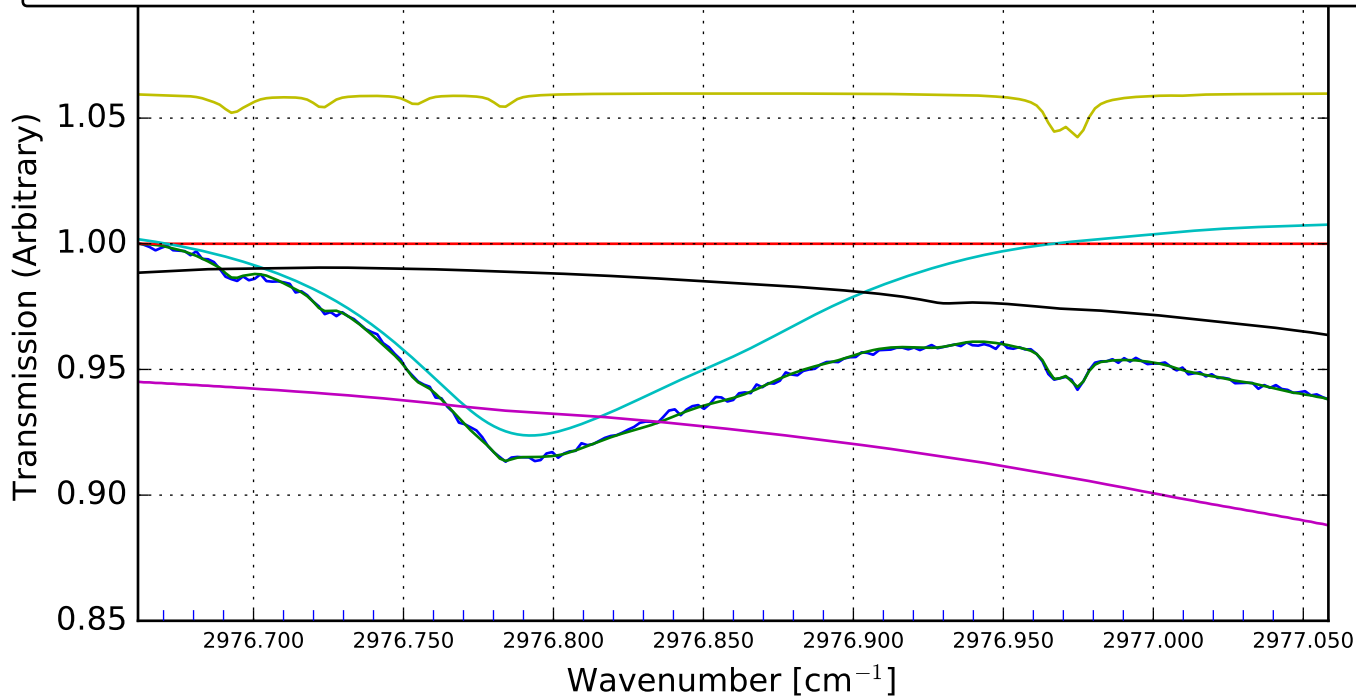
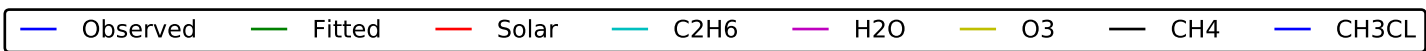
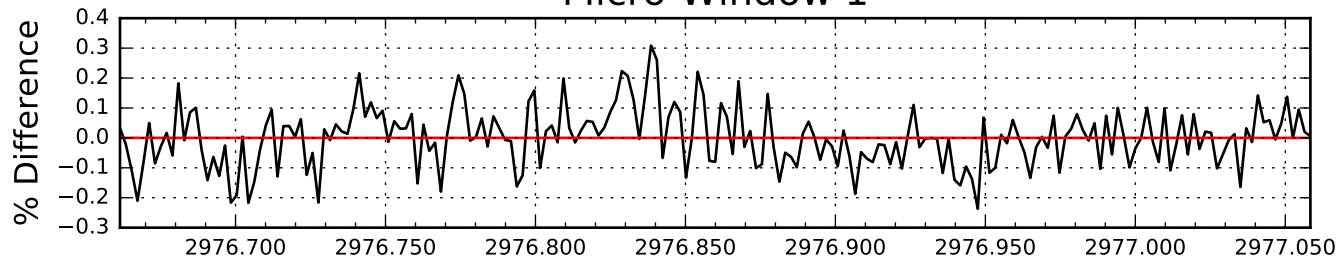
