

Theoretical study of the OH-initiated atmospheric oxidation mechanism of perfluoro methyl vinyl ether, $\text{CF}_3\text{OCF}=\text{CF}_2$

– Supporting Information –

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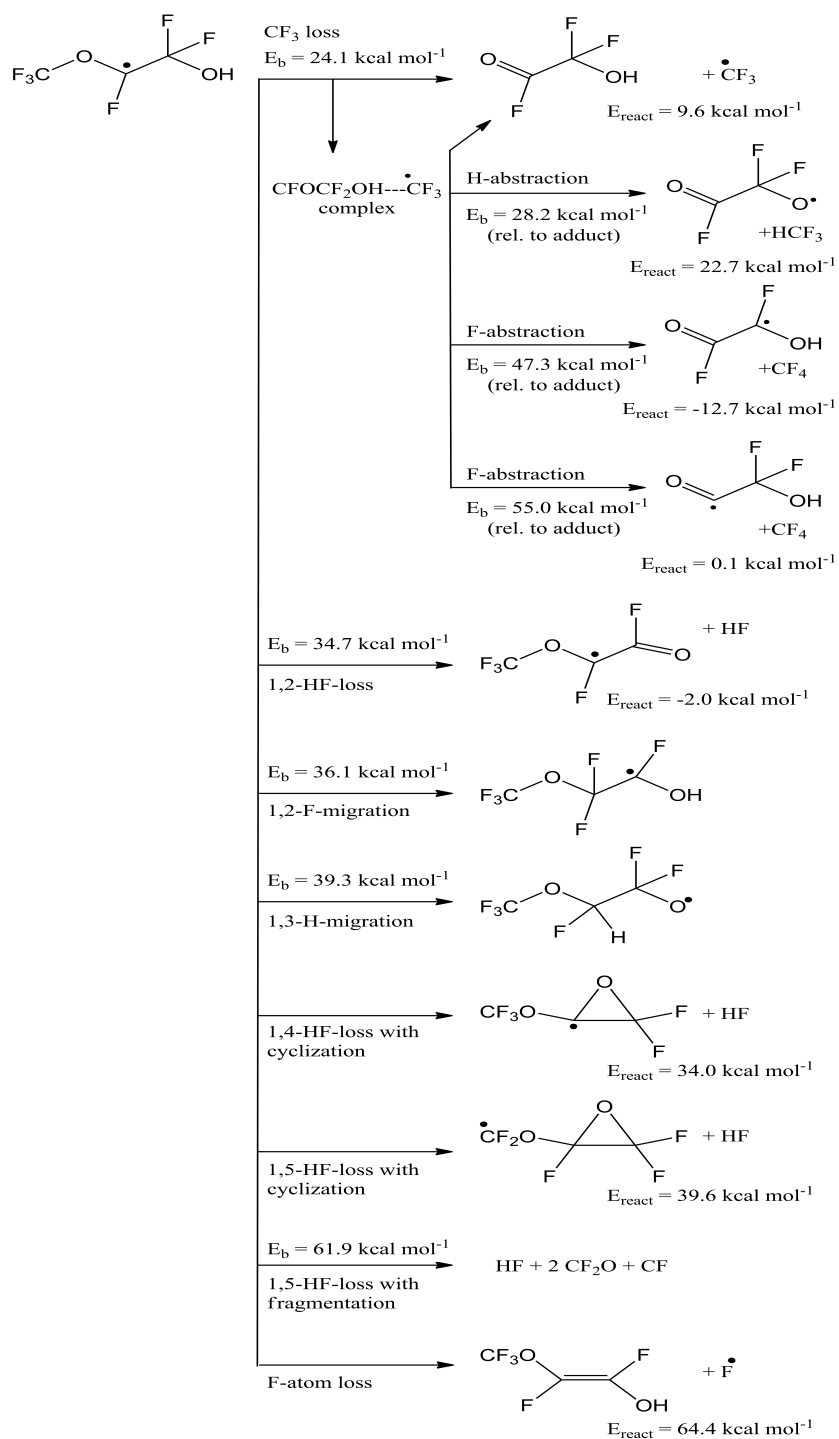


Figure SI - 1: All pathways considered for the main adduct, at the M06-2X/aug-cc-pVTZ level of theory.

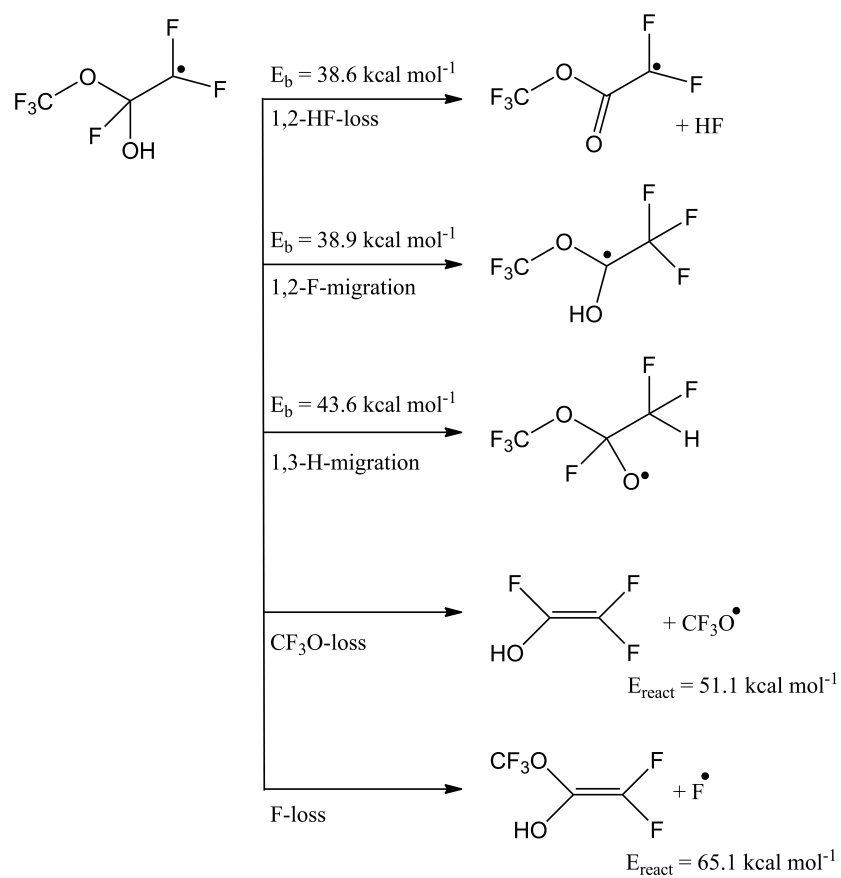


Figure SI - 2: All pathways considered for the minor adduct, at the M06-2X/aug-cc-pVTZ level of theory.

CCSD(T) *ab initio* calculations

To assess the robustness of the M06-2X/aug-cc-pVTZ calculations, we compared these against single point energy calculations at the CCSD(T)/6-311++G(d,p), CCSD(T)/aug-cc-pVDZ, and CCSD(T)-F12a/aug-cc-pVDZ levels of theory (see Table SI-1); the raw data on all compounds discussed in the paper are listed further in this supporting information. The CCSD(T) calculations were obtained using Gaussian-09, the CCSD(T)-F12 results with Molpro 2010.1. While these calculations are based on coupled cluster theory, they are not *a priori* more reliable than the M06-2X calculations, as the CC predictions still carry uncertainties of several kcal/mol due to the small size of the basis set, which limits the flexibility in accounting for electron correlation. 6-311++G(d,p) and aug-cc-pVDZ basis sets are similar in size, where the first basis set is rather small but reasonably balanced due to its use of split valence orbitals, while the latter has ~20% more Gaussian primitives and is part of a systematic series of basis sets that allows for extrapolation to the complete basis set. For basis sets of this size, errors on absolute CCSD(T) atomization energies are reported of up to several tens of kcal/mol (see e.g. D. Feller, K. A. Peterson, J. Chem. Phys., 2007, 126, 114105 ; K. A. Peterson, D. Feller, D. A. Dixon, Theor. Chem. Acc., 2012, 131, 1079). Cancellation of error leads to *relative* energies that are significantly more reliable, but non-negligible uncertainties remain especially for free products/reactants or for barrier heights for bimolecular reactions, where the cancellation of error is least effective due to significant differences in bond types and bonding partners on either side of the chemical equation. For example, the predicted relative energy of some sets of products differs as much as 7 kcal mol⁻¹ among the CC levels of theory, despite using the same geometry and a similarly sized basis set. For the unimolecular reactions listed in Table SI-1, the average unsigned difference (MUE) between the M06-2X barrier height and the CCSD(T)/6-311++G(d,p) barrier height is 2.5 kcal/mol, against CCSD(T)/aug-cc-pVDZ it is 2.1 kcal/mol. In these values, the “CF₃OC*FCF₂OH 1,5-HF-loss with fragmentation” reaction (see figure SI-1) has a disproportionate impact; it leads to 4 fragments in a single decomposition TS and thus represents an extreme change in bonding. Without this reaction M06-2X and CCSD(T) with a small basis set agree on average within ~2 kcal/mol on the barrier heights.

Preferably one would use larger basis sets in the coupled cluster calculations, such as aug-cc-pVTZ and beyond, and extrapolate the energies to the basis set limit. Unfortunately, with up to 12 non-hydrogen atoms and open shell wave functions, these calculations are beyond our computational capabilities. CCSD(T)-F12 explicitly correlated methodologies converge significantly faster to the complete basis set limit, but CCSD(T)-F12/aug-cc-pVDZ still shows sizable uncertainties even when these are up to an order of magnitude smaller than for traditional CCSD(T) calculations with a similar small basis set (See e.g. Peterson et al., 2012). Also, such calculations again strain our computational resources to the maximum, and we had to limit our calculations to the smaller fragments, and a few select reactions. Composite methods, such as G2, G3SX or CBS-QB3, have reported uncertainties of a few kcal/mol, and are likewise not a sufficient improvement over the current set of data to warrant the computational cost. Reaching chemical accuracy, i.e. with uncertainties of 1 kcal/mol, is thus computationally too costly, while more tractable calculations do not result in a meaningful increase in

the *a priori* reliability compared to the results reported here. As such, we consider higher level calculations than those considered here or available in the literature to be beyond our reach at this time.

As argued in the main paper, however, our analysis of the reaction mechanism driving the $\text{CF}_3\text{OCF}=\text{CF}_2 + \text{OH}$ reaction is not sensitive even to large uncertainties of several kcal/mol on the energy predictions, due to the sizable energy differences predicted for competing channels. Comparing M06-2X and CCSD(T) results for the critical points on the potential energy surface, we conclude that :

- The relative energy of the addition transition states indicates that the OH-addition to the outer carbon has an energetic advantage, and is expected to be the main entrance channel with a contribution of $80\pm 15\%$ (see below).
- The energy of the addition transition states is below the energy of the separated reactants, allowing for a negative temperature-dependence of the rate coefficient (see below). Protruding barriers for addition would lead to rate coefficients that are significantly lower than the experimental values, and are considered unlikely.
- The CF_3 loss channel is the only reaction channel for the adducts that has a sufficiently low barrier to allow for prompt decomposition at 1 atm. The HF elimination channel has a significantly higher barrier and concomitantly a negligible contribution.
- The reaction mechanism proposed earlier by Li et al. and Mashino et al., i.e. leading to formation of glyoxal directly by a process of HF elimination and C–O bond scission in the alkoxy radicals subsequently formed, is not viable as several steps in this mechanism have competing reactions with significantly lower energy barriers.
- The products of the reaction are predicted to be (per)fluorinated methylformate, glycolaldehyde, and formaldehyde. Direct glyoxal formation is unsupported by the available theoretical data, and glyoxal is predicted to be a secondary product.
- An *a priori* estimation of the yield of prompt decomposition has only limited reliability due to the uncertainties on the collisional energy transfer parameters. A yield of $\sim 40\%$ prompt decomposition by CF_3 loss from the $\text{CF}_3\text{OC}\cdot\text{FCF}_2\text{OH}$ radical can be reproduced using physically viable collisional parameters. A significant yield of HF-loss from the initial adducts requires physically highly unlikely collision characteristics, as well as the removal of the CF_3 loss channel from the calculations.

These main conclusions of the paper are thus supported in full within the uncertainties of the theoretical calculations available at this time, *i.e.* in the current paper and in the cited literature.

Table SI-1 : Barrier heights or (in brackets) reaction energies for critical reactions in the CF₃OCF₂ + OH reaction. The table is subdivided in sections corresponding to Figures 1 through 3 (main text), Figures SI-1 and SI-2 (supporting information), and miscellaneous reactions. The table compares the energies obtained using M06-2X/aug-cc-pVTZ (designated as M06-2X), CCSD(T)/6-311+G(d,p)//M06-2X (CC/311), CCSD(T)/aug-cc-pVDZ//M06-2X (CC/aVDZ), and ROHF-UCCSD(T)-F12a/aug-cc-pVDZ//M06-2X (CCF12/aVDZ).

Reaction	M06-2X	CC/311	CC/aVDZ	CCF12/aVDZ	
CF ₃ OC*FCFO : CF ₃ loss	30.1	25.1	25.8	29.8	Figure 1
CF ₃ OCF(O*)CFO : OCF ₃ loss	17.8	13.8	14.1		
CF ₃ OCF(O*)CFO : F-loss	24.2	24.9	24.4		
CF ₃ OCF(O*)CFO : FCO loss	0.29	2.0	3.7		
CFOCF ₂ OH : 1,2-HF-loss	45.5	44.6	43.0	48.0	
CF ₃ OCF=CF ₂ + *OH → CF ₃ OC*F-CF ₂ OH	-2.1	-0.03	-2.1		Figure 2
CF ₃ OC*FCF ₂ OH : CF ₃ -loss	24.1	21.4	22.1		
CF ₃ OC*FCF ₂ OH : 1,2-HF-loss	34.7	35.6	34.0		
CF ₃ OC*F(O*)CF ₂ OH : F-loss	26.4	20.6	20.4		
CF ₃ OC*F(O*)CF ₂ OH : OCF ₃ loss	19.5	15.1	15.2		
CF ₃ OC*F(O*)CF ₂ OH : CF ₂ OH loss	< 1.5	< 1.5	< 1.5		
CF ₃ OCF=CF ₂ + *OH → CF ₃ OCFOH-C*F ₂	-1.5	1.2	-1.3		Fig. 3
CF ₃ OCFOHC*F ₂ : 1,2-HF-loss	38.6	39.6	37.9		
CF ₃ OC*FCF ₂ OH : CF ₃ loss	24.1	21.4	22.1		Figure SI-1
CF ₃ OC*FCF ₂ OH : 1,2-HF-loss	34.7	35.6	34.0		
CF ₃ OC*FCF ₂ OH : 1,2-F-migration	36.1	37.1	35.7		
CF ₃ OC*FCF ₂ OH : 1,3-H-migration	39.3	37.5	36.7		
CF ₃ OC*FCF ₂ OH : 1,4-HF-loss with cyclization	(34.0)	(30.0)	(33.9)		
CF ₃ OC*FCF ₂ OH : 1,5-HF-loss with cyclization	(39.6)	(35.6)	(39.8)		
CF ₃ OC*FCF ₂ OH : 1,5-HF-loss with fragmentation	61.9	53.6	55.6		
CF ₃ OC*FCF ₂ OH : F-atom loss	(64.4)	(57.3)	(60.5)		
CF ₃ OCFOHC*F ₂ : 1,2-HF-loss	38.6	39.6	37.9		Figure SI-2
CF ₃ OCFOHC*F ₂ : 1,2-F-migration	38.9	40.6	38.4		
CF ₃ OCFOHC*F ₂ : 1,3-H-migration	43.6	41.9	41.0		
CF ₃ OCFOHC*F ₂ : CF ₃ O-loss	(51.1)	(46.4)	(50.1)		
CF ₃ OCFOHC*F ₂ : F-atom loss	(65.1)	(57.2)	(61.0)		
CFOCF ₂ O* : C-C scission	0.97	3.1	4.3	2.4	Misc. React.
CFOCF ₂ OH---H ₂ O : 1,2-HF-loss	14.4	15.6	12.8	15.0	
CFOCF ₂ OH + *CF ₃ -> CF ₄ + CFOC*FOH	47.3	47.5	41.6		
CFOCF ₂ OH + *CF ₃ -> HCF ₃ + CFOCF ₂ O*	28.2	20.6	21.8		
CFOCF ₂ OH + *CF ₃ -> CF ₄ + OC*CF ₂ OH	55.0	47.2	47.8		

Rate coefficient predictions

Three experimental determination for the rate coefficient of the $\text{CF}_3\text{OCF}=\text{CF}_2 + \text{OH}$ reaction are available in the literature :

$$\text{Li et al. : } k(253\text{-}348\text{K}) = (6.41 \pm 0.82) \times 10^{-11} \exp((-868 \pm 40)/T) \text{ cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$$

$$\text{Tokuhashi et al. : } k(250\text{-}430\text{K}) = (1.01 \pm 0.04) \times 10^{-12} \exp((320 \pm 10)/T) \text{ cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$$

$$\text{Mashino et al. : } k(296\text{K}) = (2.6 \pm 0.3) \times 10^{-12} \text{ cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$$

Near room temperature these rate coefficients differ by less than 30%, but the opposite sign of the temperature dependence for the expressions by Li et al. versus Tokuhashi et al. leads to larger differences at higher and lower temperatures.

The accurate prediction of the absolute rate coefficient of the $\text{CF}_3\text{OCF}=\text{CF}_2 + \text{OH}$ reaction is difficult, as the value depends critically on the absolute barrier heights of the entrance transition states, on the interplay between the initial complex formation and the OH addition step, and on the description of the degrees of freedom for internal rotation, non-harmonic vibration, etc. along the addition process. Such complex calculations are beyond the scope of the current paper. Hence, we will not attempt a highly accurate prediction of the rate coefficient, but rather limit ourselves to assess the temperature dependence, and the relative contribution of the different entrance channels leading to addition on the inner or outer carbon atom.

To this end we performed simplified canonical transition state theory calculations (CTST), where we assume that (i) the formation and redissociation of the pre-reactive complex is fast enough compared to the addition reaction to ensure canonical equilibrium between the free reactants and the complex; (ii) that the impact of anharmonicities, internal rotation, etc. is sufficiently similar between all of the reactants, complex formation TS, complex, and the addition transition states, to mostly cancel out relative to a harmonic oscillator treatment; (iii) that the predicted relative energy of the two addition transition states is sufficiently accurate; (iv) that tunneling is negligible; and (v) that the addition reaction is irreversible.

If the approximations above hold, the impact of the pre-reactive complex cancels out of the rate equations, and the absolute rate coefficient can be predicted based on the rigid-rotor, harmonic oscillator approximation for reactants and addition transition states, where we use the M06-2X/aug-cc-pVTZ rovibrational and energetic characteristics. For OH we use the experimental spin-orbit splitting value of 126.23 cm^{-1} :

$$k(T) = \frac{k_b T}{h} \times \frac{Q^\ddagger(T)}{Q_{\text{CF}_3\text{OCFCF}_2(T)} \times Q_{\text{OH}}(T)} \times \exp\left(\frac{-E_b}{k_b T}\right)$$

Figure SI-3 shows the result of this calculation, compared to the available experimental data. The submerged barriers as predicted by M06-2X//aug-cc-pVDZ and CCSD(T)/aVDZ lead to a negative temperature dependence in fair agreement with the results of Tokuhashi et al., and in strong disagreement with the results of Li et al. The excellent level of agreement of the absolute rate

coefficients between theory and experiment must be considered fortuitous, in view of the uncertainties on the underlying data and the methodology used. Raising the barrier heights by only 0.4 kcal/mol is sufficient to shift our rate coefficient predictions below the $k(T)$ values of Tokuhashi et al. and Mashino et al. To obtain a temperature-neutral rate coefficient, the barriers must be raised by 1.5 kcal/mol above the M06-2X values, while a positive temperature dependence similar to Li et al. requires barriers higher by 3.5 kcal/mol. The predicted $k(T)$ in these latter two cases are found to be a factor of 10 to 350 below the experimental rate coefficient, respectively. Using the CCSD(T)/6-311++G(d,p) addition barrier heights, which are 2 to 2.7 kcal/mol higher than the M06-2X and CCSD(T)/aVDZ results, yields a nearly temperature-neutral $k(T)$ that is a factor 25 below all experimental data.

The current rate analysis thus supports a negative temperature dependence, resulting from submerged barriers ~ 2 kcal/mol below the free reactants. The predicted absolute rate coefficient carries an *a priori* uncertainty of at least a factor of 5, and we recommend the use of the $k(T)$ expression by Tokuhashi et al. instead.

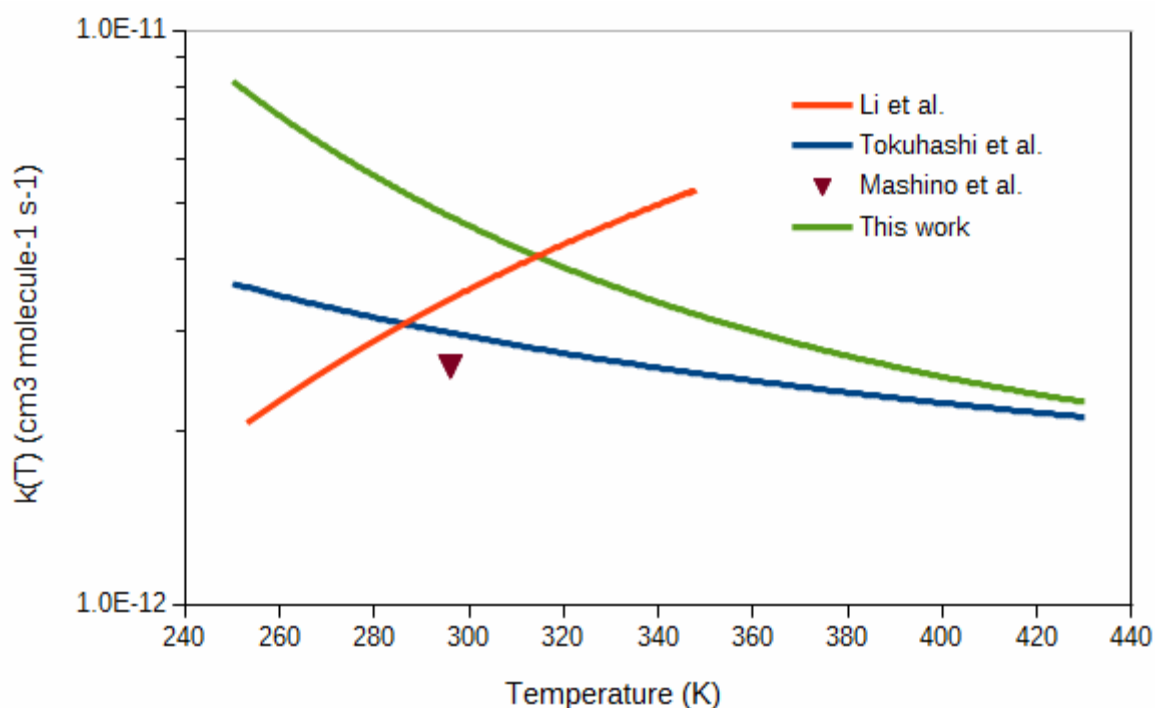


Figure SI-3 : Rate coefficients for the reaction of $\text{CF}_3\text{OCF}=\text{CF}_2 + \text{OH}$ ($\text{cm}^3 \text{ molecule}^{-1} \text{ s}^{-1}$), as reported in the literature, and as estimated in this work based on a rigid-rotor harmonic-oscillator CTST approximation using addition barrier heights of -2.1 and -1.5 kcal/mol relative to the free reactants, for addition on the outer and inner carbon atom, respectively.

The relative contribution of the two entrance channels can be estimated with fairly good accuracy, owing to the excellent error cancellation between the two addition transition states. The energy difference between the two TS at the M06-2X level of theory is 0.56 kcal/mol, leading to a ratio

of 83:17 at 300K for addition on the outer:inner carbon atom, using the RRHO approximation as detailed above, and changing only by ~4% absolute over the 240-400K T-range. Averaging the energy difference of the two TS over the levels of theory listed in Table SI-1 yields $\Delta E=0.88$ kcal/mol, leading to a ratio of 88:12. We thus estimate a contribution of $85\pm 15\%$ for addition on the outer carbon. The dominance of addition to the outer carbon is in agreement with literature data on OH addition on alkenes, where formation of the most substituted carbon radical is found to be favored. As discussed in the main text, the experimental observation of $69\pm 8\%$ of glyoxal by Mashino et al. at 10 Torr can be interpreted as indicative of a ratio of ~70:30, which remains within the uncertainties of the enthalpy and entropy calculations in the current theoretical work.

