Infrared Molecular Absorption Cross-sections

The folder IR-XSect contains files of infrared cross-sections. The definition and units have been described in articles about the HITRAN compilation. Each molecule is placed in a single file. Within that file are sets of temperature and pressure pairs. The sets have a header that provides information to programs reading the data and also points to a reference for that observation. The sets contain absorption cross-sections (ten to a line from left to right) that are in equal wavenumber (cm⁻¹) increments, and the intervals can be determined by the minimum and maximum wavenumber and the number of points, namely

$$\Delta v = \frac{v_{\text{max}} - v_{\text{min}}}{npts - 1}$$

where v_{max} is the maximum (final) wavenumber of the set, v_{min} is the minimum (initial) wavenumber of the set, and *npts* is the number of points in the set. The format of the header is given below.

	Cross-section Header Format													
	Chemical symbol	Waver Min	number Max	No. Pts.	Temp [K]	Press [torr]	Max X-section	Res.	Common Name	Not Brused	r Ref No			
20		10	10	7	7	6	10	5	15	4 3	3			
	10	20	30	40	50		60	70	80	90				

Note: **Chemical Symbol** is right adjusted; **Res**. is resolution in cm⁻¹ for FTS measurements, and **Br** indicates the broadening gas, such as air.

The \Supplemental folder contains two types of files: (1) some older, redundant cross-section data that have nonetheless been retained, and (2) data that have some small experimental negative cross-sections that were zeroed out for the files in the main directory (some users prefer these files as they do not introduce any bias). The extension for file names is ".xsc" for the former, and ".alt" for the latter.

A summary of the molecules represented with their temperature and pressures ranges and spectral coverage is given in the table on the following pages:

Summary of Molecules Represented by Infrared Cross-section Data in HITRAN

Molecule	Common Name	Temperature Range (K)	Pressure Range (torr)	Number of T,P sets	Spectral Coverage (cm ⁻¹)				
SF ₆	Sulfur hexafluoride	180-295	20-760	32	925-955				
grava		189-297	0-117	25	750-830				
ClONO ₂	Chlorine nitrate	189-297 213-296	0-117 0	25 2	1260-1320 1680-1790				
CCl ₄	Carbon tetrachloride	208-297	8-760	32	750-812				
N_2O_5	Dinitrogen pentoxide	205-293	0	5	540-1380				
HNO ₄	Peroxynitric acid	220	0	1	780-830				
C_2F_6	Hexafluoroethane,	181-296	25-760	43	1061-1165				
	CFC-116	181-296 190-296	25-760 8-760	43 55	1220-1285 810-880				
CCl ₃ F	CFC-11	190-296	8-760	55	1050-1120				
CCl_2F_2	CFC-12	190-296 190-296	8-760 8-760	52 52	850-950 1050-1200				
CCIF ₃	CFC-13	203-293 203-293	0	6 6	765-805 1065-1140				
CCIF ₃	CFC-13	203-293	0	6	1170-1235				
CF ₄	CFC-14	180-296	8-761	55	1250-1290				
$C_2Cl_2F_3$	CFC-113	203-293 203-293	0	6 6	780-995 1005-1232				
		203-293	0	6	815-860				
$C_2Cl_2F_4$	CFC-114	203-293 203-293	0	6 6	870-960 1030-1067				
		203-293	0	6	1095-1285				
C ₂ ClF ₅	CFC-115	203-293 203-293	0	6 6	955-1015 1110-1145				
	HOEG M	203-293	0	6	1167-1260				
CHCl ₂ F	HCFC-21	296 181-297	0-765	1 29	785-840 760-860				
CHCIF ₂	HCFC-22	181-296	22-761	31	1070-1195				
_		253-287 253-287	0	3	1060-1210 1275-1380				
CHCl ₂ CF ₃	HCFC-123	253-287	0	3	740-900				
		253-287 287	0	3	1080-1450 675-715				
CHCIFCF ₃	HCFC-124	287	0	1	790-920				
		287 253-287	0	3	1035-1430 710-790				
CH ₃ CCl ₂ F	HCFC-141b	253-287	0	3	895-1210				
		253-287 253-287	0	3	1325-1470 650-705				
CH ₃ CCIF ₂	HCFC-142b	253-287	0	3	875-1265				
		253-287 253-287	0	3	1360-1475 695-865				
CHCl ₂ CF ₂ CF ₃	HCFC-225ca	253-287	0	3	1010-1420				
CCIF ₂ CF ₂ CHCIF	HCFC-225cb	253-287 203-297	0 0-750	3 17	715-1375 995-1236				
CH_2F_2	HFC-32	203-297	0-750	17	1385-1475				
CHF ₂ CF ₃	HFC-125	287 287	0	1 1	700-745 840-890				
		287	0	1	1060-1465				
CHF ₂ CHF ₂	HFC-134	203-297 253-287	0-750	9	600-1700 815-865				
CFH ₂ CF ₃	HFC-134a	190-296	20-760	32	1035-1130				
CF 112CF 3		190-296 253-287	20-760 0	33 3	1135-1340 935-1485				
		203-297	0-750	9	580-630				
CF ₃ CH ₃	HFC-143a	203-297 203-297	0-750 0-750	9 9	750-1050 1100-1500				
	*****	253-287	0	3	840-995				
CH ₃ CHF ₂	HFC-152a	253-287 253-287	0 0	3 3	1050-1205 1320-1490				
		213-323	760	5	599-624				
SF ₅ CF ₃	Trifluoromethyl sulfur	213-323 213-323	760 760	5 5	676-704 740-766				
SF5CF3	pentafluoride	213-323 213-323	760 760	5 5	860-920 1150-1280				
		213-323	760	5	1280-2600				
New or modified data added after the HITRAN2004 edition									
CH ₃ C(O)OONO ₂	PAN (Peroxyacetal	<u>uded after th</u> 295	0.08	1	550-1450				
CII3C(U)UUNU2	nitrate)	295	0.08	1 2	1650-1901				
		276-324 276-324	760 760	3	624-784 867-1159				
CH ₃ CN	Acetonitrile (methyl cyanide)	276-324 276-324	760 760	3	1175-1687 2217-2343				
	(276-324	760	3	2786-3261				
C ₆ H ₆	Benzene	276-324 278-323	760 760	3	3881-4574 600-6500				
CHF ₂ CF ₃	HFC-125	203-293	0-600	16	494-1503				

Note: These data are in the main directory. Additional redundant data for CFC-11, CFC-12, HFC-125, and HFC-143a are stored in a supplemental sub-directory.