <sub>2</sub> complex	- 27
C <sub>4</sub> H <sub>9</sub> BF <sub>2</sub>	- t-butylborondifluoride	- 11	C <sub>5</sub> H <sub>5</sub> N <sub>5</sub>	- adenine	- 7
C <sub>4</sub> H <sub>9</sub> Br	- t-butyl bromide	- 13	C <sub>5</sub> H <sub>6</sub>	- cyclopentadiene	- 33
C <sub>4</sub> H <sub>9</sub> Cl	- 2-chloro-2-methylpropane	- 8	C <sub>5</sub> H <sub>6</sub> N <sub>2</sub> O <sub>2</sub>	- thymine	- 7
C <sub>4</sub> H <sub>9</sub> F	- 2-methyl-1-fluoropropane	- 14	C <sub>5</sub> H <sub>7</sub> ClO	- cyclobutylcarbonyl chloride	- 14
C <sub>4</sub> H <sub>9</sub> N	- pyrrolidine	- 13	C <sub>5</sub> H <sub>7</sub> NO	- furfurylamine	- 10
C <sub>4</sub> H <sub>9</sub> NO	- trimethylnitrosomethane	- 11	C <sub>5</sub> H <sub>8</sub>	- 1,4-pentadiene	- 24
C <sub>4</sub> H <sub>9</sub> NO	- morpholine	- 13	C <sub>5</sub> H <sub>8</sub> O	- cyclobutylcarboxaldehyde	- 14
C <sub>4</sub> H <sub>9</sub> NO	- butyraldoxime	- 58	C <sub>5</sub> H <sub>8</sub> O	- 1,4-pentadien-3-ol	- 36
C <sub>4</sub> H <sub>9</sub> NO	- isobutyraldoxime	- 58	C <sub>5</sub> H <sub>8</sub> O <sub>2</sub>	- $\alpha$ -methyl- $\gamma$ -butyrolactone	- 1
C <sub>4</sub> H <sub>9</sub> NO <sub>2</sub>	- t-butylnitrite	- 11	C <sub>5</sub> H <sub>9</sub> Cl	- chlorocyclopentane	- 14
C <sub>4</sub> H <sub>9</sub> NSi	- trimethylsilyl cyanide	- 30	C <sub>5</sub> H <sub>9</sub> N	- t-butyl cyanide	- 16
C <sub>4</sub> H <sub>10</sub>	- butane	- 13	C <sub>5</sub> H <sub>9</sub> N	- valeronitrile	- 26
C <sub>4</sub> H <sub>10</sub> Ge	- cyclobutylgermane	- 14	C <sub>5</sub> H <sub>9</sub> N	- t-butyl isocyanide	- 30
C <sub>4</sub> H <sub>10</sub> N <sub>2</sub>	- (CH <sub>3</sub> ) <sub>3</sub> N-HCN complex	- 30	C <sub>5</sub> H <sub>9</sub> NO	- cyclopentanone oxime	- 58
C <sub>4</sub> H <sub>10</sub> O <sub>2</sub>	- 3-methoxypropanol	- 36	C <sub>5</sub> H <sub>9</sub> NO	- 2-piperidone	- 58
C <sub>4</sub> H <sub>10</sub> Si	- cyclobutylsilane	- 44	C <sub>5</sub> H <sub>10</sub> FN	- (CH <sub>3</sub> ) <sub>3</sub> CCN-HF complex	- 30
C <sub>5</sub> H	- C <sub>5</sub> H radical	- 45	C <sub>5</sub> H <sub>10</sub> FN	- t-butyl cyanide-HF complex	- 35
C <sub>5</sub> H <sub>3</sub> N	- 1-cyano-but-3-en-1-yne	- 26	C <sub>5</sub> H <sub>10</sub> N <sub>2</sub> O	- N-nitrosopiperidine	- 6
C <sub>5</sub> H <sub>3</sub> N	- 1-cyano-but-1-en-3-yne	- 26	C <sub>5</sub> H <sub>10</sub> O	- pivaldehyde	- 11,27
C <sub>5</sub> H <sub>3</sub> NO	- 2-cyanofuran	- 13	C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>	- 2,2-dimethyl-1,3-dioxolane	- 43
C <sub>5</sub> H <sub>4</sub> N <sub>2</sub>	- diazocyclopentadiene	- 58	C <sub>5</sub> H <sub>10</sub> Si	- trimethylsilylacetylene	- 13,26
C <sub>5</sub> H <sub>4</sub> N <sub>2</sub>	- pyrrole-3-carbonitrile	- 58	C <sub>5</sub> H <sub>11</sub> N	- piperidine	- 13

C <sub>5</sub> H <sub>11</sub> N - (CH <sub>3</sub> ) <sub>3</sub> N-HCCH complex	- 30	C <sub>7</sub> H <sub>4</sub> FNO - m-fluorophenyl isocyanate	- 1
C <sub>6</sub> H <sub>4</sub> Cl <sub>2</sub> - 1,2-dichlorobenzene	- 13	C <sub>7</sub> H <sub>4</sub> F <sub>2</sub> O - 3-fluorobenzoyl fluoride	- 10