Quantum chemical data on reactants, products, and TS

CF3OCF=CF2 + OH : M06-2X/cc-pVDZ geometry

CF3

E(UM062X/CC-pVDZ) (Hartree): -337.47101834
Electronic state : 2-A
Cartesian coordinates (Angs):
C -0.000124 -0.000500 0.328762
F -1.104936 -0.593425 -0.073108
F 1.066832 -0.659480 -0.073081
F 0.038187 1.253238 -0.072987
Rotational constants (GHz): 10.8935400 10.8810600 5.6380500
Vibrational harmonic frequencies (cm-1):
516.3667 517.5031 705.8909
1128.4593 1335.8732 1337.8613
Zero-point correction (Hartree): 0.012626

CF3O

E(UM062X/CC-pVDZ) (Hartree): -412.67717517
Electronic state : 2-A
Cartesian coordinates (Angs):
C 0.006460 0.000842 0.035029
F -0.779589 -1.067848 -0.006813
F 0.850511 -0.027021 -0.989646
F -0.766167 1.075572 -0.051728
O 0.777306 0.021078 1.152939
Rotational constants (GHz): 6.0538100 5.8749700 5.6235700
Vibrational harmonic frequencies (cm-1):
240.4554 418.9718 588.5250
605.1949 631.2607 929.9720
1282.4225 1310.5443 1349.1661
Zero-point correction (Hartree): 0.016759

CFOCF2

E(UM062X/CC-pVDZ) (Hartree): -450.77929161
Point group : CS
Electronic state : 2-A"
Cartesian coordinates (Angs):
C -0.155187 -0.776621 0.000000
C 0.000000 0.643315 0.000000
F 1.039144 -1.385753 0.000000
O -1.195392 -1.365263 0.000000
F 1.152115 1.248269 0.000000
F -1.025230 1.439922 0.000000
Rotational constants (GHz): 5.6894300 3.4267300 2.1386300
Vibrational harmonic frequencies (cm-1):
132.1384 (A'') 217.1530 (A'') 220.6489 (A')
413.2745 (A') 537.0718 (A') 612.8043 (A')
644.8228 (A'') 829.7127 (A') 1174.1349 (A')
1467.7187 (A') 1586.2871 (A') 1848.7504 (A')
Zero-point correction (Hartree): 0.022063

CFOCF2OH

E(RM062X/CC-pVDZ) (Hartree): -526.65681872
Electronic state : 1-A
Cartesian coordinates (Angs):

```

C      -0.926111      0.130406      -0.008977
C      0.607752      0.021858      -0.004372
F      -1.509247     -1.052336     -0.016036
O      -1.519767      1.154825      0.003548
F      0.994365     -0.765495     -1.022032
F      0.969090     -0.614287      1.130834
O      1.186037      1.229197     -0.097212
H      0.492125      1.903303     -0.005491
Rotational constants (GHz):   3.8048500   2.5444800   2.0810400
Vibrational harmonic frequencies (cm-1):
  41.3475      212.8340      240.9736
 327.6125      404.0623      441.2731
 530.3607      614.9684      700.1705
 783.2171      837.9568     1117.0883
1217.7955      1280.0455     1375.8275
1485.0529      1983.0320     3779.6643
Zero-point correction (Hartree): 0.039579

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CFOCFO

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E(RM062X/CC-pVDZ) (Hartree): -426.22630863
Point group : CS
Electronic state : 1-A'
Cartesian coordinates (Angs):
  C      0.000000      0.765117      0.000000
  C     -0.000037     -0.765141      0.000000
  O     -0.955673      1.453999      0.000000
  O      0.955652     -1.453890      0.000000
  F      1.243061      1.217279      0.000000
  F     -1.243018     -1.217360      0.000000
Rotational constants (GHz):   5.9645800   3.5797300   2.2371000
Vibrational harmonic frequencies (cm-1):
  66.6815 ( A'')      245.8273 ( A')      423.0977 ( A')
 457.7500 ( A'')      535.3851 ( A')      691.0305 ( A')
 835.6805 ( A'')      841.6858 ( A')     1189.9163 ( A')
1367.2170 ( A')      1997.2040 ( A')     2001.6342 ( A')
Zero-point correction (Hartree): 0.024270

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CFO

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E(UM062X/CC-pVDZ) (Hartree): -213.03922496
Point group : CS
Electronic state : 2-A'
Cartesian coordinates (Angs):
  C      0.000000      0.423726      0.000000
  F     -1.012660     -0.432460      0.000000
  O      1.139243      0.168723      0.000000
Rotational constants (GHz):  186.9101500  11.5713600  10.8967600
Vibrational harmonic frequencies (cm-1):
  648.7703 ( A')      1106.6095 ( A')      1996.4126 ( A')
Zero-point correction (Hartree): 0.008547

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F

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E(UM062X/CC-pVDZ) (Hartree): -99.69555398
Point group : OH
Cartesian coordinates (Angs):
  F      0.000000      0.000000      0.000000
Zero-point correction (Hartree): 0.000000

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HF

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E(RM062X/CC-pVDZ) (Hartree): -100.40278281
Point group : C*v
Electronic state : 1-SG
Cartesian coordinates (Angs):
  F      0.000000      0.000000      0.092245
  H      0.000000      0.000000     -0.830203
Rotational constants (GHz):   0.0000000  620.5790284  620.5790284
Vibrational harmonic frequencies (cm-1):
  4129.9553 ( SG)

```

Zero-point correction (Hartree): 0.009409

O2

E(UM062X/CC-pVDZ) (Hartree): -150.27726911

Point group : D*H

Electronic state : 3-SGG

Cartesian coordinates (Angs):

O	0.000000	0.000000	0.595738
O	0.000000	0.000000	-0.595738

Rotational constants (GHz): 0.0000000 44.5138317 44.5138317

Vibrational harmonic frequencies (cm-1):

1777.0525 (SGG)

Zero-point correction (Hartree): 0.004048

OH

E(UM062X/CC-pVDZ) (Hartree): -75.70046613

Point group : C*V

Cartesian coordinates (Angs):

O	0.000000	0.000000	0.108808
H	0.000000	0.000000	-0.870468

Rotational constants (GHz): 0.0000000 555.8514600 555.8514600

Vibrational harmonic frequencies (cm-1):

3717.8362 (SG)

Zero-point correction (Hartree): 0.008470

FO2

E(UM062X/CC-pVDZ) (Hartree): -249.97293613

Point group : CS

Electronic state : 2-A"

Cartesian coordinates (Angs):

F	-0.993949	-0.608359	0.000000
O	0.000000	0.551603	0.000000
O	1.118193	0.132801	0.000000

Rotational constants (GHz): 80.6488800 11.3173200 9.9246100

Vibrational harmonic frequencies (cm-1):

419.2663 (A') 663.7730 (A') 1532.8411 (A')

Zero-point correction (Hartree): 0.005959

CF4

E(RM062X/CC-pVDZ) (Hartree): -437.38085819

Point group : CS

Electronic state : 1-A'

Cartesian coordinates (Angs):

C	-0.000033	0.000063	0.000000
F	-0.439754	1.243945	0.000000
F	1.319283	-0.000044	0.000000
F	-0.439754	-0.621972	1.077250
F	-0.439754	-0.621972	-1.077250

Rotational constants (GHz): 5.7310400 5.7308400 5.7304800

Vibrational harmonic frequencies (cm-1):

440.0897 (A'') 440.1489 (A') 636.6912 (A')
637.3753 (A'') 637.7396 (A') 943.4989 (A')
1358.9491 (A') 1359.7896 (A') 1359.8307 (A'')

Zero-point correction (Hartree): 0.017802

HCF3

E(RM062X/CC-pVDZ) (Hartree): -338.14538978

Electronic state : 1-A

Cartesian coordinates (Angs):

C	0.000092	0.000010	0.339330
F	0.905846	-0.858935	-0.128586
F	0.291058	1.213742	-0.128692
F	-1.196961	-0.354893	-0.128593
H	-0.000033	0.000711	1.436860

Rotational constants (GHz): 10.3655400 10.3611000 5.6902300

Vibrational harmonic frequencies (cm-1):

516.8866	517.1093	710.5043
1174.9865	1225.7947	1226.5443
1420.4163	1420.7385	3178.1038

Zero-point correction (Hartree): 0.025951

FNO

E(RM062X/CC-pVDZ) (Hartree): -229.64021922
Point group : CS
Electronic state : 1-A'
Cartesian coordinates (Angs):
F -0.945958 -0.556648 0.000000
N 0.000000 0.547666 0.000000
O 1.064203 0.147021 0.000000
Rotational constants (GHz): 97.1319400 12.5341900 11.1016100
Vibrational harmonic frequencies (cm-1):
606.1782 (A') 865.1589 (A') 1991.8551 (A')
Zero-point correction (Hartree): 0.007890

NO

E(UM062X/CC-pVDZ) (Hartree): -129.85561078
Point group : C*V
Cartesian coordinates (Angs):
N 0.000000 0.000000 -0.609646
O 0.000000 0.000000 0.533441
Rotational constants (GHz): 0.0000000 51.8018810 51.8018810
Vibrational harmonic frequencies (cm-1):
2101.9361 (SG)
Zero-point correction (Hartree): 0.004789

CF2O

E(RM062X/CC-pVDZ) (Hartree): -312.93589312
Point group : C2V
Electronic state : 1-A1
Cartesian coordinates (Angs):
C 0.000000 0.000000 0.141679
O 0.000000 0.000000 1.313238
F 0.000000 1.060818 -0.630888
F 0.000000 -1.060818 -0.630888
Rotational constants (GHz): 11.8192200 11.7736000 5.8981800
Vibrational harmonic frequencies (cm-1):
588.0028 629.2054 794.4039
1013.0912 1333.8831 2055.8938
Zero-point correction (Hartree): 0.014613

H2O

E(RM062X/CC-pVDZ) (Hartree): -76.38877424
Point group : C2V
Electronic state : 1-A1
Cartesian coordinates (Angs):
O 0.000000 0.000000 0.120151
H 0.000000 0.755011 -0.480605
H 0.000000 -0.755011 -0.480605
Rotational constants (GHz): 782.2606700 439.8404600 281.5396300
Vibrational harmonic frequencies (cm-1):
1648.4950 3842.0977 3943.2940
Zero-point correction (Hartree): 0.021492

CF3OH

E(RM062X/CC-pVDZ) (Hartree): -413.36940150
Electronic state : 1-A
Cartesian coordinates (Angs):
C -0.004235 0.023024 -0.000064
F 0.247193 -0.741912 -1.071102
F 0.251023 -0.732751 1.076428
F 0.833123 1.045907 -0.005726
O -1.261230 0.497464 0.000289

H -1.866800 -0.259049 0.001667
 Rotational constants (GHz): 5.7304900 5.6603700 5.6243900
 Vibrational harmonic frequencies (cm-1):
 254.9522 447.1666 460.1038
 608.7028 629.8676 643.5574
 934.3151 1155.9023 1266.7147
 1370.1402 1458.1360 3838.1555
 Zero-point correction (Hartree): 0.029770

CF3OCFCF2

 E(RM062X/CC-pVDZ) (Hartree): -788.35463805
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 1.672616 -0.287932 -0.068459
 C 0.694091 0.564996 -0.335089
 O -0.488561 0.168681 -0.874721
 F 0.823924 1.869870 -0.148821
 C -1.471733 -0.111240 0.043085
 F -1.683853 0.918544 0.854497
 F -1.141887 -1.159607 0.790339
 F -2.570402 -0.380072 -0.628547
 F 1.566372 -1.581829 -0.246295
 F 2.843472 0.072607 0.396666
 Rotational constants (GHz): 2.3475600 0.9395300 0.7772000
 Vibrational harmonic frequencies (cm-1):
 42.7445 67.2912 118.5448
 192.6827 210.7831 341.5436
 369.1534 452.4712 469.3770
 546.3894 554.7824 607.0276
 641.9512 652.4408 726.3140
 856.0562 946.8563 1225.8433
 1270.7324 1303.7518 1382.7947
 1388.7653 1429.0729 1961.5734
 Zero-point correction (Hartree): 0.040458

CF3OCFCF2OH.b

 E(UM062X/CC-pVDZ) (Hartree): -864.15280338
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C 1.502030 -0.252021 -0.021838
 C 0.534407 0.774244 -0.585957
 O -0.714846 0.384491 -0.972365
 F 0.556457 1.932774 0.051753
 C -1.579501 -0.094752 -0.028635
 F -1.607587 0.664494 1.058163
 F -1.223854 -1.332989 0.362502
 F -2.775371 -0.157696 -0.565521
 F 1.413833 -1.359640 -0.788736
 O 1.312176 -0.555740 1.292582
 F 2.735018 0.253939 -0.128970
 H 0.593268 -1.202761 1.354128
 Rotational constants (GHz): 1.9696800 0.8478200 0.7506500
 Vibrational harmonic frequencies (cm-1):
 34.0092 57.7252 130.6047
 166.3854 196.9071 218.3432
 300.4147 330.8038 364.7875
 430.5164 485.7505 511.8767
 558.3645 586.8435 632.2428
 637.7359 687.4308 764.8264
 824.3453 921.6583 1140.6732
 1141.4601 1226.2863 1263.6132
 1279.1272 1347.5469 1386.6135
 1410.4939 1440.3405 3841.3603
 Zero-point correction (Hartree): 0.055403

CF3OCFCF2OH

 E(UM062X/CC-pVDZ) (Hartree): -864.15271486
 Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.500508	-0.282034	-0.043294
C	0.538221	0.729093	-0.626795
O	-0.721638	0.391282	-0.972266
F	0.640371	1.923976	-0.036838
C	-1.584319	-0.096976	-0.008336
F	-1.441210	0.554633	1.138165
F	-1.375259	-1.386552	0.206145
F	-2.809080	0.070498	-0.463405
F	1.528887	-1.365365	-0.820154
O	2.758734	0.204735	0.060467
F	1.040946	-0.677631	1.180618
H	2.714872	1.055345	0.524159

Rotational constants (GHz): 1.9624500 0.8637900 0.7603500

Vibrational harmonic frequencies (cm-1):

31.1296	71.9159	109.6047
151.8585	196.0286	216.7416
280.2610	341.3362	359.5651
428.2155	473.3544	521.8439
554.3154	590.7160	632.1368
636.9499	686.3507	766.1424
819.7463	922.6191	1096.5140
1135.6155	1218.4450	1288.1493
1313.9371	1341.1435	1369.4645
1399.4842	1475.4214	3828.9544

Zero-point correction (Hartree): 0.055264

CF3OCFCFOH

E(RM062X/CC-pVDZ) (Hartree): -764.35221643

Electronic state : 1-A

Cartesian coordinates (Angs):

C	1.682358	-0.319785	-0.070948
C	0.689821	0.521001	-0.348632
O	-0.508046	0.163192	-0.880570
F	0.864540	1.841022	-0.167272
C	-1.486363	-0.096591	0.044836
F	-1.671217	0.936875	0.860975
F	-1.177893	-1.153370	0.789486
F	-2.599475	-0.338905	-0.616260
F	1.551738	-1.617250	-0.237490
O	2.889670	0.009563	0.393260
H	2.922873	0.974845	0.481991

Rotational constants (GHz): 2.3447800 0.9509600 0.7844900

Vibrational harmonic frequencies (cm-1):

45.7715	68.7031	110.0883
191.9706	209.4577	340.8296
351.6364	370.1612	437.5304
468.8086	548.0042	559.8556
624.6288	642.2215	654.0201
717.0477	855.6679	946.9498
1152.5390	1271.5632	1296.5798
1308.2966	1369.0874	1379.4009
1481.6466	1934.6331	3825.0868

Zero-point correction (Hartree): 0.052767

CF3OCFCFO

E(UM062X/CC-pVDZ) (Hartree): -763.73876235

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.826660	-0.306897	-0.011068
C	0.733990	0.588808	-0.292829
O	-0.429580	0.206207	-0.843029
F	0.902570	1.874515	-0.141163
C	-1.445694	-0.103548	0.043163
F	-1.752313	0.948446	0.789716
F	-1.082193	-1.091046	0.848178
F	-2.485288	-0.465587	-0.670284
F	1.475239	-1.576069	-0.251364
O	2.903096	0.008480	0.391610

Rotational constants (GHz): 2.3575000 0.9821800 0.8048500
 Vibrational harmonic frequencies (cm-1):
 34.6011 49.7698 74.2047
 142.3748 209.5242 327.5129
 354.0007 382.2282 466.1603
 534.6876 564.6519 638.6468
 665.7959 669.2335 702.1340
 838.0330 946.0423 1165.2849
 1239.1636 1317.6725 1390.8386
 1429.3370 1529.3187 1864.1480
 Zero-point correction (Hartree): 0.039949

CF3OCFOCF2OH

E(UM062X/CC-pVDZ) (Hartree): -939.35055168

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.351517	-0.530633	-0.022931
C	0.567815	0.867204	0.000152
O	-0.721303	0.810099	-0.541184
F	1.228040	1.676829	-0.848812
C	-1.661809	-0.048858	-0.043660
F	-1.657106	-0.107965	1.282132
F	-1.488873	-1.280349	-0.514410
F	-2.835868	0.395782	-0.453244
F	1.188648	-1.045071	-1.233854
O	2.667337	-0.371922	0.199876
F	0.773628	-1.347700	0.877370
H	2.787159	0.100392	1.038901
O	0.652902	1.255524	1.263445

Rotational constants (GHz): 1.5663200 0.7911300 0.7189700

Vibrational harmonic frequencies (cm-1):
 62.6395 79.8852 116.8089
 169.5713 200.9001 225.4360
 274.5833 315.1010 339.9183
 357.4456 381.3601 433.9029
 471.4822 517.9227 536.1759
 585.0397 624.8445 660.6677
 687.7100 734.4103 768.9161
 874.5457 977.2055 1095.3918
 1173.5591 1228.1175 1289.0952
 1300.1688 1312.2486 1335.9739
 1382.4719 1439.5061 3818.5755

Zero-point correction (Hartree): 0.058712

CF3OCFOCFO

E(UM062X/CC-pVDZ) (Hartree): -838.92025750

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.542186	-0.562552	0.161134
C	0.724132	0.727279	-0.152791
O	-0.503760	0.456203	-0.771156
F	1.409853	1.413435	-1.087698
C	-1.472993	-0.127597	-0.002901
F	-2.052376	0.741183	0.811521
F	-0.964218	-1.115729	0.742746
F	-2.377289	-0.623474	-0.819367
F	1.403667	-1.413816	-0.844425
O	2.215047	-0.749146	1.106257
O	0.596629	1.388297	1.007694

Rotational constants (GHz): 1.7516400 0.9239000 0.8419900

Vibrational harmonic frequencies (cm-1):
 39.2746 58.6082 91.7071
 165.0382 215.6815 232.5666
 334.0522 356.4422 376.2570
 433.2546 483.9399 520.5625
 622.1618 646.5482 677.6836
 714.0592 786.4851 821.2214
 942.2781 1071.3351 1186.8303
 1218.5876 1289.7746 1304.0120

1318.3689 1389.1805 2000.5364
Zero-point correction (Hartree): 0.043961

CF3OCFOHCF2

E(UM062X/CC-pVDZ) (Hartree): -864.15090431

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.468539	0.672877	0.102535
C	-0.895448	-0.724242	-0.011581
O	0.398115	-0.803145	-0.588518
F	-0.811504	-1.233597	1.243742
C	1.399895	-0.005914	-0.118907
F	1.512324	-0.046088	1.202785
F	1.208681	1.265951	-0.472754
F	2.524617	-0.431574	-0.664141
F	-1.504108	1.340980	-1.031352
F	-0.976741	1.395671	1.089820
O	-1.716115	-1.448424	-0.799620
H	-1.250861	-2.265873	-1.040083

Rotational constants (GHz): 1.5597400 1.0000600 0.8848400

Vibrational harmonic frequencies (cm-1):

52.1003	99.6120	130.5446
157.2650	204.2146	218.4213
305.6484	349.5107	365.3784
406.8178	449.6697	507.2350
550.7400	582.1924	613.4673
659.0262	683.4816	752.4771
833.6755	937.8475	1080.2086
1145.7284	1262.9880	1299.3164
1314.5774	1347.5755	1364.6519
1378.1792	1508.8686	3804.1409

Zero-point correction (Hartree): 0.055509

CF3OCFOCF2OH.a

E(UM062X/CC-pVDZ) (Hartree): -1014.49166116

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.086260	0.885992	-0.003508
C	-0.498625	-0.552614	0.091905
O	0.715372	-0.742217	-0.518433
F	-0.494074	-0.944337	1.368922
C	1.870826	-0.123141	-0.086517
F	1.842099	0.151874	1.207544
F	2.084549	0.992476	-0.763168
F	2.860159	-0.960825	-0.332803
F	-1.310368	1.118156	-1.314485
O	-2.176623	1.085897	0.753225
F	-0.143939	1.744367	0.406341
O	-1.299367	-1.485605	-0.652266
O	-2.548723	-1.423096	-0.303101
H	-2.786748	0.343351	0.602162

Rotational constants (GHz): 1.3350000 0.6682700 0.5838000

Vibrational harmonic frequencies (cm-1):

31.4599	78.3334	90.5647
120.5422	169.1334	217.1604
229.1803	287.9877	328.7751
347.7605	367.6370	399.3712
442.9937	446.5259	477.1721
543.9579	568.8980	606.2962
624.1929	683.5254	708.9546
763.6965	815.8179	951.7727
1076.3876	1171.8793	1193.4084
1236.1015	1287.8150	1310.5447
1314.9903	1336.0462	1364.9986
1399.9767	1455.9734	3771.0607

Zero-point correction (Hartree): 0.064292

CF3OCFOCF2OH.b

E(UM062X/CC-pVDZ) (Hartree): -1014.49039045
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	-0.992171	0.939396	-0.024246
C	-0.541969	-0.545378	0.031186
O	0.688170	-0.832431	-0.501728
F	-0.649540	-0.958162	1.304024
C	1.826264	-0.183361	-0.073707
F	1.870331	-0.081596	1.246036
F	1.919481	1.027385	-0.600656
F	2.842126	-0.910726	-0.491756
F	-0.749942	1.416015	-1.245293
O	-2.297579	1.115552	0.251363
F	-0.207615	1.613173	0.850007
H	-2.561557	0.510998	0.963705
O	-1.388872	-1.352621	-0.787318
O	-2.553566	-1.521717	-0.227862

Rotational constants (GHz): 1.3084300 0.6908800 0.5936000
 Vibrational harmonic frequencies (cm-1):

64.6053	82.1346	105.4384
149.8109	155.0087	210.1429
216.7138	227.3207	276.5627
331.5791	354.0653	368.8438
402.2195	438.7321	476.2832
546.8789	572.0087	598.0006
631.7293	697.3217	715.2341
765.3362	815.8198	952.8177
1079.3707	1156.6953	1194.9612
1232.1090	1270.4203	1294.6427
1312.8804	1329.2816	1366.3555
1393.0315	1453.2218	3812.5026

Zero-point correction (Hartree): 0.063903

CF3OCFOHCF2O

 E(UM062X/CC-pVDZ) (Hartree): -1014.45407928
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	1.594121	-0.730088	-0.040296
C	0.368272	0.247894	0.069703
O	-0.686573	-0.405477	-0.522742
F	0.144359	0.502683	1.366196
C	-1.960605	-0.102773	-0.091876
F	-2.221414	-0.675785	1.072848
F	-2.787211	-0.578536	-1.000766
F	-2.140996	1.204284	0.031345
F	1.703775	-1.102319	-1.331634
O	2.688275	-0.119405	0.408674
F	1.303895	-1.834177	0.669361
O	0.563465	1.415477	-0.640991
O	1.641331	2.085787	-0.016372
H	2.315618	2.053399	-0.719885

Rotational constants (GHz): 1.3195100 0.6285400 0.5485000
 Vibrational harmonic frequencies (cm-1):

42.8522	49.2126	74.7488
123.3936	161.0846	209.7715
223.9721	260.5212	306.2218
340.3914	354.9138	365.3602
385.5199	428.5558	500.7852
548.7252	562.8610	595.1564
632.2285	654.9522	715.8977
744.1175	851.1434	934.3263
1047.9762	1062.1653	1155.4691
1226.5319	1237.0957	1250.5931
1317.1305	1323.3292	1336.5852
1388.2496	1446.7462	3747.1922

Zero-point correction (Hartree): 0.062891

CF3COHCF2

 E(RM062X/CC-pVDZ) (Hartree): -764.35149549

Electronic state : 1-A
 Cartesian coordinates (Angs):

C	1.664350	-0.289644	-0.061357
C	0.700557	0.578653	-0.346123
O	-0.483577	0.102792	-0.892531
C	-1.449568	-0.135391	0.029921
F	-1.638320	0.941352	0.818407
F	-1.142101	-1.151746	0.829778
F	-2.572987	-0.402550	-0.600661
F	2.833911	0.053771	0.422852
F	1.567040	-1.582533	-0.264814
O	0.845215	1.912211	-0.231304
H	0.186973	2.233624	0.405987

Rotational constants (GHz): 2.3594500 0.9482300 0.7873400
 Vibrational harmonic frequencies (cm-1):

56.3819	73.8533	130.0879
203.8369	214.2157	325.5031
362.1713	370.6758	453.0713
480.9830	543.2033	561.7109
620.8788	645.0231	654.6573
719.0532	850.7701	940.0746
1179.7084	1239.4897	1294.3232
1324.6848	1368.6271	1387.3415
1419.5866	1945.3410	3784.9508

Zero-point correction (Hartree): 0.052740

CH3OCHCH2

 E(RM062X/CC-pVDZ) (Hartree): -193.03591470
 Electronic state : 1-A
 Cartesian coordinates (Angs):

C	1.434143	0.568648	0.000003
C	0.732832	-0.568565	-0.000005
O	-0.609584	-0.716963	0.000013
H	1.217255	-1.547108	-0.000025
C	-1.357855	0.478097	-0.000007
H	-1.138226	1.081005	-0.897020
H	-1.138307	1.080989	0.897037
H	-2.413629	0.186560	-0.000054
H	0.974108	1.555502	0.000027
H	2.520750	0.509679	-0.000013

Rotational constants (GHz): 18.2889000 6.4601700 4.9248000
 Vibrational harmonic frequencies (cm-1):

256.0947	280.8650	341.4340
596.2423	728.0034	847.8294
942.1595	1009.1841	1057.3613
1171.9543	1217.2540	1278.5429
1351.5511	1406.9691	1465.9034
1467.7104	1487.6039	1736.6951
3036.3770	3106.2853	3180.9864
3203.9581	3223.2741	3303.1601

Zero-point correction (Hartree): 0.085881

CH3OCHCH2OH

 E(UM062X/CC-pVDZ) (Hartree): -268.78646623
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	-0.950174	-0.394724	0.318576
C	-0.039447	0.740023	-0.035034
O	1.310180	0.668494	0.075055
H	-0.401361	1.754792	0.146213
C	1.903455	-0.595498	-0.139550
H	1.587391	-1.023122	-1.105581
H	1.657706	-1.303593	0.668569
H	2.987015	-0.432705	-0.153471
H	-1.055311	-0.512357	1.414269
O	-2.252888	-0.134008	-0.159411
H	-0.557859	-1.355553	-0.065077
H	-2.158923	0.097838	-1.094021