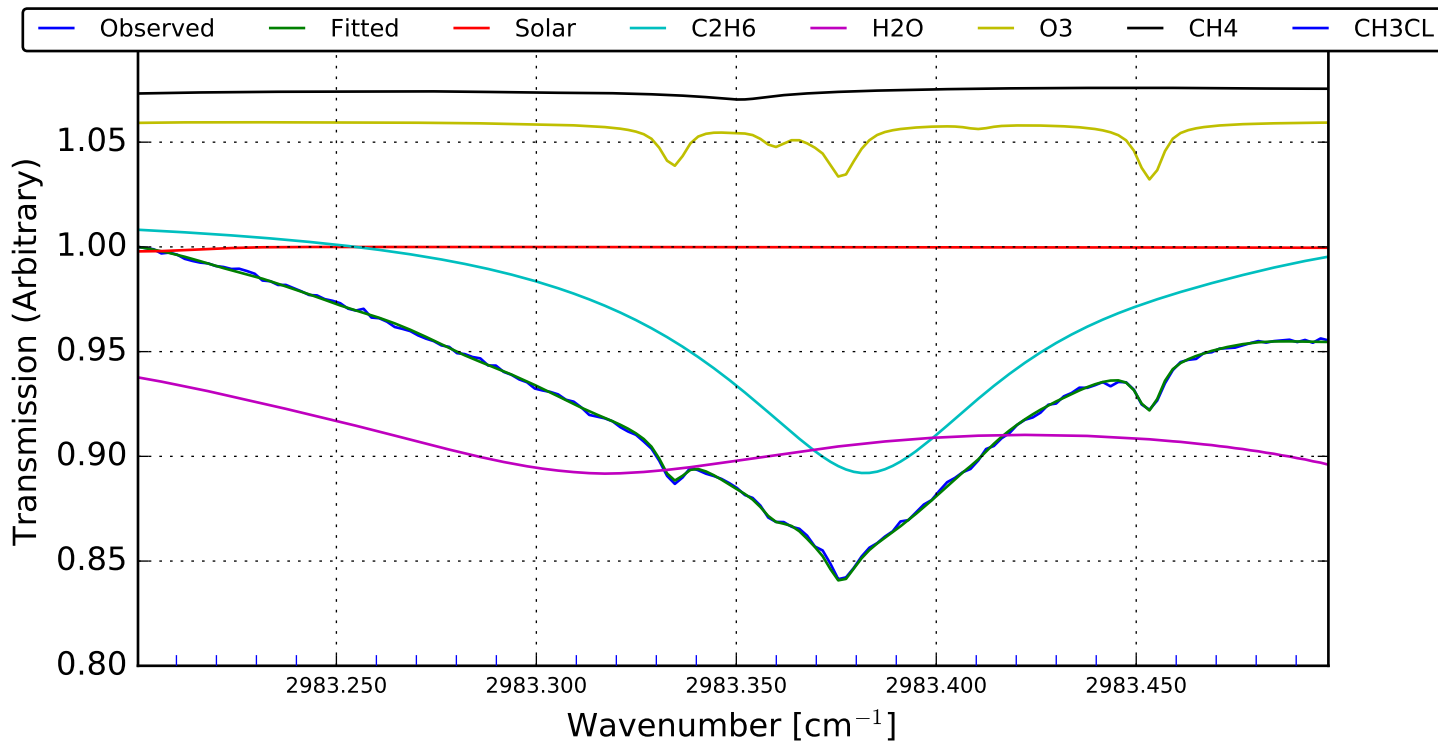
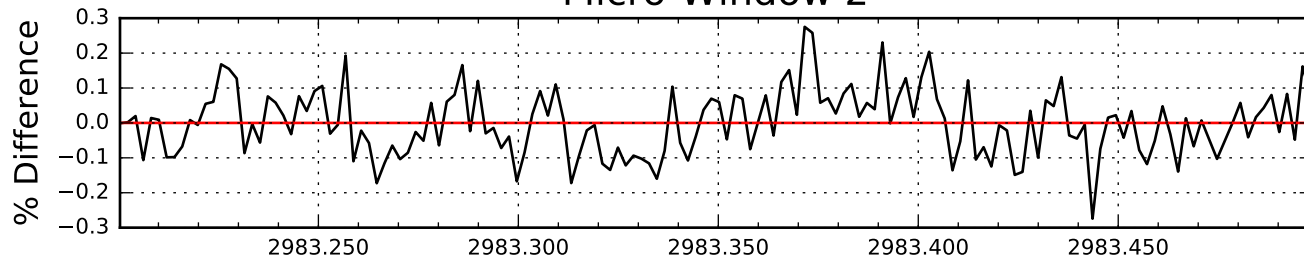