C <sub>6</sub> H <sub>4</sub> Cl <sub>2</sub> - 1,3-dichlorobenzene	- 13	C <sub>7</sub> H <sub>4</sub> F <sub>2</sub> O - 4-fluorobenzoyl fluoride	- 10
C <sub>6</sub> H <sub>4</sub> N <sub>2</sub> S - 1,2,3-benzothiadiazole	- 58	C <sub>7</sub> H <sub>4</sub> OS - benzothiet-2-one	- 58
C <sub>6</sub> H <sub>5</sub> BF <sub>2</sub> - phenyldifluoroborane	- 13	C <sub>7</sub> H <sub>5</sub> FO - o-fluorobenzaldehyde	- 1
C <sub>6</sub> H <sub>5</sub> Cl - chlorobenzene	- 13	C <sub>7</sub> H <sub>5</sub> FO - m-fluorobenzaldehyde	- 1
C <sub>6</sub> H <sub>5</sub> F - fluorobenzene	- 4	C <sub>7</sub> H <sub>5</sub> FO - p-fluorobenzaldehyde	- 1
C <sub>6</sub> H <sub>5</sub> NO - 3-formylpyridine	- 13	C <sub>7</sub> H <sub>5</sub> NO - phenyl cyanate	- 58
C <sub>6</sub> H <sub>5</sub> NO - 4-formylpyridine	- 13	C <sub>7</sub> H <sub>6</sub> OS - thiobenzoic acid	- 58
C <sub>6</sub> H <sub>5</sub> NO <sub>2</sub> - nitrobenzene	- 13	C <sub>7</sub> H <sub>6</sub> O <sub>2</sub> - tropolone	- 52
C <sub>6</sub> H <sub>5</sub> NO <sub>3</sub> - o-nitrophenol	- 13	C <sub>7</sub> H <sub>6</sub> O <sub>2</sub> - benzoic acid	- 58
C <sub>6</sub> H <sub>6</sub> - benzene	- 4,13	C <sub>7</sub> H <sub>6</sub> O <sub>3</sub> - salicylic acid	- 58
C <sub>6</sub> H <sub>6</sub> O <sub>2</sub> S - benzene-SO <sub>2</sub> complex	- 27	C <sub>7</sub> H <sub>7</sub> Cl - benzyl chloride	- 58
C <sub>6</sub> H <sub>7</sub> N - aniline	- 13	C <sub>7</sub> H <sub>7</sub> X - benzyl halides	- 54
C <sub>6</sub> H <sub>7</sub> NO - 2-picoline-N-oxide	- 13	C <sub>7</sub> H <sub>8</sub> - quadricyclane	- 4
C <sub>6</sub> H <sub>8</sub> Ge - phenyl germane	- 15	C <sub>7</sub> H <sub>8</sub> O - anisole	- 58
C <sub>6</sub> H <sub>8</sub> O - 3-cyclohexenone	- 15	C <sub>7</sub> H <sub>8</sub> O - benzyl alcohol	- 58
C <sub>6</sub> H <sub>9</sub> NO - 2-cyclohexen-1-one oxime	- 58	C <sub>7</sub> H <sub>10</sub> - 3-cyclohexene-1-methylene	- 15
C <sub>6</sub> H <sub>9</sub> NO <sub>2</sub> - 1-nitrocyclohexene	- 1	C <sub>7</sub> H <sub>10</sub> O - 3-methyl-2-cyclohexen-1-one	- 31
C <sub>6</sub> H <sub>9</sub> NSi - trimethylsilylcyanethyne	- 26	C <sub>7</sub> H <sub>12</sub> O - 8-oxabicyclo[3.2.1]octane	- 27
C <sub>6</sub> H <sub>10</sub> S - 7-thiabicyclo[2.2.1]heptane	- 24	C <sub>7</sub> H <sub>12</sub> S - 1-thiaspiro[2.5]octane	- 43
C <sub>6</sub> H <sub>11</sub> NO - 1-formylpiperidine	- 6	C <sub>7</sub> H <sub>13</sub> N - 1-azabicyclo[2.2.2]octane	- 30
C <sub>6</sub> H <sub>11</sub> NO - cyclohexanone oxime	- 58	C <sub>8</sub> H <sub>7</sub> F - o-fluorostyrene	- 1
C <sub>6</sub> H <sub>12</sub> - cyclohexane	- 4	C <sub>8</sub> H <sub>7</sub> F - m-fluorostyrene	- 1
C <sub>7</sub> H <sub>3</sub> F <sub>2</sub> N - 2,3-difluorobenzonitrile	- 37	C <sub>8</sub> H <sub>7</sub> N - benzyl cyanide	- 6,58
C <sub>7</sub> H <sub>4</sub> BrN - 3-bromobenzonitrile	- 37	C <sub>8</sub> H <sub>8</sub> O - acetophenone	- 58