Rotational constants (GHz): 14.4439400 2.6519400 2.3869200

Vibrational harmonic frequencies (cm-1):

78.0753	160.9798	249.6241
261.3442	372.0982	385.3998
548.6671	643.6833	941.1431
1012.2734	1116.1845	1131.0510
1160.5391	1190.4572	1203.5872
1318.6425	1322.3098	1388.0796
1446.9464	1473.4654	1477.6533
1487.2814	1497.1383	3003.6716
3031.3064	3041.4855	3108.8984
3172.9491	3202.9570	3823.7039

Zero-point correction (Hartree): 0.100813

CH3OCHOHCH2

E(UM062X/CC-pVDZ) (Hartree): -268.79105185
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-0.746925	1.395192	-0.111862
C	-0.464829	-0.008616	0.308975
O	0.659940	-0.540309	-0.362533
H	-0.266310	-0.050024	1.402914
C	1.874678	0.006341	0.088825
H	2.019375	-0.167479	1.171916
H	1.929814	1.093358	-0.099047
H	2.680780	-0.491940	-0.462591
H	-0.816033	1.591653	-1.182673
H	-1.093970	2.142602	0.600009
O	-1.555343	-0.803955	-0.040507
H	-1.267978	-1.721560	0.078161

Rotational constants (GHz): 8.5975000 4.0586800 3.0513600

Vibrational harmonic frequencies (cm-1):

64.4565	121.6963	208.7737
285.1803	320.1088	380.3213
472.6788	529.8539	597.2581
871.4363	1045.9035	1103.3912
1161.8142	1174.7358	1190.1219
1236.6008	1282.2993	1368.5716
1399.2550	1455.6505	1465.1359
1486.7515	1495.5145	2953.5102
3011.8079	3074.2263	3166.0398
3177.4365	3300.4756	3816.8991

Zero-point correction (Hartree): 0.098458

complex.CF3OCFCF2.OH.a

E(UM062X/CC-pVDZ) (Hartree): -864.06225948
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.508895	-0.741170	-0.098708
C	0.676475	0.217200	0.277209
O	-0.468236	0.497055	-0.411411
F	0.912975	1.009484	1.307922
C	-1.581220	-0.187689	0.018678
F	-1.791635	-0.000474	1.315632
F	-1.452800	-1.490744	-0.194850
F	-2.607631	0.271514	-0.666953
F	1.281534	-1.528877	-1.121609
F	2.638978	-1.001352	0.509560
H	0.174459	2.620018	-0.924354
O	1.139217	2.792192	-0.914344

Rotational constants (GHz): 1.2933600 0.8178700 0.6594800

Vibrational harmonic frequencies (cm-1):

25.1118	40.2045	55.5376
71.9722	89.0994	117.3545
145.5216	195.4270	215.2323
335.5326	343.8949	370.7291
460.3984	469.9556	547.1884
555.4644	609.5410	644.3089
652.8220	727.7043	855.5684
949.6319	1225.4701	1270.2020

1310.7071 1375.6679 1383.1856
 1430.0651 1969.1764 3709.0402
 Zero-point correction (Hartree): 0.050465

complex.CF3OCFCF2.OH.b

 E(UM062X/CC-pVDZ) (Hartree): -864.06520395
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C 1.580719 -0.369306 -0.371650
 C 0.583567 0.455330 -0.676246
 O -0.639913 -0.002858 -1.056408
 F 0.733549 1.765442 -0.707633
 C -1.568594 -0.110326 -0.055349
 F -1.727867 1.027812 0.600363
 F -1.205146 -1.046558 0.831888
 F -2.705125 -0.471362 -0.609008
 F 1.437861 -1.672310 -0.336831
 F 2.793910 0.025486 -0.100555
 H 0.326080 -0.279005 2.198820
 O 0.909303 0.473886 1.970987
 Rotational constants (GHz): 1.7056800 0.7952200 0.7542700
 Vibrational harmonic frequencies (cm-1):
 30.1339 61.3884 79.9992
 100.0813 114.8918 148.6306
 185.3447 199.5076 212.5133
 341.2620 367.1227 416.6684
 450.4326 470.8440 544.3322
 556.1535 589.8232 641.8362
 650.7202 722.1286 853.5707
 945.4680 1226.7566 1258.8383
 1297.4994 1380.9804 1394.1516
 1436.4112 1929.2556 3726.7267
 Zero-point correction (Hartree): 0.050879

CF2OCFCF2oxide

 E(UM062X/CC-pVDZ) (Hartree): -763.66994286
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C 1.618728 -0.267317 0.055010
 C 0.342268 0.363359 -0.214158
 O -0.739438 -0.369247 -0.607891
 F 0.264156 1.633928 -0.594234
 C -1.811541 -0.308347 0.241818
 F -2.214239 0.930430 0.486758
 F -2.783941 -1.043499 -0.241506
 F 1.803889 -1.571341 -0.034415
 O 0.814851 0.166919 1.093429
 F 2.763465 0.371867 -0.103306
 Rotational constants (GHz): 2.7697900 0.8727200 0.7390500
 Vibrational harmonic frequencies (cm-1):
 31.0534 61.5035 127.8471
 177.1115 243.9467 296.4394
 381.8529 502.1014 515.8797
 550.5544 557.1616 596.5021
 685.2540 755.5934 825.7184
 887.6839 1079.7284 1186.0441
 1218.9193 1279.1284 1329.7550
 1361.9423 1368.8636 1684.8001
 Zero-point correction (Hartree): 0.040336

CF3OCFCF2oxide

 E(UM062X/CC-pVDZ) (Hartree): -763.68149156
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C -1.914947 0.012978 -0.071451
 C -0.551255 -0.383678 -0.291268
 O 0.483133 0.470455 -0.511606
 C 1.703853 0.031279 -0.056713

F	2.059428	-1.097543	-0.658385	
F	2.582200	0.971735	-0.332870	
F	-2.287228	1.274355	0.057100	
O	-1.137969	-0.484820	0.969882	
F	-2.945167	-0.722478	-0.444706	
F	1.681075	-0.187020	1.251126	
Rotational constants (GHz):	3.2482700	0.7258200	0.6986100	
Vibrational harmonic frequencies (cm-1):				
20.4560	65.3396		113.2779	
201.2603	265.6559		358.6345	
442.2116	502.3326		528.4847	
549.4260	622.1372		632.4332	
675.5862	734.2437		809.1391	
925.7098	966.1950		1177.8388	
1224.4180	1291.7392		1311.4290	
1343.5575	1381.0109		1651.9387	
Zero-point correction (Hartree):	0.040539			

CF3OCOCF2OH

E(RM062X/CC-pVDZ) (Hartree): -839.61717822

Electronic state : 1-A

Cartesian coordinates (Angs):

C	1.854914	-0.157632	-0.002875	
C	0.436579	0.445252	-0.004516	
O	-0.500059	-0.526803	-0.009251	
C	-1.829387	-0.130295	-0.001162	
F	-2.116372	0.589773	-1.071609	
F	-2.108393	0.571514	1.083410	
F	-2.548265	-1.232811	-0.007889	
F	1.975985	-0.988816	-1.054245	
O	2.787588	0.804982	-0.045269	
F	1.988208	-0.918359	1.102918	
H	2.332952	1.664197	-0.009814	
O	0.234212	1.621839	0.003002	
Rotational constants (GHz):	2.2259100	0.6748200	0.6323600	
Vibrational harmonic frequencies (cm-1):				
7.6795	58.0155		104.5933	
132.4288	215.4772		218.0114	
329.9825	346.8529		391.8969	
443.7729	447.1603		533.7401	
567.4955	592.4113		628.7016	
678.3863	766.7280		789.3800	
916.1523	933.4619		1128.1279	
1219.0141	1230.7431		1298.9084	
1339.9056	1375.3228		1380.2128	
1488.2195	1936.3847		3764.9931	
Zero-point correction (Hartree):	0.057556			

CFOHCF2

E(RM062X/CC-pVDZ) (Hartree): -451.38638020

Electronic state : 1-A

Cartesian coordinates (Angs):

C	0.659810	0.013499	-0.011909	
C	-0.665037	-0.027700	0.009103	
O	1.440381	-1.071026	-0.106666	
F	1.290884	1.191847	0.007668	
F	-1.432303	1.042984	-0.004494	
F	-1.356974	-1.148147	0.016993	
H	1.993857	-1.126735	0.688658	
Rotational constants (GHz):	5.4523500	3.2790500	2.0583800	
Vibrational harmonic frequencies (cm-1):				
195.4942	208.9100		265.7097	
401.0679	440.7599		561.4975	
568.8550	589.5642		822.2124	
1181.7732	1306.1006		1387.9893	
1408.3588	1971.4149		3785.6745	
Zero-point correction (Hartree):	0.034390			

CFOCFOH

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E(UM062X/CC-pVDZ) (Hartree): -426.78665429
Electronic state : 2-A
Cartesian coordinates (Angs):
  C    -0.756121    0.121130    0.000383
  C     0.654128   -0.033315    0.000012
  F    -1.430369   -1.033527    0.000092
  O    -1.296256    1.203679   -0.000371
  F     1.257553   -1.192410   -0.000159
  O     1.457771    1.010350   -0.000048
  H     0.875186    1.794315    0.001579
Rotational constants (GHz):   5.6139100   3.5443500   2.1726500
Vibrational harmonic frequencies (cm-1):
  162.1495                226.0065                335.7396
  409.8361                557.5674                558.8438
  609.5359                648.9475                845.0635
  1164.8958              1280.4961              1534.9799
  1639.6829              1808.3475              3723.5835
Zero-point correction (Hartree): 0.035325

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CFOCF2O

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E(UM062X/CC-pVDZ) (Hartree): -525.96466685
Point group : CS
Electronic state : 2-A'
Cartesian coordinates (Angs):
  C     0.321360   -0.881000    0.000000
  C    -0.045881    0.626598    0.000000
  F    -0.803099   -1.582466    0.000000
  O     1.399106   -1.350222    0.000000
  F    -0.803099    0.893056    1.079353
  F    -0.803099    0.893056   -1.079353
  O     1.104743    1.311921    0.000000
Rotational constants (GHz):   3.8873600   2.5794400   2.1288200
Vibrational harmonic frequencies (cm-1):
  33.3852 ( A'')    216.2060 ( A')    223.9653 ( A'')
  374.4052 ( A')    411.2589 ( A')    473.0212 ( A'')
  591.2733 ( A')    682.6296 ( A')    768.5877 ( A'')
  824.3158 ( A')    1069.0838 ( A')   1218.1761 ( A'')
  1248.9357 ( A')   1336.0031 ( A')   2001.8490 ( A')
Zero-point correction (Hartree): 0.026138

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OCCF2OH

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E(UM062X/CC-pVDZ) (Hartree): -426.76586253
Electronic state : 2-A
Cartesian coordinates (Angs):
  C    -1.039773   -0.603678   -0.369696
  C     0.373060    0.006150   -0.018933
  O    -2.033401   -0.058667   -0.061981
  F     1.244688   -0.286037   -0.979092
  F     0.764517   -0.632185    1.105392
  O     0.336131    1.343393    0.140207
  H    -0.504399    1.571367    0.569264
Rotational constants (GHz):   5.7016900   3.0407200   2.9893800
Vibrational harmonic frequencies (cm-1):
  90.8463                251.9533                340.2675
  381.8375                421.9729                539.0082
  563.0071                666.7879                799.7310
  1097.2475              1195.4603              1313.7883
  1430.7454              2015.3724              3803.2641
Zero-point correction (Hartree): 0.033970

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TS.CF3OCFCF2OH.CF3+HOFCF2CFO

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E(UM062X/CC-pVDZ) (Hartree): -864.10683749
Electronic state : 2-A
Cartesian coordinates (Angs):
  C     1.629197   -0.310862   -0.009215
  C     0.621532    0.725049   -0.487774
  O    -0.402939    0.472511   -1.099931

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F	0.824853	1.937740	0.054701	
C	-1.857238	-0.113453	0.023013	
F	-1.488521	0.259614	1.228762	
F	-1.969980	-1.412534	-0.067506	
F	-2.941701	0.502001	-0.368663	
F	1.942317	-1.128810	-1.017292	
O	2.770159	0.207774	0.489128	
F	0.982776	-1.075009	0.918540	
H	2.553609	1.006297	0.993404	
Rotational constants (GHz):	1.8827000		0.7356500	0.6587200
Vibrational harmonic frequencies (cm-1):				
i551.1443	26.3667			41.6632
59.9623	103.3701			146.1078
197.2344	251.8859			265.8266
303.6714	400.6340			444.0886
510.6656	527.5065			539.5820
599.7575	639.9165			679.3325
720.7646	820.8845			1042.6922
1091.5365	1157.8104			1264.3768
1365.8613	1379.6566			1391.7052
1441.1681	1595.9698			3839.7409
Zero-point correction (Hartree):	0.052056			

TS.CF3OCFCF2OH.CF3OCF2CFOH

E(UM062X/CC-pVDZ) (Hartree): -864.08993652

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.657246	-0.363637	-0.273586
C	0.597298	0.463397	-0.575642
O	-0.592681	0.000751	-0.959725
F	0.759773	1.760606	-0.599711
C	-1.600205	-0.113743	0.001294
F	-1.714495	0.990319	0.709831
F	-1.378665	-1.140504	0.793635
F	-2.712302	-0.320975	-0.678369
F	1.520551	-1.645928	-0.440707
O	2.845106	0.053490	0.083761
F	0.790292	0.248350	1.463314
H	2.668182	0.623154	0.863376

Rotational constants (GHz): 1.9178600 0.8277900 0.7541200

Vibrational harmonic frequencies (cm-1):

i287.0679	21.1511		91.4147
137.2784	163.5648		206.0898
218.7353	297.8701		339.7178
354.0187	375.9234		467.0722
481.9111	540.1653		549.9658
575.5125	636.6884		654.3041
720.0576	848.6219		950.6233
1194.6503	1206.7557		1291.6838
1354.3349	1365.9412		1446.3046
1520.8888	1773.4148		3648.3167

Zero-point correction (Hartree): 0.053384

TS.CF3OCFCF2OH.CF3OCFCF2+OH

E(UM062X/CC-pVDZ) (Hartree): -864.06253826

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.407800	-0.377410	-0.336857
C	-0.483199	0.544016	-0.029448
O	0.588520	0.275364	0.748232
F	-0.645469	1.813453	-0.369235
C	1.721636	-0.086215	0.056132
F	2.065339	0.848167	-0.823455
F	1.532852	-1.221897	-0.606623
F	2.684566	-0.243565	0.936059
F	-1.268300	-1.639480	-0.058293
O	-2.307203	0.158229	1.586709
F	-2.436039	-0.112607	-1.097741
H	-2.630899	1.052258	1.355102

Rotational constants (GHz): 1.8089000 0.7346200 0.6841400
 Vibrational harmonic frequencies (cm-1):
 i292.7795 37.9959 68.1747
 79.6251 103.9954 140.7065
 190.2716 199.6665 243.2688
 342.3516 368.6890 411.5625
 465.5586 507.0406 555.5844
 558.1256 641.5671 649.4315
 664.9366 719.6689 849.7104
 946.4766 1228.5566 1262.8040
 1304.6369 1384.9430 1398.6794
 1452.5751 1846.1985 3726.2009
 Zero-point correction (Hartree): 0.050915

TS.CF3OCFCF2OH.CF2O+CF2O+CF+HF

E(UM062X/CC-pVDZ) (Hartree): -864.04373107

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.659760	-0.423680	-0.038935
C	-1.064546	1.256439	0.572911
O	0.894282	0.465073	1.059965
F	-0.726443	1.990782	-0.391549
C	1.770485	0.316736	0.267872
F	1.727884	0.734706	-0.962318
F	1.220066	-1.782023	-0.583384
F	2.940241	-0.152963	0.545489
F	-1.510972	-1.136003	1.076006
O	-1.007207	-0.665198	-1.043958
F	-2.952916	-0.076374	-0.140271
H	0.345584	-1.499101	-0.834896

Rotational constants (GHz): 1.6736400 0.7458900 0.6414700

Vibrational harmonic frequencies (cm-1):
 i121.2333 40.5500 50.4237
 72.0061 101.7826 141.2068
 156.0526 169.1996 203.6548
 213.8516 232.6536 270.0908
 369.4656 385.1384 560.4618
 593.5700 617.0532 653.8522
 716.5701 767.7607 795.0455
 853.9754 959.6299 1040.6265
 1252.6612 1387.9249 1431.3202
 1708.7091 1942.5242 3455.6154

Zero-point correction (Hartree): 0.048168

TS.CF3OCFCF2OH.CF3OCFCFO+HF

E(UM062X/CC-pVDZ) (Hartree): -864.09142693

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.535440	-0.172146	-0.350612
C	0.485968	0.796778	-0.533397
O	-0.751730	0.481182	-0.925903
F	0.648998	1.987865	-0.026365
C	-1.581072	-0.117645	0.021875
F	-1.566644	0.571164	1.149995
F	-1.193118	-1.354356	0.265881
F	-2.791130	-0.123063	-0.489800
F	1.316057	-1.295749	-0.988405
O	2.721498	0.152674	-0.013472
F	1.273892	-0.662900	1.384220
H	2.407334	-0.219429	1.028059

Rotational constants (GHz): 1.8406000 0.8516200 0.7695600

Vibrational harmonic frequencies (cm-1):
 i1442.1940 30.9921 62.5143
 79.9027 104.8496 159.6607
 206.6040 294.6185 346.2982
 373.7423 449.7006 463.7451
 538.5035 561.8552 620.8879
 631.0149 664.3307 705.6063
 790.7884 846.4672 926.3906

946.2653	1201.8459	1246.3266
1338.2892	1383.1716	1439.8977
1521.3787	1671.2167	2151.7167

Zero-point correction (Hartree): 0.049570

TS.CF3OCFCF2OH.CF3OCFHCF2O

E(UM062X/CC-pVDZ) (Hartree): -864.08853855

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.807480	-0.241396	0.005136
C	-0.458082	0.427938	0.035456
O	0.541447	-0.251413	-0.595244
F	-2.814840	0.562866	-0.350202
F	-0.440624	1.743397	-0.165095
F	-1.891239	-1.354076	-0.728696
O	-1.766299	-0.491366	1.353188
H	-0.597301	0.114028	1.312167
C	1.795591	-0.139531	-0.038174
F	1.796741	-0.596210	1.210345
F	2.207111	1.120209	-0.014545
F	2.611288	-0.859949	-0.772942

Rotational constants (GHz): 2.1052600 0.6925300 0.6379300

Vibrational harmonic frequencies (cm-1):

i1994.7735	25.8416	64.3534
91.8140	160.6811	209.5000
270.9146	332.1101	365.0041
429.9808	482.7567	531.4378
551.8734	591.5122	639.0407
658.7853	732.2696	754.9359
904.0860	937.8266	1085.5425
1182.7294	1216.6856	1247.4372
1288.3268	1311.1927	1337.6063
1390.5746	1488.5048	2032.7261

Zero-point correction (Hartree): 0.050840

TS.CF3OCFCFO.CFOCFO+CF3

E(UM062X/CC-pVDZ) (Hartree): -763.68593667

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.790748	-0.439115	0.009161
C	0.914048	0.687222	-0.434052
O	-0.123754	0.556627	-1.056845
F	1.324932	1.863740	0.008022
C	-1.639347	-0.055597	0.084879
F	-2.533931	0.880655	0.253308
F	-1.003871	-0.333381	1.202075
F	-2.131731	-1.121514	-0.479623
F	1.235963	-1.601515	-0.333068
O	2.821885	-0.349993	0.579799

Rotational constants (GHz): 2.1866000 0.9042700 0.7559300

Vibrational harmonic frequencies (cm-1):

i492.5493	39.9501	51.2726
67.9916	109.6130	144.9495
226.1791	269.6194	431.4772
459.6702	527.1826	530.8548
576.1467	664.6601	671.3077
782.1377	844.8291	1031.8328
1184.2731	1341.5574	1377.1712
1402.2187	1612.9540	1963.7628

Zero-point correction (Hartree): 0.037161

TS.CF3OCFCF2OH.CF3+HOCF2CFO.b

E(UM062X/CC-pVDZ) (Hartree): -864.10850576

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.578043	-0.338756	0.008103
C	0.613740	0.745809	-0.442607
O	-0.419402	0.511734	-1.071978

F	0.870948	1.949370	0.040126	
C	-1.849126	-0.088973	0.026421	
F	-1.317030	-0.086511	1.226706	
F	-2.140226	-1.303625	-0.372583	
F	-2.876785	0.712454	-0.058010	
F	2.772789	-0.136441	-0.600710	
O	1.090308	-1.571962	-0.246713	
F	1.803807	-0.234887	1.325928	
H	0.555285	-1.529897	-1.055084	
Rotational constants (GHz):	1.9127600	0.7309200	0.6589500	
Vibrational harmonic frequencies (cm-1):				
i557.8708	35.1067		40.6900	
51.7458	102.0955		163.8409	
207.4582	242.0278		280.0163	
325.2953	421.7072		444.3114	
511.3547	528.5437		542.2412	
603.2652	638.0907		681.6789	
716.4092	824.9535		1040.8790	
1113.7136	1180.9823		1282.8568	
1366.0766	1372.0545		1387.8648	
1454.7492	1577.8987		3816.4843	
Zero-point correction (Hartree):	0.052294			

TS.CF3OCFOCF2OH.CF2OH+CF3OCFO

E(UM062X/CC-pVDZ) (Hartree): -939.34915337

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.455216	-0.572011	-0.027808	
C	0.470431	0.989916	0.053244	
O	-0.737880	0.735443	-0.589470	
F	1.157256	1.719420	-0.845303	
C	-1.693806	-0.077919	-0.041016	
F	-1.715955	-0.054338	1.280807	
F	-1.510482	-1.336427	-0.434490	
F	-2.861251	0.335334	-0.503399	
F	1.378912	-0.978518	-1.271632	
O	2.699120	-0.276123	0.314632	
F	0.833343	-1.423385	0.777961	
H	2.664637	0.191275	1.168544	
O	0.589749	1.216932	1.261016	
Rotational constants (GHz):	1.5427200	0.7666300	0.6939300	
Vibrational harmonic frequencies (cm-1):				
i191.2982	19.9856		30.4818	
109.4788	161.5048		204.7069	
208.0590	273.6430		296.9906	
365.3265	408.5255		428.8971	
452.0698	457.4413		545.7483	
564.2543	568.3001		626.3329	
675.4160	735.3634		740.1965	
884.8728	1000.5867		1032.1709	
1192.9665	1218.8064		1290.5568	
1310.9766	1351.4774		1385.8185	
1459.2265	1593.8331		3769.0087	
Zero-point correction (Hartree):	0.057781			

TS.CF3OCFOCF2OH.CF3O+HOCF2CFO

E(UM062X/CC-pVDZ) (Hartree): -939.31671301

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.435234	-0.519615	-0.006451	
C	0.880623	0.916880	0.112298	
O	-0.818989	0.800588	-0.557271	
F	1.367435	1.680865	-0.851310	
C	-1.756448	-0.071515	-0.050436	
F	-1.739164	-0.099099	1.269135	
F	-1.566590	-1.302670	-0.511871	
F	-2.941106	0.347191	-0.474978	
F	1.288057	-0.961528	-1.249797	
O	2.757825	-0.550947	0.296411	

F	0.725405	-1.299684	0.830517	
H	2.853600	-0.420001	1.252652	
O	0.509115	1.397838	1.174562	
Rotational constants (GHz):	1.5575700	0.7348900	0.6657300	
Vibrational harmonic frequencies (cm-1):				
i632.0476		35.1574		60.6623
108.2176		131.7866		166.3456
222.0512		245.5953		264.6131
305.4907		379.7768		387.9266
429.7950		450.5346		496.5667
565.1058		590.3327		612.7000
652.2858		673.2861		695.6424
826.6604		930.9053		1139.2119
1150.1990		1226.3768		1284.6814
1316.3826		1330.8024		1367.2719
1429.7164		1608.4539		3825.6804
Zero-point correction (Hartree):	0.056750			

TS.CF3OCFOCF2OH.F+HOCHF2COOCH3

E(UM062X/CC-pVDZ) (Hartree): -939.30719846

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.800752	-0.345391	-0.024447	
C	0.391832	0.211881	0.262258	
O	-0.558815	-0.515209	-0.330466	
F	0.526627	1.747870	-0.758300	
C	-1.886715	-0.205650	-0.041757	
F	-2.153365	1.053392	-0.325325	
F	-2.152154	-0.435587	1.231722	
F	-2.618458	-0.997458	-0.794115	
F	1.947490	-0.586379	-1.328790	
O	2.763174	0.479078	0.410946	
F	1.850469	-1.545838	0.604193	
H	2.366151	1.119138	1.025012	
O	0.194786	1.010107	1.186294	
Rotational constants (GHz):	1.6596700	0.6410100	0.5965800	
Vibrational harmonic frequencies (cm-1):				
i635.6810		49.9108		63.5040
89.2412		131.4915		147.6127
201.9788		230.3687		236.8554
320.1073		350.0361		380.2717
434.6564		455.7166		519.3497
567.1603		594.8585		612.6590
652.4509		677.7699		767.8573
907.6717		932.1051		1123.2011
1202.6793		1222.4352		1308.9543
1343.2433		1379.4643		1393.5740
1479.7686		1629.2206		3784.3044
Zero-point correction (Hartree):	0.057388			

TS.CF3OCFOCFO.CF3OCFO+CFO

E(UM062X/CC-pVDZ) (Hartree): -838.91722482

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.857736	-0.559173	-0.111112	
C	0.577174	0.811581	0.163048	
O	-0.429416	0.412739	-0.706816	
F	1.216574	1.807261	-0.449126	
C	-1.559349	-0.126968	-0.134331	
F	-2.144419	0.736602	0.681322	
F	-1.294487	-1.236037	0.540724	
F	-2.379148	-0.412639	-1.127523	
F	1.334990	-1.702173	0.234743	
O	2.938802	-0.325927	-0.477604	
O	0.508743	0.726968	1.381059	
Rotational constants (GHz):	1.8674600	0.8498000	0.7293900	
Vibrational harmonic frequencies (cm-1):				
i378.6346		34.3271		66.4783
98.7611		153.6847		192.4636

199.2025	289.8169	378.8329
390.5581	419.9109	481.6961
556.2383	592.0313	630.2703
675.3103	696.2131	747.8336
901.5256	1011.6466	1153.2560
1233.0026	1290.9599	1323.4873
1365.6613	1631.5241	2036.5169

Zero-point correction (Hartree): 0.042263

TS.CF3OCFOCFO.CF3OCOCFO+F

E(UM062X/CC-pVDZ) (Hartree): -838.87849309

Electronic state : 2-A

Cartesian coordinates (Angs):

C	2.048073	0.197450	-0.230674
C	0.494052	-0.497121	0.004443
O	-0.398463	0.516318	0.066026
F	1.464165	-0.149493	1.386632
C	-1.734904	0.146342	-0.052206
F	-2.065960	-0.732044	0.875983
F	-1.979592	-0.368719	-1.243820
F	-2.432554	1.248802	0.105527
F	1.931076	1.500903	-0.223702
O	2.941697	-0.448144	-0.630835
O	0.319573	-1.640057	-0.239560

Rotational constants (GHz): 2.2024900 0.7208000 0.6727100

Vibrational harmonic frequencies (cm-1):

i445.3568	36.2862	72.4269
133.7687	160.7912	205.2082
260.6222	324.3828	378.1907
398.8609	435.3141	439.6874
556.4235	591.1264	623.6830
662.6422	678.8256	713.2768
846.0060	915.1009	1124.8898
1176.4417	1237.2151	1337.3931
1384.0305	1854.7569	1952.4264

Zero-point correction (Hartree): 0.042146

TS.CF3OCFOCFO.FCOCFO+CF3O

E(UM062X/CC-pVDZ) (Hartree): -838.88975913

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.624730	-0.622657	-0.153475
C	-0.999610	0.778646	-0.228256
O	0.503547	0.469536	0.749036
F	-1.581848	1.616811	0.609912
C	1.571019	-0.130350	0.123615
F	2.258920	0.711472	-0.624845
F	1.186232	-1.152732	-0.630164
F	2.342162	-0.574479	1.106891
F	-2.114718	-0.848576	1.055588
O	-1.668080	-1.386040	-1.045651
O	-0.397568	1.175716	-1.216853

Rotational constants (GHz): 1.7388500 0.8388600 0.7683600

Vibrational harmonic frequencies (cm-1):

i669.6136	27.4940	53.0315
91.5743	135.8066	155.0051
230.9810	259.8989	333.0859
414.6621	432.7594	483.3394
526.4895	614.4268	621.1924
666.6544	679.1187	756.3493
830.8795	934.3982	1177.1304
1236.4302	1312.9719	1329.5020
1352.8692	1595.5293	2011.9341

Zero-point correction (Hartree): 0.041607

TS.CF3OCFOHCF2.CF3OCFCF2+OH

E(UM062X/CC-pVDZ) (Hartree): -864.06137899

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.411821	0.563177	-0.210412
C	0.578794	-0.487691	-0.125044
O	-0.539000	-0.446809	0.646302
F	0.702469	-1.532234	-0.911126
C	-1.653741	0.072938	0.027938
F	-1.931570	-0.578163	-1.093366
F	-1.474679	1.355183	-0.275469
F	-2.655376	-0.041820	0.871319
F	1.259158	1.641698	0.511616
F	2.481420	0.572933	-0.950557
O	1.921931	-1.172309	1.396690
H	1.482511	-0.695972	2.129420

Rotational constants (GHz): 1.7121400 0.7791000 0.7166000

Vibrational harmonic frequencies (cm-1):

i318.5091	40.9742	70.4573
96.9759	111.7947	127.9691
196.7594	209.0177	264.8463
339.5696	359.0356	406.5740
466.5302	515.3255	546.4871
558.0087	639.1605	651.5491
702.0443	712.5300	846.8572
947.1249	1228.1840	1265.8331
1306.5812	1383.4519	1405.9371
1452.6538	1844.6972	3731.8852

Zero-point correction (Hartree): 0.051097

TS.CF3OCFOHCF2.CF3OCFOCF2H

E(UM062X/CC-pVDZ) (Hartree): -864.08151994

Electronic state : 2-A

Cartesian coordinates (Angs):

C	0.581099	-0.312717	0.167086
C	1.967578	0.269312	0.087384
F	2.858234	-0.406409	-0.613906
O	-0.401920	0.611610	-0.155025
F	2.081406	1.570169	-0.094982
F	0.408299	-1.441083	-0.548552
O	0.728547	-0.573196	1.514667
H	1.934541	-0.075213	1.366350
C	-1.700934	0.182881	-0.090145
F	-2.465115	1.251190	0.036530
F	-1.900612	-0.628154	0.945057
F	-2.052657	-0.464486	-1.194085

Rotational constants (GHz): 2.0970700 0.7205500 0.6744300

Vibrational harmonic frequencies (cm-1):

i2010.2961	41.7103	72.1776
103.8572	123.1684	209.1708
265.7493	344.6287	362.9963
399.5847	453.4509	524.1736
538.5823	612.8344	642.3207
668.3162	715.1296	779.5455
869.5129	952.8196	1074.5215
1169.6429	1193.1069	1227.3491
1308.2535	1317.6407	1355.2006
1368.4396	1504.2534	2038.0498

Zero-point correction (Hartree): 0.050658

TS.CF3OCFOHCF2.CF3OCOCF2+HF

E(UM062X/CC-pVDZ) (Hartree): -864.08520329

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.325520	0.756909	-0.183472
C	-0.766540	-0.536063	-0.455495
O	0.547711	-0.664279	-0.735124
F	-0.946929	-1.392721	1.177602
C	1.533193	0.012793	-0.039883
F	1.190624	0.237386	1.214762
F	1.799318	1.172158	-0.629300
F	2.611437	-0.739721	-0.080419

F	-2.596050	0.864898	0.039534	
F	-0.634075	1.794226	0.185189	
O	-1.540689	-1.454542	-0.917887	
H	-1.521903	-1.877315	0.130875	
Rotational constants (GHz):	1.7298600	0.9026100	0.7469200	
Vibrational harmonic frequencies (cm-1):				
i1347.2893	43.4268		97.9282	
106.7756	123.7823		174.2174	
212.5826	255.7267		358.7171	
372.3259	389.3421		462.1516	
506.5857	546.5458		598.4069	
627.2798	655.1259		728.1086	
777.6425	830.7319		924.6459	
994.5221	1219.0501		1270.9563	
1326.4221	1378.8582		1471.8888	
1509.5827	1663.1788		2194.3313	
Zero-point correction (Hartree):	0.049712			

TS.CF3OCFOHCF2.CF3OCOHC3

E(UM062X/CC-pVDZ) (Hartree): -864.08396191

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.253229	-0.572603	-0.253868	
C	0.678312	0.678579	-0.340708	
O	-0.628427	0.983919	-0.273744	
C	-1.593110	0.032835	-0.024984	
F	-1.419917	-0.540690	1.150336	
F	-1.591166	-0.902234	-0.970049	
F	-2.746941	0.663542	-0.044571	
F	2.514386	-0.730557	-0.512213	
F	0.645663	-1.672134	0.059158	
O	1.434538	1.724489	-0.628642	
H	2.011243	1.803889	0.161878	
F	1.432341	0.481626	1.514513	
Rotational constants (GHz):	1.8441400	0.8474900	0.7570000	
Vibrational harmonic frequencies (cm-1):				
i344.4161	30.2493		96.4331	
145.0258	158.9223		193.7578	
232.3611	277.2401		342.6229	
366.4745	381.7430		443.5963	
466.6893	511.6903		535.8137	
571.1586	628.5574		649.8572	
734.8371	815.8558		977.8304	
1216.1477	1262.7300		1284.4531	
1321.3611	1380.6372		1462.3312	
1486.7780	1774.3073		3635.8647	
Zero-point correction (Hartree):	0.053276			

TS.CFOCF2OH.CFOCFO+HF

E(UM062X/CC-pVDZ) (Hartree): -526.58036479

Electronic state : 1-A

Cartesian coordinates (Angs):

C	-0.523447	0.213250	-0.254719	
C	0.985572	0.077121	-0.043245	
O	1.727968	0.960271	0.190139	
F	1.344882	-1.188158	-0.181538	
F	-0.923618	1.433420	-0.031404	
O	-1.185264	-0.548049	-1.009632	
F	-1.050642	-0.687037	1.131659	
H	-1.449975	-1.064027	0.075273	
Rotational constants (GHz):	3.6475200	2.5055800	1.9820300	
Vibrational harmonic frequencies (cm-1):				
i1646.8304	63.8366		168.4411	
233.4023	324.1323		406.9839	
489.3905	637.5765		682.1138	
736.0815	829.3055		879.6634	
930.6562	1223.7996		1409.8650	
1683.6616	2009.6530		2039.5564	
Zero-point correction (Hartree):	0.033599			

TS.CF3+OCFCF2OH.CF4+OCFCFOH

E(UM062X/CC-pVDZ) (Hartree): -864.07045061

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.311918	0.684016	-0.012533
C	-1.982693	-0.629400	-0.170485
O	-2.288132	-1.182564	-1.168018
F	-2.140129	-1.182936	1.057406
C	2.102103	-0.141888	-0.004664
F	2.660478	0.932611	-0.500538
F	2.341103	-1.195030	-0.744156
F	2.461768	-0.338367	1.240927
F	-1.337908	1.447193	-1.074279
O	-1.466914	1.379967	1.105213
F	0.320776	0.135758	-0.002818
H	-1.559363	0.751358	1.839652

Rotational constants (GHz): 1.6336500 0.5942900 0.5647000

Vibrational harmonic frequencies (cm-1):

i1080.1972	12.3695	47.0225
51.5692	97.6063	132.2606
190.1267	196.8999	240.8138
299.8201	333.9486	394.7373
426.8272	512.1196	534.0525
535.5335	552.1276	642.4711
710.6455	760.0589	837.3126
1058.0286	1110.2091	1242.3046
1374.1614	1380.7016	1450.1824
1502.3223	1959.9874	3814.8382

Zero-point correction (Hartree): 0.051033

TS.CF3+OCFCF2OH.HCF3+OCFCF2O

E(UM062X/CC-pVDZ) (Hartree): -864.10213827

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.241590	0.903239	-0.111360
C	1.192367	-0.637745	-0.154407
F	2.324883	1.308292	0.534127
O	0.442537	1.636877	-0.574596
F	1.045777	-1.075743	1.111802
F	2.363231	-1.100336	-0.619552
O	0.233122	-1.121589	-0.978498
H	-0.944751	-0.457467	-0.640395
C	-1.871160	-0.039443	-0.023829
F	-2.430866	0.957557	-0.677499
F	-1.435730	0.375173	1.157683
F	-2.738106	-1.022847	0.138186

Rotational constants (GHz): 1.7577500 0.7239700 0.6307300

Vibrational harmonic frequencies (cm-1):

i810.9099	38.3740	49.5589
65.5190	91.3804	121.1996
164.9846	232.3184	241.8893
352.5742	423.5195	438.6320
518.5474	529.2110	610.4914
656.2512	710.2377	779.9419
808.7344	871.7337	1135.9724
1149.6999	1214.6565	1247.8860
1288.8744	1322.9989	1347.2995
1378.2232	1510.3796	1998.9832

Zero-point correction (Hartree): 0.048525

TS.CFOCF2OH+CF3.CF4+OCCF2OH

E(UM062X/CC-pVDZ) (Hartree): -864.05944840

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.054156	0.974810	-0.414156
C	1.726685	-0.352737	-0.011329
O	1.332864	2.041215	0.004675

F	1.769688	-1.196173	-1.044795	
F	0.928366	-0.926892	0.926436	
O	2.976502	-0.151577	0.446968	
H	3.009299	0.733101	0.847017	
F	-0.494664	0.584715	-0.669886	
C	-1.992557	-0.062071	0.032026	
F	-2.995742	0.528642	-0.555928	
F	-1.970855	-1.344723	-0.192197	
F	-1.927229	0.219963	1.303103	
Rotational constants (GHz):	1.9337200	0.6636600	0.6001200	
Vibrational harmonic frequencies (cm-1):				
i1130.7500	23.2924		38.5311	
43.7886	73.9340		150.4409	
181.4332	230.4892		257.4333	
325.6407	361.3560		379.6501	
450.2445	509.8560		537.9792	
539.8998	588.7911		635.6631	
704.0672	755.6148		892.5297	
1050.6757	1144.2081		1193.8425	
1341.1067	1402.4700		1408.2008	
1425.7377	1945.1367		3789.1528	
Zero-point correction (Hartree):	0.050988			

TS.CFOCF2O.CFO+CF2O

E(UM062X/CC-pVDZ) (Hartree): -525.95443095
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-1.039531	0.112075	-0.007761	
C	0.812832	0.189336	-0.008709	
F	-1.284151	-1.177678	-0.005331	
O	-1.797645	0.992345	0.002051	
F	1.073225	-0.646974	-1.026092	
F	1.051823	-0.494449	1.123204	
O	1.021661	1.390587	-0.092952	
Rotational constants (GHz):	3.8794700	2.4342100	2.0242400	
Vibrational harmonic frequencies (cm-1):				
i519.8411	14.3781		189.9886	
204.7994	315.8166		407.5797	
531.4701	569.2972		633.8294	
695.7399	923.2170		1094.6740	
1208.9550	1588.8385		2053.2226	
Zero-point correction (Hartree):	0.023765			

TS.CFOCF2OH+H2O.CFOCFO+HF+H2O.a

E(RM062X/CC-pVDZ) (Hartree): -603.03109006
Electronic state : 1-A
Cartesian coordinates (Angs):

C	-1.327952	0.394170	0.009209	
C	0.009839	-0.185469	-0.450234	
F	-2.117850	-0.544527	0.520156	
O	-1.631129	1.531128	-0.057210	
O	0.792385	0.515628	-1.100269	
F	-0.079502	-1.489217	-0.720886	
F	0.635827	-0.347518	1.135423	
H	2.066310	0.426815	-0.531168	
O	2.720935	0.282662	0.311275	
H	1.869993	-0.051087	0.914456	
H	2.968572	1.168080	0.620260	
Rotational constants (GHz):	3.1362200	1.3814200	1.2691300	
Vibrational harmonic frequencies (cm-1):				
i684.6103	33.6427		55.5926	
179.2564	230.1645		339.2953	
369.5372	433.1576		463.9999	
529.4191	541.7380		589.0019	
641.9971	704.7125		736.6035	
837.3321	895.7944		1166.1503	
1347.5157	1432.1526		1471.8233	
1651.6949	1754.7167		1963.3453	
2013.0355	2449.3264		3834.7792	

Zero-point correction (Hartree): 0.060749

TS.CFOCF2OH+H2O.CFOCFO+HF+H2O.b

E(RM062X/CC-pVDZ) (Hartree): -603.03151100

Electronic state : 1-A

Cartesian coordinates (Angs):

C	1.382404	-0.239831	-0.026009
C	0.032950	0.417663	-0.299537
F	1.269490	-1.561553	-0.141192
O	2.395243	0.303317	0.232485
O	-0.699984	-0.001504	-1.206104
F	0.095745	1.723851	-0.062151
F	-0.666538	-0.046317	1.191343
H	-2.002863	-0.157911	-0.711296
O	-2.692705	-0.352250	0.089467
H	-1.874789	-0.256884	0.815081
H	-2.923183	-1.292524	0.030710

Rotational constants (GHz): 3.0718100 1.4535200 1.2498600

Vibrational harmonic frequencies (cm-1):

i682.0105	32.6504	61.0537
181.1320	227.0616	344.2053
367.5631	422.4128	455.7721
528.7839	539.5117	586.5739
641.5147	706.6372	745.9918
840.0502	902.1744	1175.3860
1352.0150	1428.0299	1472.1001
1648.0759	1753.6226	1957.0270
2010.7923	2458.0927	3833.2516

Zero-point correction (Hartree): 0.060762

CF3OCF=CF2 + OH : M06-2X/aug-cc-pVTZ geometry

CF3

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -336.91604862
E(CCSD/Aug-CC-pVDZ) (Hartree): -336.89690736
T1 diagnostic: 0.016051
E(MP2/Aug-CC-pVDZ) (Hartree): -336.88703276
E(MP3/Aug-CC-pVDZ) (Hartree): -336.88631112
E(PMP2/Aug-CC-pVDZ) (Hartree): -336.88789619
E(PMP3/Aug-CC-pVDZ) (Hartree): -336.88682602
E(PUHF/Aug-CC-pVDZ) (Hartree): -336.17416151
E(UHF/Aug-CC-pVDZ) (Hartree): -336.17262351
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -336.93076910
T1 diagnostic: 0.000481
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -337.20696132
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -337.18381127
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -337.18824773
T1 diagnostic: 0.014568
D1 diagnostic: 0.048241
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -337.16509769
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -337.20664407
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -337.18349403
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -337.20849049
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -337.18534045
E(ROHF/AUG-CC-PVDZ) (Hartree): -336.17041988
E(UM062X/Aug-CC-pVTZ) (Hartree): -337.59568155
Electronic state : 2-A
Cartesian coordinates (Angs):
C -0.000050 -0.000085 0.325037
F -1.197846 -0.359473 -0.072333
F 0.910285 -0.857558 -0.072279
F 0.287595 1.217087 -0.072080
Rotational constants (GHz): 10.9547500 10.9536700 5.6693500
Vibrational harmonic frequencies (cm-1):
517.3827 517.6520 710.7345
1119.4420 1310.0219 1310.3422
Zero-point correction (Hartree): 0.012497

CF3O

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -411.98849576
E(CCSD/Aug-CC-pVDZ) (Hartree): -411.96407794
T1 diagnostic: 0.015334
E(MP2/Aug-CC-pVDZ) (Hartree): -411.94476522
E(MP3/Aug-CC-pVDZ) (Hartree): -411.95241651
E(PMP2/Aug-CC-pVDZ) (Hartree): -411.94654256
E(PMP3/Aug-CC-pVDZ) (Hartree): -411.95345976
E(PUHF/Aug-CC-pVDZ) (Hartree): -411.06670151
E(UHF/Aug-CC-pVDZ) (Hartree): -411.06361863
E(CCSD(T)/Aug-CC-pVTZ) (Hartree): -412.33420441
E(CCSD/Aug-CC-pVTZ) (Hartree): -412.29586238
T1 diagnostic: 0.014074
E(MP2/Aug-CC-pVTZ) (Hartree): -412.28571372
E(MP3/Aug-CC-pVTZ) (Hartree): -412.28768762
E(PMP2/Aug-CC-pVTZ) (Hartree): -412.28769056
E(PMP3/Aug-CC-pVTZ) (Hartree): -412.28880333
E(PUHF/Aug-CC-pVTZ) (Hartree): -411.18120075
E(UHF/Aug-CC-pVTZ) (Hartree): -411.17773468
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -411.99440504
T1 diagnostic: 0.000215
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -412.34303404
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -412.31374557
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -412.31930684
T1 diagnostic: 0.013568
D1 diagnostic: 0.044179
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -412.29001836
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -412.34260727

E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -412.31331880
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -412.34474919
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -412.31546072
E(ROHF/AUG-CC-PVDZ) (Hartree): -411.05905363
E(UM062X/Aug-CC-pVTZ) (Hartree): -412.82107529
Electronic state : 2-A
Cartesian coordinates (Angs):
C 0.006571 0.000343 0.033248
F -0.772785 -1.069473 -0.014373
F 0.843088 -0.010899 -0.994138
F -0.770794 1.070309 -0.035049
O 0.783124 0.011063 1.149069
Rotational constants (GHz): 6.0660100 5.8857900 5.6437400
Vibrational harmonic frequencies (cm-1):
267.1286 417.7304 588.9537
606.4694 629.1854 926.4697
1254.4370 1273.1352 1310.9422
Zero-point correction (Hartree): 0.016572

CFOCF2

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -449.99813181
E(CCSD/Aug-CC-pVDZ) (Hartree): -449.96486152
T1 diagnostic: 0.021485
E(MP2/Aug-CC-pVDZ) (Hartree): -449.94576231
E(MP3/Aug-CC-pVDZ) (Hartree): -449.94619591
E(PMP2/Aug-CC-pVDZ) (Hartree): -449.95357447
E(PMP3/Aug-CC-pVDZ) (Hartree): -449.95125759
E(PUHF/Aug-CC-pVDZ) (Hartree): -448.93111686
E(UHF/Aug-CC-pVDZ) (Hartree): -448.92124577
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -450.00806528
T1 diagnostic: 0.000468
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -450.38991069
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -450.35663922
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -450.35709843
T1 diagnostic: 0.017211
D1 diagnostic: 0.063304
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -450.32382695
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -450.38926694
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -450.35599546
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -450.39278884
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -450.35951736
E(ROHF/AUG-CC-PVDZ) (Hartree): -448.91221652
E(UM062X/Aug-CC-pVTZ) (Hartree): -450.93515226
Point group : CS
Electronic state : 2-A"
Cartesian coordinates (Angs):
C -0.155605 -0.774946 0.000000
C 0.000000 0.642754 0.000000
F 1.035812 -1.385601 0.000000
O -1.190350 -1.362945 0.000000
F 1.146662 1.248069 0.000000
F -1.020648 1.437167 0.000000
Rotational constants (GHz): 5.7371300 3.4339200 2.1481500
Vibrational harmonic frequencies (cm-1):
128.0493 (A") 205.8804 (A") 227.2438 (A')
415.7058 (A') 539.2289 (A') 615.8811 (A')
645.1764 (A") 827.1702 (A') 1156.3462 (A')
1447.0106 (A') 1565.4192 (A') 1828.4957 (A')
Zero-point correction (Hartree): 0.021874

CFOCF2OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -525.74929926
E(CCSD/Aug-CC-pVDZ) (Hartree): -525.70880050
T1 diagnostic: 0.015331
E(MP2/Aug-CC-pVDZ) (Hartree): -525.69623314
E(MP3/Aug-CC-pVDZ) (Hartree): -525.69243515
E(RHF/Aug-CC-pVDZ) (Hartree): -524.44082680
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -526.21072651
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -526.17054324

E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -526.17163048
 T1 diagnostic: 0.013498
 D1 diagnostic: 0.048046
 E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -526.13144721
 E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -526.20995940
 E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -526.16977613
 E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -526.21358603
 E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -526.17340276
 E(RHF/AUG-CC-PVDZ) (Hartree): -524.44082679
 E(RM062X/Aug-CC-pVTZ) (Hartree): -526.84003187
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 0.929047 0.134498 -0.000171
 C -0.607494 0.021071 -0.000078
 F 1.509983 -1.046438 -0.000278
 O 1.519779 1.153804 0.000040
 F -0.978647 -0.691862 1.078483
 F -0.979141 -0.694302 -1.076682
 O -1.192509 1.226434 -0.001584
 H -0.517234 1.918099 0.000146
 Rotational constants (GHz): 3.8203600 2.5446700 2.0827900
 Vibrational harmonic frequencies (cm-1):
 33.3716 210.5089 241.1353
 305.5466 397.5701 441.0268
 529.8981 613.8764 702.6013
 787.4974 836.8157 1102.3224
 1187.8995 1248.4435 1341.6037
 1456.1852 1962.9970 3803.7008
 Zero-point correction (Hartree): 0.039191

CFOCFO

 E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -425.47323729
 E(CCSD/Aug-CC-pVDZ) (Hartree): -425.43606739
 T1 diagnostic: 0.016755
 E(MP2/Aug-CC-pVDZ) (Hartree): -425.43116557
 E(MP3/Aug-CC-pVDZ) (Hartree): -425.42066496
 E(RHF/Aug-CC-pVDZ) (Hartree): -424.39234723
 E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -425.84644472
 E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -425.81388459
 E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -425.81073389
 T1 diagnostic: 0.015542
 D1 diagnostic: 0.046093
 E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -425.77817376
 E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -425.84556309
 E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -425.81300296
 E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -425.84950321
 E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -425.81694308
 E(RHF/AUG-CC-PVDZ) (Hartree): -424.39234723
 E(RM062X/Aug-CC-pVTZ) (Hartree): -426.36655918
 Point group : CS
 Electronic state : 1-A'
 Cartesian coordinates (Angs):
 C 0.000000 0.766517 0.000000
 C -0.000005 -0.766518 0.000000
 O -0.950757 1.452662 0.000000
 O 0.950759 -1.452655 0.000000
 F 1.239417 1.218915 0.000000
 F -1.239416 -1.218919 0.000000
 Rotational constants (GHz): 6.0128900 3.5767100 2.2426700
 Vibrational harmonic frequencies (cm-1):
 52.2591 (A'') 250.5281 (A') 422.5548 (A')
 458.8021 (A'') 536.7023 (A') 693.7135 (A')
 835.7030 (A'') 839.3270 (A') 1173.8611 (A')
 1335.5356 (A') 1976.6718 (A') 1977.9391 (A')
 Zero-point correction (Hartree): 0.024043

CFO

 E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -212.66782019
 E(CCSD/Aug-CC-pVDZ) (Hartree): -212.64990301