C <sub>7</sub> H <sub>4</sub> ClN - 2-chlorobenzonitrile	- 58	C <sub>8</sub> H <sub>8</sub> O <sub>2</sub> - p-anisaldehyde	- 6,54

$C_8H_9NSi$ - trimethylsilylcyanobutadiyne - 26	$Cl_2Si$ - dichlorosilylene - 24,53
$C_8H_{14}O$ - 4-ethylcyclohexanone - 31	$CsH$ - cesium hydride - 44
$C_8H_{14}O_2$ - 1,4-dioxaspiro[4.5]decane - 43	$FHN_2O$ - hydrogen fluoride-nitrous oxide complex - 5,28
$C_8H_{14}S_2$ - 1,4-dithiaspiro[4.5]decane - 43	$FH_3Si$ - silyl fluoride - 13
$C_9H_{14}O$ - bicyclo[3.3.1]nonan-9-one - 31	$FI$ - iodine fluoride - 44
$C_{10}H_{12}O$ - p-isopropyl benzaldehyde - 6	$F_3HSi$ - trifluorosilane - 52
$C_{10}H_{15}Cl$ - 1-chloroadamantane - 30	$F_3H_2OP$ - water-PF <sub>3</sub> complex - 27
$C_{12}H_9F$ - 4-fluorobiphenyl - 58	$F_3KrP$ - krypton-PF <sub>3</sub> complex - 6,27
$CaS$ - calcium sulfide - 45	$F_3NeP$ - neon-PF <sub>3</sub> complex - 27
$ClFH_2$ - hydrogen fluoride hydrogen chloride complex - 32	$F_3OP$ - phosphoryl trifluoride - 4,52
$ClFO_3$ - perchloryl fluoride - 13	$F_3Si$ - silicon trifluoride - 24,53
$ClFSi$ - chlorofluorosilylene - 24	$F_6OS$ - fluoroxysulfurpentafuoride - 9
$ClF_3Si$ - chlorotrifluorosilane - 8	$GaH$ - gallium hydride - 44
$ClGeH_3$ - germyl chloride - 13,33	$GeH_3N_3$ - germylazide - 14
$ClHO_2S$ - SO <sub>2</sub> -HCl complex - 30	$HIn$ - indium hydride - 44
$ClH_3O$ - H <sub>2</sub> O-HCl complex - 30	$HNO$ - nitroxyl - 24
$ClH_4N$ - H <sub>3</sub> N-HCl complex - 30	$HN_3$ - hydrazoic acid - 13,17
$ClH_4P$ - H <sub>3</sub> P-HCl complex - 30	$HPb$ - lead hydride - 44
$ClIn$ - indium monochloride - 25	$HRb$ - rubidium hydride - 44
$ClMg$ - magnesium monochloride - 33	$HS_2$ - HS <sub>2</sub> radical - 45
$ClOS$ - ClSO radical - 24	$HSr$ - strontium hydride - 44
$ClSi$ - silicon monochloride - 53	$HTl$ - thallium hydride - 44
$Cl_2OS$ - thionyl chloride - 13	$HZn$ - zinc hydride - 44
$Cl_2O_2$ - chlorine peroxide - 39	$H_2N$ - NH <sub>2</sub> radical - 45
$Cl_2O_2S$ - sulfuryl chloride - 13	$H_2NO$ - H <sub>2</sub> NO radical - 45
$Cl_2S$ - sulfur dichloride - 13	$H_2NS$ - H <sub>2</sub> NS radical - 45
	$H_2O$ - water - 33

$H_2O_2$	- hydrogen peroxide	- 13	$O_2S$	- sulfur dioxide	- 53
$H_2O_3S$	- water-sulfur dioxide complex	- 32			
$H_2O_4$	- ozone-water complex	- 18,32			
$H_2S_2$	- hydrogen disulfide	- 57			
$H_2S_3$	- hydrogen trisulfide	- 57			
$H_2S_4$	- hydrogen tetrasulfide	- 57			
$H_3IS$	- $H_2S$ -HI complex	- 30			
$H_3N$	- ammonia	- 13,39,47,48			
$H_3P$	- phosphine	- 44			
$H_3Sb$	- stibine	- 12			
$H_4IP$	- $H_3P$ -HI complex	- 30			
$H_4N_2$	- hydrazine	- 55			
$H_4O_2$	- water dimer	- 32			
$H_4Si$	- silane	- 13			
$H_4Sn$	- stannane	- 50			
KO	- potassium monoxide	- 24			
LiO	- lithium monoxide	- 24			
MgS	- magnesium sulfide	- 45			
N	- nitrogen atom	- 33			
NO	- nitric oxide	- 57			
NS <sub>2</sub>	- nitrogen disulfide	- 2			
N <sub>2</sub> O	- nitrous oxide	- 13			
N <sub>2</sub> O <sub>3</sub>	- dinitrogen trioxide	- 11			
N <sub>2</sub> O <sub>5</sub>	- nitrogen pentoxide	- 39			
NaO	- sodium monoxide	- 24			
ORb	- rubidium monoxide	- 24			
OY	- yttrium monoxide	- 25			