T1 diagnostic: 0.022821
E(MP2/Aug-CC-pVDZ) (Hartree): -212.64389186
E(MP3/Aug-CC-pVDZ) (Hartree): -212.63876694
E(PMP2/Aug-CC-pVDZ) (Hartree): -212.64672252
E(PMP3/Aug-CC-pVDZ) (Hartree): -212.64064141
E(PUHF/Aug-CC-pVDZ) (Hartree): -212.14496698
E(UHF/Aug-CC-pVDZ) (Hartree): -212.14114534
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -212.67245338
T1 diagnostic: 0.000793
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -212.85268665
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -212.83667697
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -212.83520252
T1 diagnostic: 0.023178
D1 diagnostic: 0.068730
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -212.81919284
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -212.85218889
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -212.83617920
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -212.85482458
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -212.83881490
E(ROHF/AUG-CC-PVDZ) (Hartree): -212.13632136
E(UM062X/Aug-CC-pVTZ) (Hartree): -213.11139379
Point group : CS
Electronic state : 2-A'
Cartesian coordinates (Angs):
C 0.000000 0.418366 0.000000
F -1.008812 -0.433873 0.000000
O 1.134913 0.174333 0.000000
Rotational constants (GHz): 192.1730700 11.6329400 10.9689500
Vibrational harmonic frequencies (cm-1):
650.9356 (A') 1091.2606 (A') 1982.8252 (A')
Zero-point correction (Hartree): 0.008486

F

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -99.55006944
E(CCSD/Aug-CC-pVDZ) (Hartree): -99.54774893
T1 diagnostic: 0.009722
E(MP2/Aug-CC-pVDZ) (Hartree): -99.53569906
E(MP3/Aug-CC-pVDZ) (Hartree): -99.54577504
E(PMP2/Aug-CC-pVDZ) (Hartree): -99.53706645
E(PMP3/Aug-CC-pVDZ) (Hartree): -99.54652791
E(PUHF/Aug-CC-pVDZ) (Hartree): -99.38373494
E(UHF/Aug-CC-pVDZ) (Hartree): -99.38109179
E(UM062X/Aug-CC-pVTZ) (Hartree): -99.73232189
Point group : OH
Cartesian coordinates (Angs):
F 0.000000 0.000000 0.000000
Zero-point correction (Hartree): 0.000000

HF

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -100.26362258
E(CCSD/Aug-CC-pVDZ) (Hartree): -100.25947091
T1 diagnostic: 0.012484
E(MP2/Aug-CC-pVDZ) (Hartree): -100.25577890
E(MP3/Aug-CC-pVDZ) (Hartree): -100.25644635
E(RHF/Aug-CC-pVDZ) (Hartree): -100.03333500
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -100.34962473
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -100.34247742
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -100.34547449
T1 diagnostic: 0.008822
D1 diagnostic: 0.014652
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -100.33832718
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -100.34965408
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -100.34250677
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -100.34978276
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -100.34263545
E(RHF/AUG-CC-PVDZ) (Hartree): -100.03333500
E(RM062X/Aug-CC-pVTZ) (Hartree): -100.45301335
Point group : C*V
Electronic state : 1-SG

Cartesian coordinates (Angs):
F 0.000000 0.000000 0.091997
H 0.000000 0.000000 -0.827973
Rotational constants (GHz): 0.0000000 623.9250717 623.9250717
Vibrational harmonic frequencies (cm-1):
4159.7924 (SG)
Zero-point correction (Hartree): 0.009477

O2

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -150.01974010
E(CCSD/Aug-CC-pVDZ) (Hartree): -150.00798598
T1 diagnostic: 0.017237
E(MP2/Aug-CC-pVDZ) (Hartree): -150.00187910
E(MP3/Aug-CC-pVDZ) (Hartree): -149.99995982
E(PMP2/Aug-CC-pVDZ) (Hartree): -150.00860413
E(PMP3/Aug-CC-pVDZ) (Hartree): -150.00309231
E(PUHF/Aug-CC-pVDZ) (Hartree): -149.65562822
E(UHF/Aug-CC-pVDZ) (Hartree): -149.64482689
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -150.03256052
T1 diagnostic: 0.001007
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -150.14485938
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -150.13379610
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -150.13259123
T1 diagnostic: 0.009707
D1 diagnostic: 0.016332
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -150.12152796
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -150.14458503
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -150.13352175
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -150.14450572
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -150.13344245
E(ROHF/AUG-CC-PVDZ) (Hartree): -149.62284252
E(UM062X/Aug-CC-pVTZ) (Hartree): -150.32479538

Point group : D*H

Electronic state : 3-SGG

Cartesian coordinates (Angs):
O 0.000000 0.000000 0.594925
O 0.000000 0.000000 -0.594925
Rotational constants (GHz): 0.0000000 44.6355313 44.6355313
Vibrational harmonic frequencies (cm-1):
1754.4580 (SGG)
Zero-point correction (Hartree): 0.003997

OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -75.58403081
E(CCSD/Aug-CC-pVDZ) (Hartree): -75.58066519
T1 diagnostic: 0.012131
E(MP2/Aug-CC-pVDZ) (Hartree): -75.56556372
E(MP3/Aug-CC-pVDZ) (Hartree): -75.57786120
E(PMP2/Aug-CC-pVDZ) (Hartree): -75.56732589
E(PMP3/Aug-CC-pVDZ) (Hartree): -75.57892392
E(PUHF/Aug-CC-pVDZ) (Hartree): -75.40650427
E(UHF/Aug-CC-pVDZ) (Hartree): -75.40357810
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -75.57183096
T1 diagnostic: 0.000130
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -75.64685987
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -75.64057201
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -75.64358240
T1 diagnostic: 0.008897
D1 diagnostic: 0.015262
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -75.63729455
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -75.64687263
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -75.64058478
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -75.64693076
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -75.64064291
E(ROHF/AUG-CC-PVDZ) (Hartree): -75.39925405
E(UM062X/Aug-CC-pVTZ) (Hartree): -75.73381011

Point group : C*V

Cartesian coordinates (Angs):
O 0.000000 0.000000 0.108021

```

H      0.000000      0.000000      -0.864170
Rotational constants (GHz):      0.0000000  563.9825141  563.9825141
Vibrational harmonic frequencies (cm-1):
  3766.2002 ( SG)
Zero-point correction (Hartree): 0.008580

FO2
-----
E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -249.57802139
E(CCSD/Aug-CC-pVDZ) (Hartree): -249.55129982
  T1 diagnostic: 0.040624
E(MP2/Aug-CC-pVDZ) (Hartree): -249.53506491
E(MP3/Aug-CC-pVDZ) (Hartree): -249.52903895
E(PMP2/Aug-CC-pVDZ) (Hartree): -249.53879473
E(PMP3/Aug-CC-pVDZ) (Hartree): -249.53124506
E(PUHF/Aug-CC-pVDZ) (Hartree): -248.94171267
E(UHF/Aug-CC-pVDZ) (Hartree): -248.93634014
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -249.56305878
  T1 diagnostic: 0.000845
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -249.78791795
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -249.76938471
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -249.76217378
  T1 diagnostic: 0.042129
  D1 diagnostic: 0.138980
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -249.74364053
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -249.78758819
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -249.76905495
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -249.79135691
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -249.77282367
E(ROHF/AUG-CC-PVDZ) (Hartree): -248.92908639
E(UM062X/Aug-CC-pVTZ) (Hartree): -250.06141196
Point group : CS
Electronic state : 2-A"
Cartesian coordinates (Angs):
  F      -0.990182      -0.605894      0.000000
  O       0.000000      0.548845      0.000000
  O       1.113954      0.132785      0.000000
Rotational constants (GHz):      81.4651100  11.4030800  10.0029200
Vibrational harmonic frequencies (cm-1):
  436.8240 ( A')      671.6464 ( A')      1523.0971 ( A')
Zero-point correction (Hartree): 0.005995

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CF4
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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -436.66486364
E(CCSD/Aug-CC-pVDZ) (Hartree): -436.64021317
  T1 diagnostic: 0.013058
E(MP2/Aug-CC-pVDZ) (Hartree): -436.63214204
E(MP3/Aug-CC-pVDZ) (Hartree): -436.62942215
E(RHF/Aug-CC-pVDZ) (Hartree): -435.69485173
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -437.04452276
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -437.01456001
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -437.02052419
  T1 diagnostic: 0.010845
  D1 diagnostic: 0.027692
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -436.99056143
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -437.04412895
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -437.01416620
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -437.04602440
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -437.01606165
E(RHF/AUG-CC-PVDZ) (Hartree): -435.69485172
E(RM062X/Aug-CC-pVTZ) (Hartree): -437.53902683
Point group : CS
Electronic state : 1-A'
Cartesian coordinates (Angs):
  C      -0.000011      0.000013      0.000000
  F      -1.075299      0.760407      0.000000
  F       1.075327      0.760333      0.000000
  F      -0.000011      -0.760374      1.075325
  F      -0.000011      -0.760374      -1.075325
Rotational constants (GHz):      5.7513200  5.7512700  5.7511800

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Vibrational harmonic frequencies (cm-1):

438.8158 (A')	439.1734 (A'')	636.8057 (A')
637.0667 (A'')	637.0717 (A')	941.5475 (A')
1320.6264 (A'')	1320.7493 (A')	1321.2840 (A')

Zero-point correction (Hartree): 0.017526

HCF3

E(CCS(T)/Aug-CC-pVDZ) (Hartree): -337.59215799
E(CCS(D)/Aug-CC-pVDZ) (Hartree): -337.57256318
T1 diagnostic: 0.013587
E(MP2/Aug-CC-pVDZ) (Hartree): -337.56049378
E(MP3/Aug-CC-pVDZ) (Hartree): -337.56277752
E(RHF/Aug-CC-pVDZ) (Hartree): -336.81510309
E(CCS(D)(T)-F12a/AUG-CC-PVDZ) (Hartree): -337.88713544
E(CCS(D)(T)-F12b/AUG-CC-PVDZ) (Hartree): -337.86291101
E(CCS(D)-F12a/AUG-CC-PVDZ) (Hartree): -337.86802468
T1 diagnostic: 0.011165
D1 diagnostic: 0.027248
E(CCS(D)-F12b/AUG-CC-PVDZ) (Hartree): -337.84380025
E(CCS(D)-T-F12a/AUG-CC-PVDZ) (Hartree): -337.88684220
E(CCS(D)-T-F12b/AUG-CC-PVDZ) (Hartree): -337.86261776
E(CCS(D)[T]-F12a/AUG-CC-PVDZ) (Hartree): -337.88837738
E(CCS(D)[T]-F12b/AUG-CC-PVDZ) (Hartree): -337.86415295
E(RHF/AUG-CC-PVDZ) (Hartree): -336.81510309
E(RM062X/Aug-CC-pVTZ) (Hartree): -338.27266965
Electronic state : 1-A
Cartesian coordinates (Angs):

C	0.000087	-0.000124	0.340583
F	-0.798904	-0.956222	-0.128595
F	1.227729	-0.213570	-0.128578
F	-0.428866	1.169883	-0.128547
H	-0.000157	-0.000077	1.427986

Rotational constants (GHz): 10.4024900 10.3988600 5.7107400
Vibrational harmonic frequencies (cm-1):

515.0152	515.2700	712.2913
1172.3039	1200.9083	1201.6049
1414.1147	1414.2983	3178.9846

Zero-point correction (Hartree): 0.025800

FNO

E(CCS(D)(T)/Aug-CC-pVDZ) (Hartree): -229.26335692
E(CCS(D)/Aug-CC-pVDZ) (Hartree): -229.23913560
T1 diagnostic: 0.022157
E(MP2/Aug-CC-pVDZ) (Hartree): -229.23987876
E(MP3/Aug-CC-pVDZ) (Hartree): -229.22602153
E(RHF/Aug-CC-pVDZ) (Hartree): -228.65800609
E(CCS(D)(T)-F12a/AUG-CC-PVDZ) (Hartree): -229.45853503
E(CCS(D)(T)-F12b/AUG-CC-PVDZ) (Hartree): -229.44115566
E(CCS(D)-F12a/AUG-CC-PVDZ) (Hartree): -229.43538718
T1 diagnostic: 0.021528
D1 diagnostic: 0.060852
E(CCS(D)-F12b/AUG-CC-PVDZ) (Hartree): -229.41800781
E(CCS(D)-T-F12a/AUG-CC-PVDZ) (Hartree): -229.45773443
E(CCS(D)-T-F12b/AUG-CC-PVDZ) (Hartree): -229.44035505
E(CCS(D)[T]-F12a/AUG-CC-PVDZ) (Hartree): -229.46128411
E(CCS(D)[T]-F12b/AUG-CC-PVDZ) (Hartree): -229.44390474
E(RHF/AUG-CC-PVDZ) (Hartree): -228.65800609
E(RM062X/Aug-CC-pVTZ) (Hartree): -229.72011450
Point group : CS
Electronic state : 1-A'
Cartesian coordinates (Angs):

F	-0.940989	-0.579405	0.000000
N	0.000000	0.551202	0.000000
O	1.058613	0.169529	0.000000

Rotational constants (GHz): 97.6744100 12.4354500 11.0310300
Vibrational harmonic frequencies (cm-1):

580.8086 (A')	827.3994 (A')	1981.8837 (A')
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Zero-point correction (Hartree): 0.007723

NO

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -129.62340013
E(CCSD/Aug-CC-pVDZ) (Hartree): -129.61043015
T1 diagnostic: 0.022643
E(MP2/Aug-CC-pVDZ) (Hartree): -129.60198234
E(MP3/Aug-CC-pVDZ) (Hartree): -129.60094805
E(PMP2/Aug-CC-pVDZ) (Hartree): -129.60591473
E(PMP3/Aug-CC-pVDZ) (Hartree): -129.60340171
E(PUHF/Aug-CC-pVDZ) (Hartree): -129.27660045
E(UHF/Aug-CC-pVDZ) (Hartree): -129.27121024
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -129.61991150
T1 diagnostic: 0.000669
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -129.73304160
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -129.72263170
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -129.72039356
T1 diagnostic: 0.020809
D1 diagnostic: 0.045520
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -129.70998366
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -129.73269619
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -129.72228629
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -129.73424813
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -129.72383822
E(ROHF/AUG-CC-PVDZ) (Hartree): -129.26419668
E(UM062X/Aug-CC-pVTZ) (Hartree): -129.89335138
Point group : C*v
Cartesian coordinates (Angs):
N 0.000000 0.000000 -0.606410
O 0.000000 0.000000 0.530609
Rotational constants (GHz): 0.0000000 52.3561998 52.3561998
Vibrational harmonic frequencies (cm-1):
2066.0391 (SG)
Zero-point correction (Hartree): 0.004707

CF2O

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -312.40224471
E(CCSD/Aug-CC-pVDZ) (Hartree): -312.37961809
T1 diagnostic: 0.015319
E(MP2/Aug-CC-pVDZ) (Hartree): -312.37639437
E(MP3/Aug-CC-pVDZ) (Hartree): -312.36978218
E(RHF/Aug-CC-pVDZ) (Hartree): -311.65523696
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -312.67659354
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -312.65401106
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -312.65478245
T1 diagnostic: 0.014042
D1 diagnostic: 0.045636
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -312.63219997
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -312.67611355
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -312.65353107
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -312.67831177
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -312.65572929
E(RHF/AUG-CC-PVDZ) (Hartree): -311.65523696
E(RM062X/Aug-CC-pVTZ) (Hartree): -313.04496676
Point group : C2v
Electronic state : 1-A1
Cartesian coordinates (Angs):
C 0.000000 0.000000 0.141685
O 0.000000 0.000000 1.308682
F 0.000000 1.057949 -0.628865
F 0.000000 -1.057949 -0.628865
Rotational constants (GHz): 11.8834000 11.8530400 5.9341000
Vibrational harmonic frequencies (cm-1):
589.8658 631.4124 797.9617
1010.4306 1300.3034 2020.6839
Zero-point correction (Hartree): 0.014468

H2O

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -76.27377604
E(CCSD/Aug-CC-pVDZ) (Hartree): -76.26855934

T1 diagnostic: 0.012284
E(MP2/Aug-CC-pVDZ) (Hartree): -76.26079349
E(MP3/Aug-CC-pVDZ) (Hartree): -76.26553289
E(RHF/Aug-CC-pVDZ) (Hartree): -76.04133056
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -76.34547905
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -76.33832353
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -76.34048177
T1 diagnostic: 0.009595
D1 diagnostic: 0.017559
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -76.33332624
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -76.34547247
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -76.33831694
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -76.34566631
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -76.33851078
E(RHF/AUG-CC-PVDZ) (Hartree): -76.04133056
E(RM062X/Aug-CC-pVTZ) (Hartree): -76.43010673
Point group : C2V
Electronic state : 1-A1
Cartesian coordinates (Angs):
O 0.000000 0.000000 0.116390
H 0.000000 0.762470 -0.465561
H 0.000000 -0.762470 -0.465561
Rotational constants (GHz): 833.6308100 431.2772300 284.2309300
Vibrational harmonic frequencies (cm-1):
1619.5204 3869.6600 3972.6125
Zero-point correction (Hartree): 0.021556

CF3OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -412.68034333
E(CCSD/Aug-CC-pVDZ) (Hartree): -412.65448293
T1 diagnostic: 0.013152
E(MP2/Aug-CC-pVDZ) (Hartree): -412.64325767
E(MP3/Aug-CC-pVDZ) (Hartree): -412.64377238
E(RHF/Aug-CC-pVDZ) (Hartree): -411.70621449
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -413.04287317
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -413.01274678
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -413.01779508
T1 diagnostic: 0.010914
D1 diagnostic: 0.029502
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -412.98766869
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -413.04249503
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -413.01236864
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -413.04432387
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -413.01419748
E(RHF/AUG-CC-PVDZ) (Hartree): -411.70621449
E(RM062X/Aug-CC-pVTZ) (Hartree): -413.51933257
Electronic state : 1-A
Cartesian coordinates (Angs):
C -0.004770 0.022374 0.000027
F 0.264983 -0.731883 -1.071617
F 0.269124 -0.728063 1.073263
F 0.807798 1.062014 -0.003307
O -1.271624 0.466482 0.001623
H -1.875528 -0.284719 0.001796
Rotational constants (GHz): 5.7367500 5.6817300 5.6378500
Vibrational harmonic frequencies (cm-1):
251.7883 445.3654 455.5775
608.3323 630.1886 642.8287
931.0605 1132.7543 1223.7762
1326.0300 1432.5993 3857.2159
Zero-point correction (Hartree): 0.029474

CF3OCF2

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -787.00837175
E(CCSD/Aug-CC-pVDZ) (Hartree): -786.95343314
T1 diagnostic: 0.013967
E(MP2/Aug-CC-pVDZ) (Hartree): -786.93380512
E(MP3/Aug-CC-pVDZ) (Hartree): -786.93357391
E(RHF/Aug-CC-pVDZ) (Hartree): -785.14914490

E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -787.69440603
 E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -787.63710232
 E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -787.64116203
 T1 diagnostic: 0.012171
 D1 diagnostic: 0.036034
 E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -787.58385832
 E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -787.69354502
 E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -787.63624131
 E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -787.69767868
 E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -787.64037497
 E(RHF/AUG-CC-PVDZ) (Hartree): -785.14914490
 E(RM062X/Aug-CC-pVTZ) (Hartree): -788.62828972
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 1.688942 -0.285791 -0.062739
 C 0.700416 0.547437 -0.322426
 O -0.480816 0.152285 -0.855667
 F 0.829571 1.851845 -0.147668
 C -1.486193 -0.110863 0.040109
 F -1.715822 0.926726 0.833412
 F -1.185912 -1.150646 0.807160
 F -2.568138 -0.380648 -0.654268
 F 1.609467 -1.579624 -0.233086
 F 2.856115 0.096460 0.385080
 Rotational constants (GHz): 2.3722300 0.9247100 0.7685300
 Vibrational harmonic frequencies (cm-1):
 37.5078 74.7467 120.8240
 198.0165 213.2699 344.4227
 369.4525 445.6470 470.1532
 544.6117 557.0763 614.0070
 643.3737 651.5821 727.8200
 849.3273 945.4895 1219.7324
 1240.0152 1268.3436 1345.5522
 1364.9966 1401.8671 1937.4241
 Zero-point correction (Hartree): 0.040062

CF3OCFCF2OH.b

 E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.67940195
 E(CCSD/Aug-CC-pVDZ) (Hartree): -862.61985276
 T1 diagnostic: 0.015011
 E(MP2/Aug-CC-pVDZ) (Hartree): -862.59333768
 E(MP3/Aug-CC-pVDZ) (Hartree): -862.59527210
 E(PMP2/Aug-CC-pVDZ) (Hartree): -862.59481043
 E(PMP3/Aug-CC-pVDZ) (Hartree): -862.59616170
 E(PUHF/Aug-CC-pVDZ) (Hartree): -860.62487051
 E(UHF/Aug-CC-pVDZ) (Hartree): -860.62247052
 E(UM062X/Aug-CC-pVTZ) (Hartree): -864.45363994
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C 1.518734 -0.252099 -0.020518
 C 0.533475 0.767904 -0.565023
 O -0.712525 0.387214 -0.947776
 F 0.570982 1.928007 0.060637
 C -1.599449 -0.095831 -0.027759
 F -1.644657 0.650552 1.064778
 F -1.267976 -1.337953 0.353610
 F -2.783372 -0.136948 -0.589199
 F 1.450197 -1.344777 -0.808124
 O 1.345658 -0.587610 1.287134
 F 2.740312 0.275692 -0.126404
 H 0.629000 -1.227850 1.367257
 Rotational constants (GHz): 1.9736900 0.8326600 0.7393400
 Vibrational harmonic frequencies (cm-1):
 30.0127 73.9088 118.6704
 161.3775 201.9144 217.8709
 249.0447 325.6431 356.8960
 430.3517 479.6864 510.7112
 554.5476 586.4402 630.1259
 638.1139 688.3116 756.0145
 820.2984 918.7688 1113.7329

1133.7301	1201.5724	1230.6017
1244.8910	1331.3237	1348.5507
1380.4784	1422.2351	3850.6365

Zero-point correction (Hartree): 0.054691

CF3OCFCF2OH

E (CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.67938565
E (CCSD/Aug-CC-pVDZ) (Hartree): -862.61998113
T1 diagnostic: 0.014933
E (MP2/Aug-CC-pVDZ) (Hartree): -862.59347533
E (MP3/Aug-CC-pVDZ) (Hartree): -862.59540270
E (PMP2/Aug-CC-pVDZ) (Hartree): -862.59495995
E (PMP3/Aug-CC-pVDZ) (Hartree): -862.59629719
E (PUHF/Aug-CC-pVDZ) (Hartree): -860.62565487
E (UHF/Aug-CC-pVDZ) (Hartree): -860.62324094
E (UM062X/Aug-CC-pVTZ) (Hartree): -864.45383120
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.510545	-0.281409	-0.045150
C	0.536932	0.730124	-0.609555
O	-0.721333	0.397795	-0.950609
F	0.651089	1.923102	-0.032890
C	-1.599479	-0.095959	-0.008830
F	-1.529678	0.589442	1.122669
F	-1.353475	-1.368017	0.254568
F	-2.808913	0.011401	-0.513900
F	1.456386	-1.392563	-0.775455
O	2.787224	0.167227	-0.049836
F	1.132515	-0.608677	1.223468
H	2.853567	0.971113	0.478627

Rotational constants (GHz): 1.9671300 0.8517200 0.7523600
Vibrational harmonic frequencies (cm-1):

56.1864	64.2653	100.0778
145.6761	195.6141	217.2406
231.8838	333.9885	355.5384
430.3594	473.7716	518.6328
551.6080	588.2199	631.7813
635.5767	688.9757	756.4629
817.0888	919.8916	1076.6466
1115.4015	1195.0133	1252.5595
1279.5839	1324.2073	1334.9484
1380.0539	1445.4786	3851.2739

Zero-point correction (Hartree): 0.054603

CF3OCFCFOH

E (CCSD(T)/Aug-CC-pVDZ) (Hartree): -763.03065280
E (CCSD/Aug-CC-pVDZ) (Hartree): -762.97447156
T1 diagnostic: 0.014030
E (MP2/Aug-CC-pVDZ) (Hartree): -762.95275856
E (MP3/Aug-CC-pVDZ) (Hartree): -762.95492937
E (RHF/Aug-CC-pVDZ) (Hartree): -761.16788487
E (CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -763.70066848
E (CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -763.64327573
E (CCSD-F12a/AUG-CC-PVDZ) (Hartree): -763.64634146
T1 diagnostic: 0.012242
D1 diagnostic: 0.039870
E (CCSD-F12b/AUG-CC-PVDZ) (Hartree): -763.58894871
E (CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -763.69982780
E (CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -763.64243505
E (CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -763.70385615
E (CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -763.64646340
E (RHF/AUG-CC-PVDZ) (Hartree): -761.16788486
E (RM062X/Aug-CC-pVTZ) (Hartree): -764.61669024
Electronic state : 1-A
Cartesian coordinates (Angs):

C	1.697303	-0.314127	-0.065234
C	0.696456	0.508023	-0.332742
O	-0.498166	0.147337	-0.860605
F	0.864571	1.827593	-0.167327

C	-1.498558	-0.097117	0.042375	
F	-1.698955	0.942592	0.843706	
F	-1.218999	-1.146717	0.803783	
F	-2.595628	-0.337827	-0.641731	
F	1.582652	-1.611350	-0.229121	
O	2.906087	0.023870	0.381457	
H	2.962662	0.981061	0.483004	
Rotational constants (GHz):	2.3678000	0.9378600	0.7770100	
Vibrational harmonic frequencies (cm-1):				
39.3526		75.9167		113.6979
196.8963		210.8595		333.3735
343.3296		369.5487		430.9537
470.1321		546.0914		561.8221
633.2306		642.1482		653.3715
718.7025		848.4312		945.1177
1141.9258		1241.8175		1261.1224
1275.7995		1340.5050		1347.0802
1453.5124		1912.3119		3852.1868
Zero-point correction (Hartree):	0.052305			

CF3OCFCFO

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-----
E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -762.41668276
E(CCSD/Aug-CC-pVDZ) (Hartree): -762.36085840
  T1 diagnostic: 0.018660
E(MP2/Aug-CC-pVDZ) (Hartree): -762.33613262
E(MP3/Aug-CC-pVDZ) (Hartree): -762.33453882
E(PMP2/Aug-CC-pVDZ) (Hartree): -762.34370037
E(PMP3/Aug-CC-pVDZ) (Hartree): -762.33949247
E(PUHF/Aug-CC-pVDZ) (Hartree): -760.59888672
E(UHF/Aug-CC-pVDZ) (Hartree): -760.58936602
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -762.43563794
  T1 diagnostic: 0.000371
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -763.08010394
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -763.02364568
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -763.02552988
  T1 diagnostic: 0.015132
  D1 diagnostic: 0.062102
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -762.96907162
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -763.07905991
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -763.02260164
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -763.08425775
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -763.02779948
E(ROHF/AUG-CC-PVDZ) (Hartree): -760.58090571
E(UM062X/Aug-CC-pVTZ) (Hartree): -763.99861019

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Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.873488	-0.283003	-0.005999	
C	0.732139	0.553543	-0.268754	
O	-0.410369	0.124131	-0.815363	
F	0.856461	1.843711	-0.142807	
C	-1.471029	-0.111747	0.038604	
F	-1.800357	0.988057	0.697154	
F	-1.163206	-1.046276	0.922696	
F	-2.479154	-0.508567	-0.697243	
F	1.582952	-1.570207	-0.221660	
O	2.938138	0.086718	0.364567	
Rotational constants (GHz):	2.4015400	0.9527600	0.7876000	
Vibrational harmonic frequencies (cm-1):				
30.8096		54.9156		79.5883
148.8373		212.8933		327.6208
357.1783		380.5919		465.8821
530.7936		567.1727		638.9214
665.2909		669.9304		703.4067
832.0471		943.3547		1149.2375
1208.2625		1280.0836		1354.9931
1409.3401		1503.0235		1845.4521
Zero-point correction (Hartree):	0.039548			

CF3OCFOCF2OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -937.74545787
E(CCSD/Aug-CC-pVDZ) (Hartree): -937.67914431
T1 diagnostic: 0.017289
E(MP2/Aug-CC-pVDZ) (Hartree): -937.64401159
E(MP3/Aug-CC-pVDZ) (Hartree): -937.65164978
E(PMP2/Aug-CC-pVDZ) (Hartree): -937.64616276
E(PMP3/Aug-CC-pVDZ) (Hartree): -937.65300845
E(PUHF/Aug-CC-pVDZ) (Hartree): -935.50218141
E(UHF/Aug-CC-pVDZ) (Hartree): -935.49864713
E(UM062X/Aug-CC-pVTZ) (Hartree): -939.67178872
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.380403	-0.521443	-0.023440
C	0.565405	0.857434	-0.000780
O	-0.721756	0.778985	-0.534941
F	1.203447	1.683136	-0.848147
C	-1.680708	-0.058606	-0.042483
F	-1.699462	-0.104059	1.281207
F	-1.527345	-1.295254	-0.497423
F	-2.840391	0.401564	-0.471954
F	1.241341	-1.033102	-1.236549
O	2.689990	-0.343433	0.214143
F	0.811390	-1.355972	0.863406
H	2.823771	0.113820	1.053303
O	0.642367	1.258830	1.262304

Rotational constants (GHz): 1.5690500 0.7777300 0.7075900
Vibrational harmonic frequencies (cm-1):

41.5964	65.2281	117.6917
158.8565	197.5291	220.4649
248.5576	317.1541	337.4952
349.6156	373.1564	432.1009
467.0354	521.1334	536.1744
585.5829	625.3223	668.9336
692.3582	731.0070	778.8964
867.7167	974.3913	1091.4635
1149.4329	1203.4483	1255.8633
1272.0385	1277.2491	1300.5400
1342.0130	1413.1134	3840.4384

Zero-point correction (Hartree): 0.057988

CF3OCFOCFO

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -837.47280392
E(CCSD/Aug-CC-pVDZ) (Hartree): -837.41044764
T1 diagnostic: 0.016370
E(MP2/Aug-CC-pVDZ) (Hartree): -837.38394495
E(MP3/Aug-CC-pVDZ) (Hartree): -837.38469304
E(PMP2/Aug-CC-pVDZ) (Hartree): -837.38594288
E(PMP3/Aug-CC-pVDZ) (Hartree): -837.38592337
E(PUHF/Aug-CC-pVDZ) (Hartree): -835.46096970
E(UHF/Aug-CC-pVDZ) (Hartree): -835.45766747
E(UM062X/Aug-CC-pVTZ) (Hartree): -839.20080563
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.555107	-0.553596	0.169987
C	0.720907	0.725950	-0.137611
O	-0.504024	0.459298	-0.754731
F	1.405239	1.422662	-1.063895
C	-1.483177	-0.139469	-0.013081
F	-2.090763	0.717269	0.789941
F	-0.985749	-1.126159	0.734756
F	-2.363411	-0.636134	-0.852945
F	1.519219	-1.354387	-0.881463
O	2.157596	-0.771354	1.148906
O	0.581698	1.386236	1.024160

Rotational constants (GHz): 1.7520800 0.9117100 0.8377000
Vibrational harmonic frequencies (cm-1):

48.7130	56.8937	87.6194
161.1413	214.2614	231.8856
327.8106	357.1793	375.8044
435.9079	486.7333	516.3118

621.8567	643.5234	679.5603
716.0841	791.8045	817.8198
940.6982	1073.4457	1169.4076
1200.0343	1262.6999	1276.5884
1285.8286	1351.7603	1977.0995

Zero-point correction (Hartree): 0.043532

CF3OCFOHCF2

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.67823901
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.61849154
T1 diagnostic: 0.015042
E(MP2/Aug-CC-pVDZ) (Hartree): -862.59266732
E(MP3/Aug-CC-pVDZ) (Hartree): -862.59369205
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.59389342
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.59443402
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.62203596
E(UHF/Aug-CC-pVDZ) (Hartree): -860.61996547
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.45173426
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-1.489195	0.654573	0.099178
C	-0.883855	-0.731683	-0.002240
O	0.415719	-0.800292	-0.558508
F	-0.816477	-1.234597	1.254094
C	1.417062	0.008786	-0.116461
F	1.555906	-0.016787	1.201120
F	1.216253	1.272691	-0.480181
F	2.533853	-0.416748	-0.674872
F	-1.564058	1.306421	-1.037024
F	-1.022788	1.406063	1.070653
O	-1.686380	-1.470557	-0.795468
H	-1.222976	-2.276653	-1.055168

Rotational constants (GHz): 1.5627000 0.9837300 0.8717900
Vibrational harmonic frequencies (cm-1):

56.5679	85.7806	130.8114
157.6260	206.2864	223.6211
296.7746	350.1205	367.4183
400.5865	449.6121	503.7303
552.3728	582.8158	612.7878
659.3554	682.7016	751.8294
822.5844	930.1136	1073.3260
1123.2349	1239.6282	1270.2941
1275.8182	1327.0924	1334.4584
1349.5862	1490.0589	3828.4442

Zero-point correction (Hartree): 0.054985

CF3OCOHC2F2

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -763.02985823
E(CCSD/Aug-CC-pVDZ) (Hartree): -762.97351005
T1 diagnostic: 0.013960
E(MP2/Aug-CC-pVDZ) (Hartree): -762.95168113
E(MP3/Aug-CC-pVDZ) (Hartree): -762.95394099
E(RHF/Aug-CC-pVDZ) (Hartree): -761.16563384
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -763.69972079
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -763.64233447
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -763.64523972
T1 diagnostic: 0.012156
D1 diagnostic: 0.035790
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -763.58785340
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -763.69887598
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -763.64148966
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -763.70287446
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -763.64548814
E(RHF/AUG-CC-PVDZ) (Hartree): -761.16563383
E(RM062X/Aug-CC-pVTZ) (Hartree): -764.61538532
Electronic state : 1-A
Cartesian coordinates (Angs):

C	1.678096	-0.291577	-0.055468
C	0.710327	0.567035	-0.323176

O	-0.473649	0.100157	-0.866355	
C	-1.465067	-0.130761	0.028176	
F	-1.678083	0.942631	0.807627	
F	-1.188354	-1.146419	0.836749	
F	-2.568124	-0.396081	-0.635380	
F	2.850439	0.057041	0.409667	
F	1.589966	-1.582927	-0.254277	
O	0.869225	1.899256	-0.231098	
H	0.242670	2.268319	0.402951	
Rotational constants (GHz):	2.3772400	0.9333700	0.7776800	
Vibrational harmonic frequencies (cm-1):				
47.4401	76.4699		127.1506	
204.4254	216.0984		297.9877	
355.5515	370.6517		444.6534	
479.0097	541.0661		562.9832	
625.1109	646.9510		653.3066	
720.3848	844.9401		939.5104	
1171.1641	1212.0809		1262.7407	
1294.9651	1346.2558		1350.8466	
1391.3225	1919.8654		3817.5206	
Zero-point correction (Hartree):	0.052217			

complex.CF3OCFCF2.OH.a

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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.59807137
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.53797614
  T1 diagnostic: 0.016079
E(MP2/Aug-CC-pVDZ) (Hartree): -862.50298195
E(MP3/Aug-CC-pVDZ) (Hartree): -862.51390028
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.50576865
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.51582359
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.55092244
E(UHF/Aug-CC-pVDZ) (Hartree): -860.54721272
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.36843635
Electronic state : 2-A
Cartesian coordinates (Angs):
  C    1.418486    0.500640   -0.330002
  C    0.517285   -0.464442   -0.227240
  O   -0.541192   -0.389634    0.618146
  F    0.637440   -1.600055   -0.883083
  C   -1.717468    0.068731    0.075703
  F   -2.083036   -0.655524   -0.971429
  F   -1.602458    1.329406   -0.318686
  F   -2.635448   -0.013519    1.011633
  F    1.324506    1.637659    0.304185
  F    2.478816    0.415204   -1.078209
  H    1.606075   -0.612321    2.277499
  O    2.291909   -0.864840    1.635857
Rotational constants (GHz):  1.6907800    0.7131500    0.6825900
Vibrational harmonic frequencies (cm-1):
  34.2782          45.7880          63.8957
  78.2298          121.4535          135.6739
  164.3050         200.6411          216.1293
  344.2503         368.2670          453.0312
  464.7562         473.9792          544.9255
  558.4325         602.9293          644.2961
  652.0892         725.0745          848.4207
  946.5575         1221.9930         1237.9249
  1272.8852        1344.6454         1371.4613
  1417.4671        1893.5131         3758.5733
Zero-point correction (Hartree): 0.050589

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complex.CF3OCFCF2.OH.b

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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.59820785
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.53817586
  T1 diagnostic: 0.015977
E(MP2/Aug-CC-pVDZ) (Hartree): -862.50282845
E(MP3/Aug-CC-pVDZ) (Hartree): -862.51411269
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.50592076
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.51628754

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E(PUHF/Aug-CC-pVDZ) (Hartree): -860.55195866
E(UHF/Aug-CC-pVDZ) (Hartree): -860.54803000
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.36919563
Electronic state : 2-A
Cartesian coordinates (Angs):
C 1.586220 -0.171660 -0.488051
C 0.565543 0.670581 -0.459746
O -0.655646 0.365890 -0.966548
F 0.701298 1.910311 -0.033530
C -1.596676 -0.082121 -0.081446
F -1.739527 0.735201 0.949257
F -1.267410 -1.284038 0.399233
F -2.733310 -0.182973 -0.729610
F 1.486623 -1.400739 -0.920015
F 2.793714 0.149499 -0.127497
H 0.313624 -1.090926 2.008692
O 1.053567 -0.460292 2.007326
Rotational constants (GHz): 1.6995100 0.7728000 0.7418400
Vibrational harmonic frequencies (cm-1):
48.0406 53.7971 75.4763
105.9341 124.4629 141.9200
195.7721 211.7467 218.5406
344.4485 368.4928 411.6841
446.6154 472.3467 545.1325
558.7250 604.8006 643.6548
651.1163 722.6584 848.4529
945.5485 1220.0265 1235.1844
1262.6165 1347.5944 1368.8487
1414.4823 1899.8467 3769.7304
Zero-point correction (Hartree): 0.050707

CF2OCFCF2oxide

E(CCS(D)/Aug-CC-pVDZ) (Hartree): -762.34736815
E(CCS(D)/Aug-CC-pVDZ) (Hartree): -762.29155883
T1 diagnostic: 0.015476
E(MP2/Aug-CC-pVDZ) (Hartree): -762.27324443
E(MP3/Aug-CC-pVDZ) (Hartree): -762.26916913
E(PMP2/Aug-CC-pVDZ) (Hartree): -762.27414350
E(PMP3/Aug-CC-pVDZ) (Hartree): -762.26971430
E(PUHF/Aug-CC-pVDZ) (Hartree): -760.51813262
E(UHF/Aug-CC-pVDZ) (Hartree): -760.51652241
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -762.36705355
T1 diagnostic: 0.000342
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -763.01222617
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -762.95573571
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -762.95816991
T1 diagnostic: 0.013927
D1 diagnostic: 0.051838
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -762.90167945
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -763.01121960
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -762.95472914
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -763.01597650
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -762.95948604
E(ROHF/AUG-CC-PVDZ) (Hartree): -760.51420455
E(UM062X/Aug-CC-pVTZ) (Hartree): -763.93270461
Electronic state : 2-A
Cartesian coordinates (Angs):
C 1.665535 -0.235154 -0.001502
C 0.355397 0.375006 -0.062484
O -0.721466 -0.316469 -0.520953
F 0.211460 1.688877 -0.179776
C -1.865512 -0.170195 0.208155
F -2.557483 0.904881 -0.110769
F -2.602517 -1.245800 0.102795
F 1.890295 -1.485551 -0.349794
O 0.928535 -0.039089 1.148952
F 2.770572 0.473874 -0.116791
Rotational constants (GHz): 2.8170800 0.8481800 0.7059000
Vibrational harmonic frequencies (cm-1):
23.4862 54.7258 120.8768

189.3076	247.9181	291.8713
399.6678	464.2976	524.7385
557.1793	580.4822	591.7129
676.9716	754.6718	818.6343
889.2836	1073.3178	1168.9634
1223.0893	1243.9420	1310.3014
1330.9024	1335.4954	1659.4276

Zero-point correction (Hartree): 0.039939

CF3OCCF2oxide

E(CCS(T)/Aug-CC-pVDZ) (Hartree): -762.35692294
E(CCS(T)/Aug-CC-pVDZ) (Hartree): -762.30086983
T1 diagnostic: 0.016113
E(MP2/Aug-CC-pVDZ) (Hartree): -762.28211803
E(MP3/Aug-CC-pVDZ) (Hartree): -762.27828136
E(PMP2/Aug-CC-pVDZ) (Hartree): -762.28366849
E(PMP3/Aug-CC-pVDZ) (Hartree): -762.27925679
E(PUHF/Aug-CC-pVDZ) (Hartree): -760.52736845
E(UHF/Aug-CC-pVDZ) (Hartree): -760.52492002
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -762.37617071
T1 diagnostic: 0.000424
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -763.02262505
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -762.96612639
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -762.96833382
T1 diagnostic: 0.015178
D1 diagnostic: 0.070706
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -762.91183515
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -763.02160517
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -762.96510650
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -763.02651683
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -762.97001817
E(ROHF/AUG-CC-PVDZ) (Hartree): -760.52154358
E(UM062X/Aug-CC-pVTZ) (Hartree): -763.94181699
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-1.929430	0.011476	-0.068472
C	-0.548650	-0.339502	-0.197303
O	0.479595	0.522828	-0.370075
C	1.723486	0.031340	-0.061111
F	2.052733	-0.972269	-0.861393
F	2.581219	1.013707	-0.218257
F	-2.352541	1.257836	0.024335
O	-1.201452	-0.452380	1.023170
F	-2.909313	-0.758319	-0.495813
F	1.772616	-0.405785	1.188523

Rotational constants (GHz): 3.2260600 0.7128200 0.6888800
Vibrational harmonic frequencies (cm-1):

29.5396	59.8188	110.7674
205.4997	262.9486	360.7505
440.7598	500.5218	527.3636
542.8031	622.4714	631.4558
673.4118	733.0941	805.9305
921.0847	957.7983	1170.2819
1211.1854	1265.9956	1276.5725
1309.1788	1346.3577	1630.7166

Zero-point correction (Hartree): 0.040087

CF3OCCF2OH

E(CCS(T)/Aug-CC-pVDZ) (Hartree): -838.16777797
E(CCS(T)/Aug-CC-pVDZ) (Hartree): -838.10449444
T1 diagnostic: 0.014858
E(MP2/Aug-CC-pVDZ) (Hartree): -838.08658535
E(MP3/Aug-CC-pVDZ) (Hartree): -838.08037577
E(RHF/Aug-CC-pVDZ) (Hartree): -836.10817264
E(RM062X/Aug-CC-pVTZ) (Hartree): -839.90423913
Electronic state : 1-A
Cartesian coordinates (Angs):

C	-1.854856	-0.159767	-0.000011
C	-0.438818	0.455911	0.000000

O	0.500736	-0.509017	0.000034	
C	1.831717	-0.128909	0.000000	
F	2.125130	0.576524	1.075987	
F	2.125026	0.576573	-1.075978	
F	2.538438	-1.237262	-0.000079	
F	-1.977278	-0.955777	1.077806	
O	-2.796308	0.792654	-0.000322	
F	-1.977171	-0.956220	-1.077485	
H	-2.374405	1.663323	0.000029	
O	-0.249572	1.628704	0.000010	
Rotational constants (GHz):	2.2283300	0.6736300	0.6310900	
Vibrational harmonic frequencies (cm-1):				
32.8480	61.5489		110.0044	
134.2569	219.0461		226.7349	
343.4511	353.8895		390.5974	
443.1185	445.6198		535.5075	
566.9201	593.4068		630.0337	
678.5275	767.6590		794.5522	
906.4158	932.3305		1117.0631	
1186.4887	1205.3410		1275.5785	
1303.3657	1337.7853		1349.8449	
1463.0118	1915.2829		3788.5830	
Zero-point correction (Hartree):	0.057202			

CFOHCF2

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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -450.60736301
E(CCSD/Aug-CC-pVDZ) (Hartree): -450.57401707
  T1 diagnostic: 0.014209
E(MP2/Aug-CC-pVDZ) (Hartree): -450.55676005
E(MP3/Aug-CC-pVDZ) (Hartree): -450.56202051
E(RHF/Aug-CC-pVDZ) (Hartree): -449.49611634
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -451.00406231
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -450.96988970
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -450.97185401
  T1 diagnostic: 0.012657
  D1 diagnostic: 0.036188
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -450.93768139
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -451.00362064
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -450.96944803
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -451.00594509
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -450.97177247
E(RHF/AUG-CC-PVDZ) (Hartree): -449.49611633
E(RM062X/Aug-CC-pVTZ) (Hartree): -451.54768494
Electronic state : 1-A
Cartesian coordinates (Angs):
  C    0.656130    0.013701   -0.006502
  C   -0.661490   -0.030323    0.009164
  O    1.446122   -1.060012   -0.107734
  F    1.280076    1.192610    0.007123
  F   -1.429429    1.035357   -0.004390
  F   -1.353870   -1.146394    0.015487
  H    1.992182   -1.154327    0.681922
Rotational constants (GHz):    5.4919500    3.2940600    2.0698000
Vibrational harmonic frequencies (cm-1):
  194.0972    216.8521    264.6386
  400.8565    431.6542    564.5436
  571.4981    602.0626    820.2360
  1172.3342    1287.8268    1362.9885
  1380.5251    1948.8876    3814.8780
Zero-point correction (Hartree): 0.034250

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CFOCFOH

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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -426.02904005
E(CCSD/Aug-CC-pVDZ) (Hartree): -425.99420552
  T1 diagnostic: 0.020264
E(MP2/Aug-CC-pVDZ) (Hartree): -425.97529967
E(MP3/Aug-CC-pVDZ) (Hartree): -425.97757400
E(PMP2/Aug-CC-pVDZ) (Hartree): -425.98123547
E(PMP3/Aug-CC-pVDZ) (Hartree): -425.98126355

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E(PUHF/Aug-CC-pVDZ) (Hartree): -424.95311725
 E(UHF/Aug-CC-pVDZ) (Hartree): -424.94538756
 E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -426.03132137
 T1 diagnostic: 0.000396
 E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -426.40556281
 E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -426.37222904
 E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -426.37143085
 T1 diagnostic: 0.017109
 D1 diagnostic: 0.064312
 E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -426.33809709
 E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -426.40494249
 E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -426.37160872
 E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -426.40829172
 E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -426.37495795
 E(ROHF/AUG-CC-PVDZ) (Hartree): -424.93771413
 E(UM062X/Aug-CC-pVTZ) (Hartree): -426.93307043
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C 0.754553 0.128359 0.000367
 C -0.651790 -0.033475 0.000070
 F 1.430825 -1.023298 -0.000033
 O 1.293598 1.205334 -0.000041
 F -1.240602 -1.195519 -0.000078
 O -1.468606 0.994490 -0.000164
 H -0.928519 1.801454 0.000004
 Rotational constants (GHz): 5.6555000 3.5529700 2.1821000
 Vibrational harmonic frequencies (cm-1):
 157.5606 225.7205 328.2947
 408.4199 535.6237 557.3604
 612.8479 653.8507 841.1988
 1148.6338 1250.4243 1508.4537
 1610.2012 1790.5511 3755.1269
 Zero-point correction (Hartree): 0.035048

CFOCF2O

 E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -525.05836483
 E(CCSD/Aug-CC-pVDZ) (Hartree): -525.01929689
 T1 diagnostic: 0.017057
 E(MP2/Aug-CC-pVDZ) (Hartree): -524.99814725
 E(MP3/Aug-CC-pVDZ) (Hartree): -525.00159692
 E(PMP2/Aug-CC-pVDZ) (Hartree): -525.00012655
 E(PMP3/Aug-CC-pVDZ) (Hartree): -525.00281169
 E(PUHF/Aug-CC-pVDZ) (Hartree): -523.80160897
 E(UHF/Aug-CC-pVDZ) (Hartree): -523.79834834
 E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -525.05863601
 T1 diagnostic: 0.000285
 E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -525.51096034
 E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -525.47167219
 E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -525.47318757
 T1 diagnostic: 0.015868
 D1 diagnostic: 0.051619
 E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -525.43389942
 E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -525.51016783
 E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -525.47087968
 E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -525.51399971
 E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -525.47471156
 E(ROHF/AUG-CC-PVDZ) (Hartree): -523.79361420
 E(UM062X/Aug-CC-pVTZ) (Hartree): -526.14217031
 Point group : CS
 Electronic state : 2-A'
 Cartesian coordinates (Angs):
 C 0.318830 -0.884516 0.000000
 C -0.047256 0.625293 0.000000
 F -0.800310 -1.588483 0.000000
 O 1.393234 -1.348090 0.000000
 F -0.800310 0.896921 1.077851
 F -0.800310 0.896921 -1.077851
 O 1.104130 1.311478 0.000000
 Rotational constants (GHz): 3.9044600 2.5750200 2.1287000
 Vibrational harmonic frequencies (cm-1):

43.6510 (A")	223.1176 (A')	227.8454 (A")
376.3144 (A')	412.0765 (A')	471.7187 (A")
591.1972 (A')	685.9632 (A')	772.8299 (A")
824.5014 (A')	1064.9203 (A')	1188.8857 (A")
1229.0920 (A')	1309.9565 (A')	1979.7835 (A')

Zero-point correction (Hartree): 0.025975

OCCF2OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -426.01734478
E(CCSD/Aug-CC-pVDZ) (Hartree): -425.98232323
T1 diagnostic: 0.018175
E(MP2/Aug-CC-pVDZ) (Hartree): -425.96617013
E(MP3/Aug-CC-pVDZ) (Hartree): -425.96587666
E(PMP2/Aug-CC-pVDZ) (Hartree): -425.96865245
E(PMP3/Aug-CC-pVDZ) (Hartree): -425.96733430
E(PUHF/Aug-CC-pVDZ) (Hartree): -424.93616796
E(UHF/Aug-CC-pVDZ) (Hartree): -424.93238472
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -426.01665548
T1 diagnostic: 0.000562
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -426.39048335
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -426.35695254
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -426.35645981
T1 diagnostic: 0.017714
D1 diagnostic: 0.066500
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -426.32292900
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -426.38977793
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -426.35624712
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -426.39343365
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -426.35990284
E(ROHF/AUG-CC-PVDZ) (Hartree): -424.92717822
E(UM062X/Aug-CC-pVTZ) (Hartree): -426.91125170
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-1.049838	-0.584821	-0.365029
C	0.372313	0.004091	-0.021205
O	-2.039488	-0.048544	-0.059200
F	1.245081	-0.355275	-0.953054
F	0.729001	-0.593934	1.134673
O	0.379505	1.346198	0.078770
H	-0.421722	1.646023	0.526285

Rotational constants (GHz): 5.7013100 3.0374700 2.9896300
Vibrational harmonic frequencies (cm-1):

84.2093	234.1835	322.9303
377.1291	422.6526	537.6628
560.6715	667.3818	799.1284
1088.4371	1156.8923	1272.6315
1401.8200	2009.1179	3826.4575

Zero-point correction (Hartree): 0.033629

TS.CF3OCFCF2OH.CF3+HOFC2CFO

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.64053884
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.57751906
T1 diagnostic: 0.021698
E(MP2/Aug-CC-pVDZ) (Hartree): -862.54113787
E(MP3/Aug-CC-pVDZ) (Hartree): -862.54154177
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.55622720
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.55218426
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.58615424
E(UHF/Aug-CC-pVDZ) (Hartree): -860.56834725
E(ROHF/CC-PVDZ) (Hartree): -860.48484834
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.41078610
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.675179	-0.301957	-0.020319
C	0.601874	0.679554	-0.466645
O	-0.427502	0.409188	-1.057597
F	0.791182	1.893652	0.066706
C	-1.902316	-0.114266	0.024439
F	-1.567591	0.212967	1.247567

F	-2.066878	-1.401137	-0.102985	
F	-2.937743	0.554336	-0.395947	
F	1.712072	-1.327230	-0.865208	
O	2.913542	0.232442	0.052739	
F	1.288419	-0.795856	1.190007	
H	2.888084	1.056401	0.552753	
Rotational constants (GHz):	1.9154300	0.7076500	0.6410400	
Vibrational harmonic frequencies (cm-1):				
i554.3051	17.6969		29.9469	
60.0456	106.4047		159.4376	
193.3809	228.9851		253.3864	
303.7389	398.2146		441.5904	
509.4854	529.9397		542.0100	
588.3353	640.8728		684.9633	
720.1813	821.6012		1037.7657	
1087.9941	1114.6423		1230.4715	
1337.5576	1354.5967		1363.3952	
1415.4877	1566.2214		3853.5012	
Zero-point correction (Hartree):	0.051468			

TS.CF3OCFCF2OH.CF3OCFCFOH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.62089616
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.55628508
T1 diagnostic: 0.017062
E(MP2/Aug-CC-pVDZ) (Hartree): -862.53597953
E(MP3/Aug-CC-pVDZ) (Hartree): -862.53031313
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.53927138
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.53214927
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.54326571
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53815295
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.39472005
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.711898	-0.336597	-0.266718
C	0.594830	0.423750	-0.531622
O	-0.564049	-0.101718	-0.902386
F	0.693846	1.724209	-0.560208
C	-1.638371	-0.119125	-0.010078
F	-1.747004	1.022736	0.632788
F	-1.519375	-1.107347	0.844149
F	-2.710604	-0.313017	-0.750733
F	1.645994	-1.620026	-0.414891
O	2.891437	0.135433	0.017250
F	0.813663	0.192743	1.497945
H	2.782080	0.828425	0.690152
Rotational constants (GHz):	1.9562300	0.7969100	0.7328400

Vibrational harmonic frequencies (cm-1):

i288.5897	39.9094	71.1590
122.4052	141.5482	198.9146
220.6904	293.9334	310.8448
352.1449	374.8366	453.0971
468.9650	534.7445	558.8619
576.5838	635.7790	656.6258
719.1082	845.1584	949.7707
1175.1196	1178.4222	1277.9653
1322.9519	1331.8935	1430.5431
1516.0337	1744.2349	3734.9404
Zero-point correction (Hartree):	0.052938	

TS.CF3OCFCF2OH.CF3OCFCF2+OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.59765570
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.53604781
T1 diagnostic: 0.022139
E(MP2/Aug-CC-pVDZ) (Hartree): -862.49515544
E(MP3/Aug-CC-pVDZ) (Hartree): -862.50566040
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.50376341
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.51202356
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.54923808
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53923177

E(UM062X/Aug-CC-pVTZ) (Hartree): -864.36729051
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	-1.430017	-0.389354	-0.342970
C	-0.493943	0.520273	-0.057218
O	0.577187	0.264575	0.716942
F	-0.639776	1.777369	-0.433494
C	1.730628	-0.086846	0.056047
F	2.087416	0.843581	-0.819237
F	1.576661	-1.227432	-0.602013
F	2.670646	-0.225187	0.959978
F	-1.320795	-1.642314	-0.031017
O	-2.256083	0.235902	1.631870
F	-2.462013	-0.123902	-1.090793
H	-2.618092	1.112699	1.423514

Rotational constants (GHz): 1.8013900 0.7260600 0.6812700
 Vibrational harmonic frequencies (cm-1):

i269.7903	35.5640	75.2433
80.9655	96.5386	143.5073
193.0727	205.3903	234.8082
345.5931	368.7527	416.3289
467.0634	514.0674	558.6978
559.9480	630.7137	644.3130
652.2667	720.9117	843.0935
946.2057	1223.8594	1231.1164
1269.6484	1348.6497	1377.9124
1431.2420	1803.7253	3772.7118

Zero-point correction (Hartree): 0.050557

TS.CF3OCFCF2OH.CF2O+CF2O+CF+HF

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.58433156
 E(CCSD/Aug-CC-pVDZ) (Hartree): -862.51846768
 T1 diagnostic: 0.018586
 E(MP2/Aug-CC-pVDZ) (Hartree): -862.49719885
 E(MP3/Aug-CC-pVDZ) (Hartree): -862.48651034
 E(PMP2/Aug-CC-pVDZ) (Hartree): -862.50060635
 E(PMP3/Aug-CC-pVDZ) (Hartree): -862.48867594
 E(PUHF/Aug-CC-pVDZ) (Hartree): -860.51216120
 E(UHF/Aug-CC-pVDZ) (Hartree): -860.50737212
 E(UM062X/Aug-CC-pVTZ) (Hartree): -864.34864816
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	-1.679507	-0.400523	-0.016743
C	-1.046212	1.204728	0.536700
O	0.889177	0.463150	1.010832
F	-0.762172	1.959309	-0.416475
C	1.814145	0.335115	0.278699
F	1.857814	0.773474	-0.938003
F	1.235130	-1.818077	-0.644830
F	2.950707	-0.161798	0.619384
F	-1.642567	-1.104947	1.114722
O	-1.018363	-0.738867	-0.988259
F	-2.953887	0.008852	-0.196581
H	0.337709	-1.541497	-0.816473

Rotational constants (GHz): 1.6701500 0.7216100 0.6239500
 Vibrational harmonic frequencies (cm-1):

i115.4814	32.1340	54.5247
88.8987	101.8802	131.4933
158.8324	171.8931	187.1125
222.4105	234.4065	290.6940
383.1021	418.6631	563.0746
597.7023	620.3676	657.7788
734.7096	777.5463	830.7555
877.2384	952.7793	1043.5024
1190.9469	1384.0759	1427.6633
1640.0739	1905.8151	3367.2170

Zero-point correction (Hartree): 0.047949

TS.CF3OCFCF2OH.CF3OCFCFO+HF

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.62002013
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.55533439
T1 diagnostic: 0.017866
E(MP2/Aug-CC-pVDZ) (Hartree): -862.53576873
E(MP3/Aug-CC-pVDZ) (Hartree): -862.52686098
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.53782573
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.52814138
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.54313662
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53989298
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.39324575
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.556082	-0.175304	-0.376269
C	0.479334	0.755273	-0.514002
O	-0.745688	0.430512	-0.912185
F	0.642071	1.959233	-0.057421
C	-1.619128	-0.119969	0.026731
F	-1.621320	0.598691	1.132480
F	-1.272191	-1.355312	0.315304
F	-2.810145	-0.108620	-0.521387
F	1.339248	-1.318588	-0.962025
O	2.745198	0.171731	-0.081777
F	1.387507	-0.602490	1.444578
H	2.519656	-0.134170	0.969177

Rotational constants (GHz): 1.8474200 0.8202600 0.7500100
Vibrational harmonic frequencies (cm-1):

i1374.2634	38.9989	61.4906
83.4007	110.1025	157.3347
211.8273	278.0787	349.3974
370.1721	451.4480	463.6846
524.6076	554.5607	579.1052
632.8354	656.9980	699.8217
796.6602	842.8946	913.8695
943.2971	1170.9947	1241.5041
1309.4966	1350.2878	1432.0946
1492.3127	1648.5753	2223.0076

Zero-point correction (Hartree): 0.049183

TS.CF3OCFCF2OH.CF3OCFHCF2O

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.61672615
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.55411607
T1 diagnostic: 0.019024
E(MP2/Aug-CC-pVDZ) (Hartree): -862.52060322
E(MP3/Aug-CC-pVDZ) (Hartree): -862.52494024
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.52743790
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.52909278
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.55284052
E(UHF/Aug-CC-pVDZ) (Hartree): -860.54339114
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.38698276
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-1.804584	-0.246054	0.005822
C	-0.462313	0.445265	0.028629
O	0.545804	-0.210806	-0.603969
F	-2.827331	0.537295	-0.340948
F	-0.468615	1.757231	-0.168416
F	-1.878282	-1.356963	-0.726917
O	-1.746009	-0.494039	1.353106
H	-0.586774	0.133735	1.306255
C	1.796478	-0.137223	-0.038498
F	1.798170	-0.682814	1.170723
F	2.207568	1.115582	0.077505
F	2.614148	-0.799987	-0.820287

Rotational constants (GHz): 2.1031100 0.6925800 0.6373700
Vibrational harmonic frequencies (cm-1):

i1996.4691	22.5775	63.9986
86.7921	162.5658	208.0786
276.0575	332.5992	365.4500
429.7050	481.7408	531.8349
551.9575	591.4560	638.1031

657.9074	728.9378	758.0103
899.1180	935.4871	1089.4735
1173.2517	1189.6807	1222.2583
1258.2395	1277.8000	1317.2779
1353.5748	1463.1174	2014.8314

Zero-point correction (Hartree): 0.050306

TS.CF3OCFCFO.CFOCFO+CF3

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -762.37291173
E(CCSD/Aug-CC-pVDZ) (Hartree): -762.31353776
T1 diagnostic: 0.024691
E(MP2/Aug-CC-pVDZ) (Hartree): -762.28306046
E(MP3/Aug-CC-pVDZ) (Hartree): -762.27742277
E(PMP2/Aug-CC-pVDZ) (Hartree): -762.29903778
E(PMP3/Aug-CC-pVDZ) (Hartree): -762.28901160
E(PUHF/Aug-CC-pVDZ) (Hartree): -760.55094787
E(UHF/Aug-CC-pVDZ) (Hartree): -760.53247703
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -762.39039281
T1 diagnostic: 0.000789
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -763.03266543
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -762.97609684
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -762.97438622
T1 diagnostic: 0.021568
D1 diagnostic: 0.119084
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -762.91781763
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -763.03175773
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -762.97518913
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -763.03792702
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -762.98135842
E(ROHF/AUG-CC-PVDZ) (Hartree): -760.51923147
E(UM062X/Aug-CC-pVTZ) (Hartree): -763.94798878
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.882104	-0.383208	0.020853
C	0.890233	0.639866	-0.434929
O	-0.125158	0.408804	-1.054218
F	1.189742	1.856701	-0.021705
C	-1.681629	-0.085399	0.087392
F	-2.544698	0.887351	0.128964
F	-1.113469	-0.290508	1.249035
F	-2.180694	-1.173108	-0.417159
F	1.454279	-1.602365	-0.291486
O	2.901323	-0.175078	0.570625

Rotational constants (GHz): 2.2250000 0.8638900 0.7306400
Vibrational harmonic frequencies (cm-1):

i469.8749	29.9762	40.4820
60.0664	98.7742	147.1805
218.5968	269.6410	432.5613
460.9035	527.0947	531.1295
576.9237	672.4402	678.7719
782.1003	844.4182	1030.0366
1169.8016	1316.6842	1353.2941
1375.1597	1591.0766	1938.2571

Zero-point correction (Hartree): 0.036782

TS.CF3OCFCF2OH.CF3+HOFC2CFO.b

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.64143690
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.57833043
T1 diagnostic: 0.021296
E(MP2/Aug-CC-pVDZ) (Hartree): -862.54305868
E(MP3/Aug-CC-pVDZ) (Hartree): -862.54315242
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.55766115
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.55337565
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.58559720
E(UHF/Aug-CC-pVDZ) (Hartree): -860.56828990
E(ROHF/AUG-CC-PVDZ) (Hartree): -860.53719978
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.41264558
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.643618	-0.317813	0.007686	
C	0.592362	0.686344	-0.435467	
O	-0.428319	0.391852	-1.050652	
F	0.786050	1.907760	0.024192	
C	-1.895640	-0.104331	0.032360	
F	-1.403107	-0.125308	1.243888	
F	-2.269831	-1.290869	-0.362798	
F	-2.847050	0.774481	-0.096808	
F	2.827480	0.037615	-0.545708	
O	1.306590	-1.581765	-0.315238	
F	1.818142	-0.245982	1.335655	
H	0.726653	-1.585167	-1.086147	
Rotational constants (GHz):	1.9442600	0.7014200	0.6381900	
Vibrational harmonic frequencies (cm-1):				
i554.9889	30.7411		40.2811	
52.2054	102.6362		170.8503	
203.8192	224.4021		252.8621	
287.5366	416.7374		441.2303	
510.9703	530.2345		543.7918	
603.1876	643.0244		686.5523	
716.5144	825.8487		1036.4248	
1097.3670	1150.3383		1251.7802	
1339.0225	1351.7626		1361.6419	
1430.9888	1552.9938		3834.9776	
Zero-point correction (Hartree):	0.051693			

TS.CF3OCFOCF2OH.CF2OH+CF3OCFO

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -937.74202802
E(CCSD/Aug-CC-pVDZ) (Hartree): -937.67388820
T1 diagnostic: 0.021478
E(MP2/Aug-CC-pVDZ) (Hartree): -937.63715236
E(MP3/Aug-CC-pVDZ) (Hartree): -937.64267252
E(PMP2/Aug-CC-pVDZ) (Hartree): -937.64083195
E(PMP3/Aug-CC-pVDZ) (Hartree): -937.64508852
E(PUHF/Aug-CC-pVDZ) (Hartree): -935.48985728
E(UHF/Aug-CC-pVDZ) (Hartree): -935.48447937
E(UM062X/Aug-CC-pVTZ) (Hartree): -939.67113747
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.423577	-0.537545	-0.032357	
C	0.524255	0.910984	0.047453	
O	-0.730656	0.759311	-0.543005	
F	1.174813	1.719825	-0.809178	
C	-1.695078	-0.065255	-0.040411	
F	-1.726008	-0.092379	1.281715	
F	-1.536173	-1.309415	-0.477289	
F	-2.852315	0.385085	-0.488265	
F	1.293460	-0.958347	-1.272796	
O	2.707097	-0.309166	0.234017	
F	0.848204	-1.409986	0.792731	
H	2.794963	0.114579	1.098074	
O	0.632395	1.177764	1.285430	
Rotational constants (GHz):	1.5704300	0.7679200	0.6989100	
Vibrational harmonic frequencies (cm-1):				
i332.3175	43.4372		75.3512	
118.8831	167.9631		196.3794	
227.2453	275.4721		332.6196	
340.9035	361.0959		389.5071	
429.9941	461.1038		544.2182	
559.9386	579.3278		627.3729	
675.8678	732.9627		741.3972	
863.7859	970.7471		1053.9986	
1159.9015	1187.5700		1263.0971	
1277.1348	1308.8452		1316.7117	
1391.0107	1438.9925		3823.7297	
Zero-point correction (Hartree):	0.056810			

TS.CF3OCFOCF2OH.CF3O+HOFCF2CFO

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -937.71943602

E (CCSD/Aug-CC-pVDZ) (Hartree): -937.64922839
 T1 diagnostic: 0.021029
 E (MP2/Aug-CC-pVDZ) (Hartree): -937.60911516
 E (MP3/Aug-CC-pVDZ) (Hartree): -937.61365277
 E (PMP2/Aug-CC-pVDZ) (Hartree): -937.62395586
 E (PMP3/Aug-CC-pVDZ) (Hartree): -937.62379711
 E (PUHF/Aug-CC-pVDZ) (Hartree): -935.47239880
 E (UHF/Aug-CC-pVDZ) (Hartree): -935.45430877
 E (UM062X/Aug-CC-pVTZ) (Hartree): -939.63900777
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	1.466486	-0.510372	-0.011660
C	0.866124	0.911549	0.115324
O	-0.831182	0.791376	-0.533606
F	1.337547	1.693055	-0.836794
C	-1.778961	-0.080304	-0.047648
F	-1.817105	-0.089910	1.268798
F	-1.570140	-1.315568	-0.481429
F	-2.944991	0.328536	-0.525151
F	1.323347	-0.953977	-1.252681
O	2.790918	-0.488665	0.280771
F	0.799006	-1.318774	0.826258
H	2.912051	-0.421763	1.235664
O	0.492400	1.373073	1.182489

Rotational constants (GHz): 1.5631000 0.7213300 0.6548900
 Vibrational harmonic frequencies (cm-1):

i628.7103	40.6591	62.2319
115.2759	129.0157	161.0747
216.4083	244.9752	250.9064
294.7989	379.6499	386.5131
431.6183	450.4395	496.0167
565.1062	591.5541	613.9707
652.3146	675.3761	698.8486
826.7784	929.1785	1123.6266
1135.5794	1195.3604	1253.2480
1283.2263	1297.8371	1341.1639
1401.9115	1574.4071	3844.1724

Zero-point correction (Hartree): 0.056187

TS.CF3OCFOCF2OH.F+HOCHF2COOCH3

 E (CCSD(T)/Aug-CC-pVDZ) (Hartree): -937.71185877
 E (CCSD/Aug-CC-pVDZ) (Hartree): -937.64208949
 T1 diagnostic: 0.021073
 E (MP2/Aug-CC-pVDZ) (Hartree): -937.60199907
 E (MP3/Aug-CC-pVDZ) (Hartree): -937.60717835
 E (PMP2/Aug-CC-pVDZ) (Hartree): -937.61411966
 E (PMP3/Aug-CC-pVDZ) (Hartree): -937.61552018
 E (PUHF/Aug-CC-pVDZ) (Hartree): -935.46831054
 E (UHF/Aug-CC-pVDZ) (Hartree): -935.45360121
 E (UM062X/Aug-CC-pVTZ) (Hartree): -939.62871153
 Electronic state : 2-A
 Cartesian coordinates (Angs):

C	1.797299	-0.357837	-0.024446
C	0.394143	0.223616	0.263938
O	-0.559790	-0.484117	-0.334062
F	0.557327	1.791532	-0.747198
C	-1.891841	-0.204733	-0.041423
F	-2.187932	1.047922	-0.310492
F	-2.151156	-0.455370	1.227252
F	-2.607134	-1.002080	-0.800031
F	1.941340	-0.596897	-1.327946
O	2.776898	0.436665	0.421449
F	1.821497	-1.560892	0.596334
H	2.414389	1.132145	0.986166
O	0.210707	1.032907	1.173131

Rotational constants (GHz): 1.6432800 0.6407200 0.5926400
 Vibrational harmonic frequencies (cm-1):

i580.4901	51.2225	64.6886
87.4923	128.4582	153.5557
194.8202	228.0485	234.7415

321.3706	343.3728	377.8342
435.4696	452.3041	522.0948
566.7010	594.7533	617.1754
657.0881	678.1758	768.1261
898.6162	930.4179	1108.0816
1177.1383	1192.6682	1283.9011
1310.2182	1344.4806	1367.1172
1453.0168	1608.3332	3809.0531

Zero-point correction (Hartree): 0.056864

TS.CF3OCFOCFO.CF3OCFO+CFO

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -837.46524178
E(CCSD/Aug-CC-pVDZ) (Hartree): -837.39834265
T1 diagnostic: 0.023844
E(MP2/Aug-CC-pVDZ) (Hartree): -837.37162475
E(MP3/Aug-CC-pVDZ) (Hartree): -837.36400140
E(PMP2/Aug-CC-pVDZ) (Hartree): -837.38168932
E(PMP3/Aug-CC-pVDZ) (Hartree): -837.37073540
E(PUHF/Aug-CC-pVDZ) (Hartree): -835.43430294
E(UHF/Aug-CC-pVDZ) (Hartree): -835.42138532
E(UM062X/Aug-CC-pVTZ) (Hartree): -839.19865386

Electronic state : 2-A

Cartesian coordinates (Angs):

C	-1.879207	0.546995	-0.115421
C	-0.591487	-0.779432	0.150928
O	0.431770	-0.382353	-0.695201
F	-1.191844	-1.801725	-0.454384
C	1.574872	0.123791	-0.124111
F	2.132703	-0.751298	0.695389
F	1.347212	1.242918	0.543153
F	2.404375	0.378072	-1.115463
F	-1.357463	1.699910	0.174938
O	-2.963318	0.289452	-0.434282
O	-0.548442	-0.689476	1.371849

Rotational constants (GHz): 1.8991800 0.8400000 0.7230500

Vibrational harmonic frequencies (cm-1):

i370.8262	42.7355	64.9691
89.2924	159.4585	186.9478
208.2233	297.4666	376.8775
390.2583	418.8619	481.9112
555.8845	607.3861	631.3872
679.9620	696.7828	747.4453
889.4183	1007.4192	1136.9004
1203.0303	1263.4581	1287.5968
1329.2013	1571.1631	2022.8479

Zero-point correction (Hartree): 0.041797

TS.CF3OCFOCFO.CF3OCOCFO+F

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -837.43219275
E(CCSD/Aug-CC-pVDZ) (Hartree): -837.36003044
T1 diagnostic: 0.021182
E(MP2/Aug-CC-pVDZ) (Hartree): -837.35010221
E(MP3/Aug-CC-pVDZ) (Hartree): -837.32777222
E(PMP2/Aug-CC-pVDZ) (Hartree): -837.35888663
E(PMP3/Aug-CC-pVDZ) (Hartree): -837.33323747
E(PUHF/Aug-CC-pVDZ) (Hartree): -835.37222406
E(UHF/Aug-CC-pVDZ) (Hartree): -835.36068024
E(UM062X/Aug-CC-pVTZ) (Hartree): -839.16044089

Electronic state : 2-A

Cartesian coordinates (Angs):

C	2.065619	0.204204	-0.230711
C	0.486934	-0.505151	-0.000707
O	-0.403153	0.500511	0.059028
F	1.460745	-0.157452	1.379908
C	-1.743650	0.146473	-0.051731
F	-2.081639	-0.721503	0.880000
F	-2.001338	-0.367288	-1.238136
F	-2.426088	1.256074	0.105014
F	1.944933	1.501271	-0.219434

O	2.954348	-0.441787	-0.619895	
O	0.333439	-1.642858	-0.247542	
Rotational constants (GHz):	2.2089000	0.7157400	0.6676600	
Vibrational harmonic frequencies (cm-1):				
i441.4778	36.9800		71.6160	
129.2402	159.6948		198.1710	
256.6988	322.1329		377.8939	
396.1222	430.4005		437.5991	
557.0824	590.6606		622.6895	
660.1061	678.0296		706.8826	
838.5195	907.3406		1110.8786	
1159.1472	1211.7602		1304.6640	
1348.7953	1853.4532		1948.1994	
Zero-point correction (Hartree):	0.041724			

TS.CF3OCFOCFO.FCOCFO+CF3O

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -837.44812455
E(CCSD/Aug-CC-pVDZ) (Hartree): -837.38134344
T1 diagnostic: 0.022634
E(MP2/Aug-CC-pVDZ) (Hartree): -837.34757892
E(MP3/Aug-CC-pVDZ) (Hartree): -837.34565704
E(PMP2/Aug-CC-pVDZ) (Hartree): -837.36389217
E(PMP3/Aug-CC-pVDZ) (Hartree): -837.35697153
E(PUHF/Aug-CC-pVDZ) (Hartree): -835.43202413
E(UHF/Aug-CC-pVDZ) (Hartree): -835.41233663
E(UM062X/Aug-CC-pVTZ) (Hartree): -839.17012119
Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.615557	-0.611459	0.167332
C	1.010145	0.802606	0.143083
O	-0.557765	0.531430	-0.734433
F	1.596647	1.548703	-0.773627
C	-1.563351	-0.131506	-0.075447
F	-2.257598	0.657755	0.717746
F	-1.095564	-1.145694	0.641700
F	-2.353915	-0.600421	-1.027964
F	1.822958	-1.054577	-1.060247
O	1.879264	-1.206063	1.139665
O	0.455144	1.298416	1.108733

Rotational constants (GHz):	1.7192600	0.8504700	0.7798900
Vibrational harmonic frequencies (cm-1):			
i648.6765	23.7542		53.6856
92.7687	132.3501		160.0546
231.5583	264.8513		335.0105
408.8784	432.9679		469.9908
531.8542	617.6893		632.7302
665.5211	682.2925		764.5716
829.1010	931.1272		1156.2570
1209.1234	1274.5178		1295.0814
1323.6037	1581.3031		1988.3552
Zero-point correction (Hartree):	0.041210		

TS.CF3OCFOHCF2.CF3OCFCF2+OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.59673149
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.53431566
T1 diagnostic: 0.023363
E(MP2/Aug-CC-pVDZ) (Hartree): -862.49349129
E(MP3/Aug-CC-pVDZ) (Hartree): -862.50288816
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.50215836
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.50924109
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.54255594
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53245923
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.36683411
Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.436578	0.525811	-0.240306
C	0.563514	-0.483445	-0.106182
O	-0.547336	-0.386869	0.658071
F	0.685200	-1.578384	-0.812154

C	-1.682096	0.076497	0.032129	
F	-1.998175	-0.670863	-1.013818	
F	-1.518767	1.323532	-0.390273	
F	-2.652741	0.037621	0.913973	
F	1.315673	1.656418	0.390785	
F	2.496295	0.450046	-0.979172	
O	1.994672	-1.004276	1.432824	
H	1.565972	-0.549355	2.174937	
Rotational constants (GHz):	1.7590800	0.7611000	0.7045100	
Vibrational harmonic frequencies (cm-1):				
i208.6375	34.5023		79.1320	
105.1782	118.6201		143.5370	
202.4421	213.5280		277.9692	
344.0591	363.3976		422.4983	
467.5007	512.1654		545.7766	
562.7722	641.5639		651.2857	
705.1343	713.5189		841.3200	
946.9926	1221.2347		1237.4507	
1273.2712	1348.5922		1389.0336	
1438.6167	1803.3672		3784.1755	
Zero-point correction (Hartree):	0.051005			

TS.CF3OCFOHCF2.CF3OCFOCF2H

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.60971245
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.54705959
T1 diagnostic: 0.019146
E(MP2/Aug-CC-pVDZ) (Hartree): -862.51333376
E(MP3/Aug-CC-pVDZ) (Hartree): -862.51764804
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.52035324
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.52192476
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.54543759
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53576565
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.37996583
Electronic state : 2-A
Cartesian coordinates (Angs):

C	0.584079	-0.319081	0.166980	
C	1.971163	0.269459	0.084270	
F	2.860488	-0.390354	-0.626697	
O	-0.405099	0.596571	-0.142219	
F	2.079851	1.569393	-0.082849	
F	0.414327	-1.443539	-0.549266	
O	0.749695	-0.585351	1.512433	
H	1.955479	-0.084164	1.358693	
C	-1.706731	0.182334	-0.087925	
F	-2.457953	1.254972	0.060719	
F	-1.920645	-0.647007	0.925947	
F	-2.065325	-0.432560	-1.205672	
Rotational constants (GHz):	2.0980500	0.7174500	0.6717500	
Vibrational harmonic frequencies (cm-1):				
i2015.9899	42.6400		69.2004	
99.6316	121.9042		207.3065	
270.9150	346.1038		362.3987	
400.5431	453.6452		524.0217	
536.6286	613.4970		641.1175	
669.4857	713.1116		781.4713	
861.0625	951.0369		1083.7002	
1143.3029	1179.7979		1211.4550	
1272.3321	1292.3885		1330.0886	
1331.9661	1480.4397		2022.9537	
Zero-point correction (Hartree):	0.050152			

TS.CF3OCFOHCF2.CF3OCOCF2+HF

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.61401792
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.54899858
T1 diagnostic: 0.018543
E(MP2/Aug-CC-pVDZ) (Hartree): -862.52915280
E(MP3/Aug-CC-pVDZ) (Hartree): -862.52017158
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.53110680
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.52140711

E(PUHF/Aug-CC-pVDZ) (Hartree): -860.53479003
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53171880
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.38724771
Electronic state : 2-A
Cartesian coordinates (Angs):

C	-1.312258	0.760795	-0.164717
C	-0.758554	-0.511710	-0.477459
O	0.549150	-0.672335	-0.702377
F	-1.070041	-1.392942	1.206577
C	1.554121	0.000751	-0.031792
F	1.243507	0.226919	1.226136
F	1.816906	1.153574	-0.630045
F	2.620940	-0.761173	-0.097674
F	-2.582001	0.885309	-0.003054
F	-0.630389	1.793135	0.206504
O	-1.533504	-1.426639	-0.946334
H	-1.615323	-1.850619	0.057490

Rotational constants (GHz): 1.7230700 0.8863000 0.7393900
Vibrational harmonic frequencies (cm-1):

i1223.2278	37.9054	86.3594
112.7170	126.5213	163.5551
210.2271	250.8651	360.5204
375.2722	393.6866	459.4113
504.3533	510.2285	566.2184
627.0775	648.9168	723.6016
788.0626	836.9924	911.0038
986.5642	1201.7558	1260.3519
1293.3931	1350.2492	1465.4356
1474.7543	1661.4034	2314.2668

Zero-point correction (Hartree): 0.049440

TS.CF3OCFOHCF2.CF3OCOHC3

E(CCS(T)/Aug-CC-pVDZ) (Hartree): -862.61664555
E(CCS(D)/Aug-CC-pVDZ) (Hartree): -862.55182261
T1 diagnostic: 0.017516
E(MP2/Aug-CC-pVDZ) (Hartree): -862.53166476
E(MP3/Aug-CC-pVDZ) (Hartree): -862.52541472
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.53497081
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.52736079
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.53732842
E(UHF/Aug-CC-pVDZ) (Hartree): -860.53233584
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.39021090
Electronic state : 2-A
Cartesian coordinates (Angs):

C	1.248349	-0.568850	-0.260649
C	0.673860	0.686173	-0.264747
O	-0.614678	0.975595	-0.075765
C	-1.613908	0.031073	-0.018360
F	-1.563172	-0.666071	1.097302
F	-1.547144	-0.797425	-1.052973
F	-2.747865	0.691339	-0.067818
F	2.473231	-0.728017	-0.638377
F	0.670858	-1.658911	0.108418
O	1.380495	1.739705	-0.610560
H	2.125039	1.787772	0.014780
F	1.591715	0.447913	1.524377

Rotational constants (GHz): 1.8324500 0.8303900 0.7467900
Vibrational harmonic frequencies (cm-1):

i325.1781	33.4947	89.7242
146.9702	152.1645	192.9073
235.0765	268.2588	320.7191
366.9892	385.8086	433.2069
463.7913	508.6374	547.4901
575.3980	628.2495	652.0262
733.9534	813.0123	974.4926
1206.6632	1230.8965	1274.7873
1289.2789	1345.0598	1459.0618
1477.1835	1737.7700	3705.4499

Zero-point correction (Hartree): 0.052964

TS.CFOCF2OH.CFOCF2OH+HF

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -525.67501740
E(CCSD/Aug-CC-pVDZ) (Hartree): -525.62938469
T1 diagnostic: 0.017762
E(MP2/Aug-CC-pVDZ) (Hartree): -525.62524209
E(MP3/Aug-CC-pVDZ) (Hartree): -525.60970230
E(RHF/Aug-CC-pVDZ) (Hartree): -524.34365141
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -526.13420652
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -526.09368003
E(CCSD-F12a/AUG-CC-PVDZ) (Hartree): -526.09008279
T1 diagnostic: 0.015939
D1 diagnostic: 0.062387
E(CCSD-F12b/AUG-CC-PVDZ) (Hartree): -526.04955631
E(CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -526.13317517
E(CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -526.09264868
E(CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -526.13823069
E(CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -526.09770420
E(RHF/AUG-CC-PVDZ) (Hartree): -524.34365140
E(RM062X/Aug-CC-pVTZ) (Hartree): -526.76180267
Electronic state : 1-A
Cartesian coordinates (Angs):
C -0.521730 0.218288 -0.256601
C 0.990286 0.073453 -0.038949
O 1.727696 0.951575 0.200869
F 1.350120 -1.185959 -0.187213
F -0.913873 1.432775 -0.028416
O -1.186420 -0.532654 -1.012257
F -1.067457 -0.695876 1.127884
H -1.460660 -1.060285 0.054098
Rotational constants (GHz): 3.6577100 2.4959500 1.9794400
Vibrational harmonic frequencies (cm-1):
i1718.0616 59.8729 169.9897
239.4132 315.8077 407.4263
491.5837 637.4505 659.1799
723.6792 823.8337 873.0426
917.9924 1210.9887 1386.1750
1642.7519 1987.6056 2033.5102
Zero-point correction (Hartree): 0.033216

TS.CF3+OCFCF2OH.CF4+OCFCFOH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.60928919
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.54248391
T1 diagnostic: 0.021610
E(MP2/Aug-CC-pVDZ) (Hartree): -862.51465178
E(MP3/Aug-CC-pVDZ) (Hartree): -862.50636818
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.52848288
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.51602616
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.53522157
E(UHF/Aug-CC-pVDZ) (Hartree): -860.51870478
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.37462523
Electronic state : 2-A
Cartesian coordinates (Angs):
C -1.320575 0.685812 -0.015120
C -1.977503 -0.634834 -0.163911
O -2.272130 -1.192534 -1.155896
F -2.130870 -1.185508 1.060254
C 2.105478 -0.141260 -0.003165
F 2.518447 0.051810 1.220336
F 2.618760 0.725365 -0.829737
F 2.309447 -1.368058 -0.394799
F -1.340311 1.431340 -1.083131
O -1.470709 1.396218 1.088526
F 0.318605 0.134282 0.002727
H -1.548391 0.809139 1.851308
Rotational constants (GHz): 1.6359200 0.5969900 0.5664500
Vibrational harmonic frequencies (cm-1):
i1064.4289 10.2986 47.6263
51.4001 98.9299 133.4582
187.5836 199.2136 236.2362

300.6501	330.2411	388.8668
418.1217	513.5960	534.6885
536.8198	552.7081	648.2857
714.4174	761.1712	832.7928
1043.2207	1101.4942	1207.7526
1351.6877	1358.7800	1424.5491
1479.7072	1935.0598	3835.1276

Zero-point correction (Hartree): 0.050654

TS.CF3+OCFCF2OH.HCF3+OCFCF2O

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.63813342
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.57501365
T1 diagnostic: 0.020325
E(MP2/Aug-CC-pVDZ) (Hartree): -862.54160471
E(MP3/Aug-CC-pVDZ) (Hartree): -862.54336663
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.54575705
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.54595130
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.57613376
E(UHF/Aug-CC-pVDZ) (Hartree): -860.56986068
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.40226723

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.287707	0.902589	-0.109434
C	1.228456	-0.641211	-0.133815
F	2.407610	1.306273	0.464526
O	0.465996	1.631851	-0.522222
F	1.156978	-1.056976	1.145321
F	2.363450	-1.118368	-0.660409
O	0.219075	-1.130316	-0.888205
H	-0.941948	-0.458569	-0.524408
C	-1.922161	-0.033622	-0.011129
F	-2.767482	-1.037070	0.128362
F	-1.610751	0.458920	1.174133
F	-2.450098	0.900528	-0.770368

Rotational constants (GHz): 1.7554700 0.6916500 0.6065400

Vibrational harmonic frequencies (cm-1):

i944.6941	33.9637	45.9322
52.0862	91.1267	108.0265
160.6220	235.8734	245.2956
345.5741	421.3763	438.1111
518.0687	526.0069	609.1327
646.6749	709.9752	783.4166
806.4170	870.0984	1128.1046
1142.4991	1192.5069	1232.1537
1265.9506	1289.3734	1332.7500
1358.0416	1422.5710	1979.9759

Zero-point correction (Hartree): 0.047823

TS.CFOCF2OH+CF3.CF4+OCCF2OH

E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -862.59903132
E(CCSD/Aug-CC-pVDZ) (Hartree): -862.53181085
T1 diagnostic: 0.021581
E(MP2/Aug-CC-pVDZ) (Hartree): -862.50284981
E(MP3/Aug-CC-pVDZ) (Hartree): -862.49495069
E(PMP2/Aug-CC-pVDZ) (Hartree): -862.51698383
E(PMP3/Aug-CC-pVDZ) (Hartree): -862.50465381
E(PUHF/Aug-CC-pVDZ) (Hartree): -860.52186956
E(UHF/Aug-CC-pVDZ) (Hartree): -860.50486614
E(UM062X/Aug-CC-pVTZ) (Hartree): -864.36185024

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.073557	0.938723	-0.302073
C	1.893291	-0.339558	-0.013227
O	1.356232	2.029302	-0.004132
F	1.690322	-1.243401	-0.966721
F	1.403048	-0.850408	1.138355
O	3.212563	-0.084822	0.068536
H	3.356729	0.777952	0.477902
F	-0.533755	0.472006	-0.409471

C	-2.163997	-0.066748	0.016449	
F	-2.667357	-0.629599	-1.044343	
F	-2.026939	-0.910107	0.998941	
F	-2.834675	0.991698	0.372124	
Rotational constants (GHz):	1.9659000	0.5803500	0.5323400	
Vibrational harmonic frequencies (cm-1):				
i1098.4268	12.9482		34.1881	
43.2479	70.8067		141.9630	
172.2737	215.5960		241.4977	
307.6501	318.4633		360.9121	
445.1286	500.1687		538.5659	
538.9492	581.7717		618.6026	
690.1078	750.0476		880.8564	
1049.3430	1125.4035		1169.2282	
1294.2123	1363.7877		1365.5017	
1394.2881	1985.0453		3821.2553	
Zero-point correction (Hartree):	0.050192			

TS.CFOCF2O.CFO+CF2O

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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -525.04961433
E(CCSD/Aug-CC-pVDZ) (Hartree): -525.00575542
  T1 diagnostic: 0.027180
E(MP2/Aug-CC-pVDZ) (Hartree): -524.98542696
E(MP3/Aug-CC-pVDZ) (Hartree): -524.97941667
E(PMP2/Aug-CC-pVDZ) (Hartree): -524.99602995
E(PMP3/Aug-CC-pVDZ) (Hartree): -524.98651874
E(PUHF/Aug-CC-pVDZ) (Hartree): -523.77194651
E(UHF/Aug-CC-pVDZ) (Hartree): -523.75842729
E(RHF-RMP2/AUG-CC-PVDZ) (Hartree): -525.06953066
  T1 diagnostic: 0.000268
E(RHF-UCCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -525.50712744
E(RHF-UCCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -525.46790236
E(RHF-UCCSD-F12a/AUG-CC-PVDZ) (Hartree): -525.46349363
  T1 diagnostic: 0.016960
  D1 diagnostic: 0.056838
E(RHF-UCCSD-F12b/AUG-CC-PVDZ) (Hartree): -525.42426856
E(RHF-UCCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -525.50610936
E(RHF-UCCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -525.46688428
E(RHF-UCCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -525.51085236
E(RHF-UCCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -525.47162728
E(ROHF/AUG-CC-PVDZ) (Hartree): -523.73968032
E(UM062X/Aug-CC-pVTZ) (Hartree): -526.13876657
Electronic state : 2-A

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Cartesian coordinates (Angs):

C	1.034634	0.162510	-0.045069	
C	-0.835679	-0.057742	0.137164	
F	1.577481	-1.016995	-0.049748	
O	1.536504	1.207259	-0.069259	
F	-1.077515	-0.874796	-0.892121	
F	-1.270166	1.148662	-0.201935	
O	-0.819245	-0.449815	1.286967	
Rotational constants (GHz):	3.9225000	2.3782900	1.9750900	
Vibrational harmonic frequencies (cm-1):				
i391.6888	55.5061		168.5440	
191.1711	303.7387		425.2446	
536.0196	599.4151		643.9384	
697.6817	928.6932		1147.5914	
1224.1061	1622.1894		2020.4906	
Zero-point correction (Hartree):	0.024067			

TS.CFOCF2OH+H2O.CFOCFO+HF+H2O.a

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E(CCSD(T)/Aug-CC-pVDZ) (Hartree): -602.00307406
E(CCSD/Aug-CC-pVDZ) (Hartree): -601.95060972
  T1 diagnostic: 0.016785
E(MP2/Aug-CC-pVDZ) (Hartree): -601.94110899
E(MP3/Aug-CC-pVDZ) (Hartree): -601.92921718
E(RHF/Aug-CC-pVDZ) (Hartree): -600.43006680
E(CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -602.53227995
E(CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -602.48427164

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E (CCSD-F12a/AUG-CC-PVDZ) (Hartree): -602.48159532
 T1 diagnostic: 0.014827
 D1 diagnostic: 0.059287
 E (CCSD-F12b/AUG-CC-PVDZ) (Hartree): -602.43358701
 E (CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -602.53127136
 E (CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -602.48326305
 E (CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -602.53641579
 E (CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -602.48840748
 E (RHF/AUG-CC-PVDZ) (Hartree): -600.43006678
 E (RM062X/Aug-CC-pVTZ) (Hartree): -603.24766211
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C -0.814999 0.738943 -0.154037
 C -0.171044 -0.648419 -0.301181
 F -2.058281 0.697710 0.282002
 O -0.248261 1.747395 -0.384948
 O 0.740999 -0.808852 -1.109902
 F -1.018641 -1.630518 -0.075016
 F 0.531945 -0.515102 1.262377
 H 1.831870 -0.029684 -0.641654
 O 2.385765 0.440331 0.146763
 H 1.703336 0.091561 0.879984
 H 2.257823 1.395189 0.053410
 Rotational constants (GHz): 2.5596500 1.7880700 1.3795400
 Vibrational harmonic frequencies (cm-1):
 i634.5895 58.6533 99.6759
 201.9683 246.3559 318.5067
 359.3176 404.6756 502.9935
 536.6809 551.2843 570.4571
 641.3512 714.4591 737.5793
 823.1345 860.3918 1186.6084
 1344.6038 1453.5953 1526.9168
 1615.6893 1726.2058 1952.4157
 2077.5790 2525.7216 3816.3529
 Zero-point correction (Hartree): 0.061176

TS.CFOCF2OH+H2O.CFOCFO+HF+H2O.b

 E (CCSD(T)/Aug-CC-pVDZ) (Hartree): -602.00221312
 E (CCSD/Aug-CC-pVDZ) (Hartree): -601.94998022
 T1 diagnostic: 0.016664
 E (MP2/Aug-CC-pVDZ) (Hartree): -601.94016463
 E (MP3/Aug-CC-pVDZ) (Hartree): -601.92849928
 E (RHF/Aug-CC-pVDZ) (Hartree): -600.42936206
 E (CCSD(T)-F12a/AUG-CC-PVDZ) (Hartree): -602.53116844
 E (CCSD(T)-F12b/AUG-CC-PVDZ) (Hartree): -602.48317884
 E (CCSD-F12a/AUG-CC-PVDZ) (Hartree): -602.48071017
 T1 diagnostic: 0.014746
 D1 diagnostic: 0.058358
 E (CCSD-F12b/AUG-CC-PVDZ) (Hartree): -602.43272056
 E (CCSD-T-F12a/AUG-CC-PVDZ) (Hartree): -602.53016264
 E (CCSD-T-F12b/AUG-CC-PVDZ) (Hartree): -602.48217303
 E (CCSD[T]-F12a/AUG-CC-PVDZ) (Hartree): -602.53522266
 E (CCSD[T]-F12b/AUG-CC-PVDZ) (Hartree): -602.48723305
 E (RHF/AUG-CC-PVDZ) (Hartree): -600.42936204
 E (RM062X/Aug-CC-pVTZ) (Hartree): -603.24685084
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 1.073937 -0.599170 -0.058424
 C 0.204028 0.645199 -0.258120
 F 0.302512 -1.693366 -0.184836
 O 2.224301 -0.659527 0.143898
 O -0.647909 0.678940 -1.150882
 F 0.885128 1.735323 0.019541
 F -0.606116 0.436021 1.233275
 H -1.791649 -0.016085 -0.728148
 O -2.438286 -0.405476 0.036256
 H -1.753059 -0.104516 0.801414
 H -2.461642 -1.368882 -0.020004
 Rotational constants (GHz): 2.5726800 1.7783000 1.3496700
 Vibrational harmonic frequencies (cm-1):

i680.6466	70.7687	74.8132
205.3512	239.9946	329.6123
353.3892	401.9789	500.4400
519.0699	540.4093	578.2923
659.1031	702.6761	733.0959
827.7661	856.6118	1131.0343
1324.0065	1411.1517	1459.8003
1608.1318	1714.7500	1999.3620
2035.1924	2480.1621	3849.2532

Zero-point correction (Hartree): 0.060613

 CF3CF=CF2 + OH : M06-2X/cc-pVDZ geometry

CF3CFCF2

 E(RM062X/CC-pVDZ) (Hartree): -713.14454132
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 1.402535 -0.157078 -0.000002
 C 0.241372 0.488409 -0.000029
 C -1.119024 -0.134950 -0.000001
 F 0.223975 1.823237 -0.000021
 F 2.553768 0.463613 0.000017
 F 1.543737 -1.455671 -0.000006
 F -1.046686 -1.464353 -0.000102
 F -1.812307 0.251056 1.074853
 F -1.812409 0.251196 -1.074720
 Rotational constants (GHz): 2.5537100 1.2667000 0.9927100
 Vibrational harmonic frequencies (cm-1):
 37.6737 129.4577 175.5998
 247.7172 248.0135 367.1500
 373.9281 474.5601 518.7231
 586.9552 612.2294 669.6819
 684.1958 790.8516 1075.6346
 1260.1232 1271.3807 1295.6839
 1408.1577 1472.8593 1912.1661
 Zero-point correction (Hartree): 0.035568

CF2CFCF2OH

 E(RM062X/CC-pVDZ) (Hartree): -689.13327169
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 1.130387 -0.155677 -0.013341
 C -0.237318 0.471407 0.012577
 C -1.411410 -0.144718 -0.006079
 F -0.194039 1.815783 0.032421
 F 1.667032 -0.092489 1.234137
 F 1.050162 -1.447904 -0.344061
 F -1.572666 -1.441633 -0.024193
 F -2.554524 0.495258 -0.004832
 O 1.955663 0.450639 -0.905135
 H 1.901059 1.407668 -0.759111
 Rotational constants (GHz): 2.5271100 1.2794000 1.0000900
 Vibrational harmonic frequencies (cm-1):
 39.3905 129.6122 163.7252
 237.8446 260.9031 293.8011
 372.3207 380.1755 478.6689
 514.6170 583.7531 609.4401
 666.6879 686.9422 786.5067
 1058.5628 1153.0456 1218.6495
 1289.1613 1373.2415 1403.4459
 1477.1091 1919.5303 3830.8603
 Zero-point correction (Hartree): 0.047677

CF3CFCF2OH

 E(UM062X/CC-pVDZ) (Hartree): -788.93888163
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C -1.303451 -0.125659 -0.031349
 C -0.012458 0.591152 -0.341223
 C 1.286070 -0.107464 -0.053253
 F -0.017123 1.886270 -0.079736
 F -2.313601 0.450548 -0.687155
 F -1.164890 -1.396558 -0.483118
 F 1.227696 -0.716933 1.160464
 F 1.545173 -1.055425 -0.945956
 F 2.293452 0.753572 -0.023522

O	-1.631398	-0.113679	1.288385	
H	-0.906144	-0.532003	1.779079	
Rotational constants (GHz):	2.0845200		1.0701800	0.9353500
Vibrational harmonic frequencies (cm-1):				
48.3876		68.2928		150.0408
172.5356		253.2285		295.2930
326.2258		347.4436		364.4615
460.5885		514.9848		540.8673
554.7354		619.6182		691.2520
710.6764		801.7906		1006.1738
1160.7549		1173.7163		1224.9472
1264.1820		1315.2839		1390.6640
1434.8583		1480.1216		3820.5567
Zero-point correction (Hartree): 0.050556				

CF3CFOHCF2

E(UM062X/CC-pVDZ) (Hartree): -788.92967318

Electronic state : 2-A

Cartesian coordinates (Angs):

C	1.492087	0.117492	-0.010636	
C	0.093616	0.696578	-0.072352	
C	-0.996979	-0.381197	0.052457	
F	-0.042576	1.250952	-1.315383	
F	1.828803	-0.668544	-1.012126	
F	1.818916	-0.409933	1.155231	
F	-0.879622	-1.016590	1.216775	
F	-2.205770	0.161420	-0.022284	
F	-0.865974	-1.271564	-0.930083	
O	-0.129021	1.596847	0.924706	
H	0.615840	2.216318	0.956367	
Rotational constants (GHz):	1.7887100		1.1952500	1.1640100
Vibrational harmonic frequencies (cm-1):				
29.8252		88.1596		147.0810
211.2630		228.0769		278.2800
323.2286		348.1811		379.4977
400.5734		514.3736		524.6588
557.5784		613.5478		642.0341
700.0667		792.8711		1007.5044
1164.5180		1264.7544		1279.1357
1298.8124		1306.6746		1359.8392
1430.6224		1453.2474		3814.6291
Zero-point correction (Hartree): 0.050482				

CF2CFCF2OH

E(RM062X/CC-pVDZ) (Hartree): -689.13327169

Electronic state : 1-A

Cartesian coordinates (Angs):

C	1.130387	-0.155677	-0.013341	
C	-0.237318	0.471407	0.012577	
C	-1.411410	-0.144718	-0.006079	
F	-0.194039	1.815783	0.032421	
F	1.667032	-0.092489	1.234137	
F	1.050162	-1.447904	-0.344061	
F	-1.572666	-1.441633	-0.024193	
F	-2.554524	0.495258	-0.004832	
O	1.955663	0.450639	-0.905135	
H	1.901059	1.407668	-0.759111	
Rotational constants (GHz):	2.5271100		1.2794000	1.0000900
Vibrational harmonic frequencies (cm-1):				
39.3905		129.6122		163.7252
237.8446		260.9031		293.8011
372.3207		380.1755		478.6689
514.6170		583.7531		609.4401
666.6879		686.9422		786.5067
1058.5628		1153.0456		1218.6495
1289.1613		1373.2415		1403.4459
1477.1091		1919.5303		3830.8603
Zero-point correction (Hartree): 0.047677				

 CF3OCF=CF2 + Cl : M06-2X/cc-pVDZ geometry

CF3OCFCF2

 E(RM062X/CC-pVDZ) (Hartree): -788.35463805
 Electronic state : 1-A
 Cartesian coordinates (Angs):
 C 1.672616 -0.287932 -0.068459
 C 0.694091 0.564996 -0.335089
 O -0.488561 0.168681 -0.874721
 F 0.823924 1.869870 -0.148821
 C -1.471733 -0.111240 0.043085
 F -1.683853 0.918544 0.854497
 F -1.141887 -1.159607 0.790339
 F -2.570402 -0.380072 -0.628547
 F 1.566372 -1.581829 -0.246295
 F 2.843472 0.072607 0.396666
 Rotational constants (GHz): 2.3475600 0.9395300 0.7772000
 Vibrational harmonic frequencies (cm-1):
 42.7445 67.2912 118.5448
 192.6827 210.7831 341.5436
 369.1534 452.4712 469.3770
 546.3894 554.7824 607.0276
 641.9512 652.4408 726.3140
 856.0562 946.8563 1225.8433
 1270.7324 1303.7518 1382.7947
 1388.7653 1429.0729 1961.5734
 Zero-point correction (Hartree): 0.040458

Cl

 E(UM062X/CC-pVDZ) (Hartree): -460.12141082
 Point group : OH
 Cartesian coordinates (Angs):
 Cl 0.000000 0.000000 0.000000
 Zero-point correction (Hartree): 0.000000

CF3OCFCF2Cl

 E(UM062X/CC-pVDZ) (Hartree): -1248.52010876
 Electronic state : 2-A
 Cartesian coordinates (Angs):
 C 1.169693 -0.358406 -0.014712
 C 0.279604 0.718444 -0.580233
 O -0.976187 0.397586 -0.970858
 F 0.385550 1.889082 0.023170
 C -1.870659 -0.059709 -0.021859
 F -1.799266 0.657819 1.091094
 F -1.627993 -1.327747 0.283720
 F -3.074623 0.038479 -0.541604
 F 1.012173 -1.471282 -0.730603
 F 0.841879 -0.638347 1.260388
 Cl 2.864600 0.157959 -0.059281
 Rotational constants (GHz): 1.9522300 0.6495700 0.5887900
 Vibrational harmonic frequencies (cm-1):
 41.2676 60.3426 100.3693
 154.3055 177.1877 209.7793
 272.9062 339.4980 367.3236
 436.4009 456.5008 499.3195
 557.3173 602.5320 627.1769
 646.2689 733.0174 766.8344
 907.1656 1010.7153 1129.5933
 1227.7668 1244.5698 1313.5165
 1360.0125 1381.4091 1421.3548
 Zero-point correction (Hartree): 0.041108