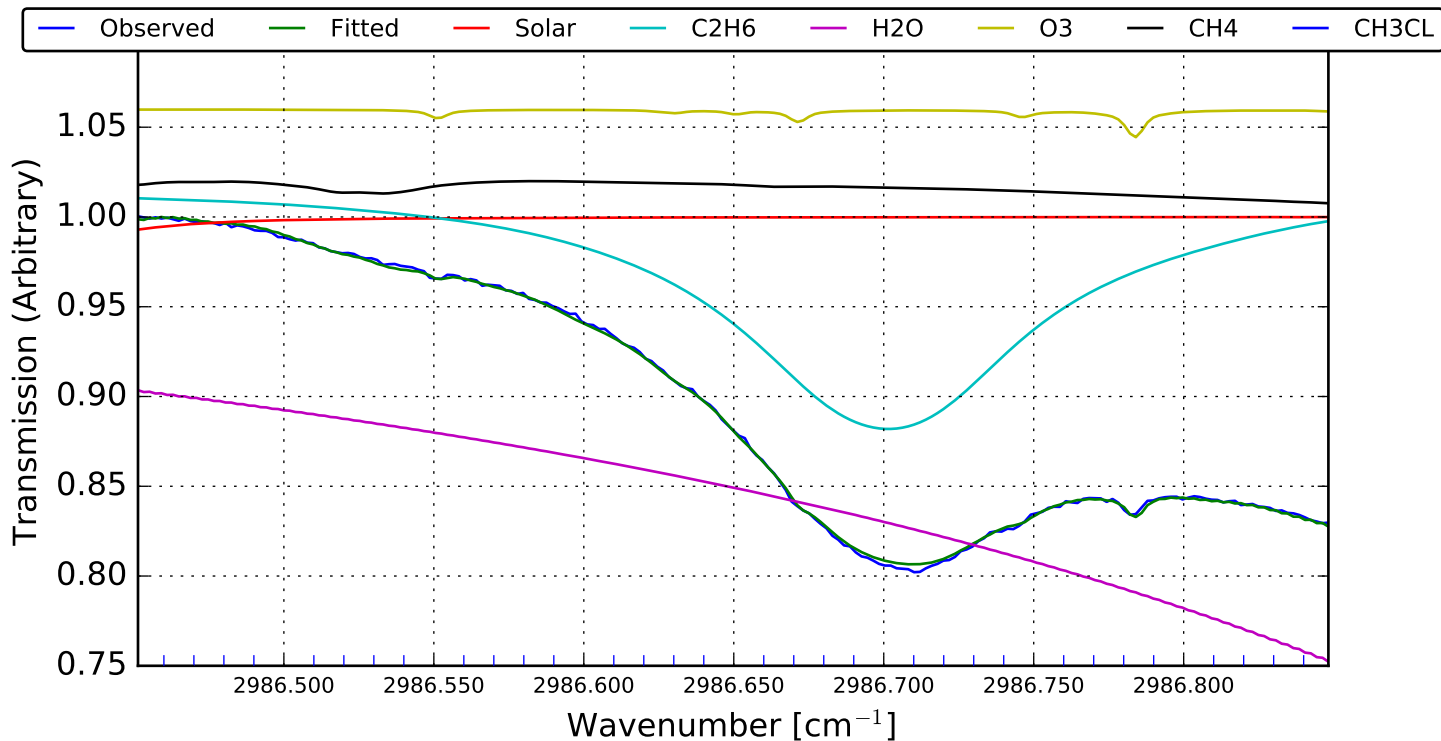
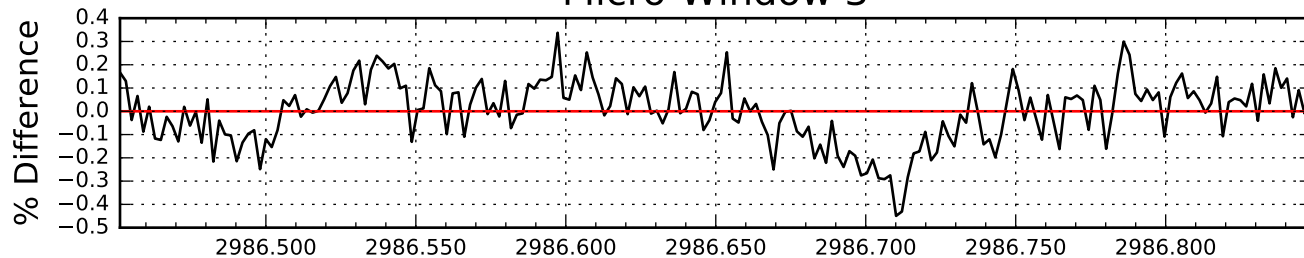
# Micro-Window 1



### Micro-Window 2



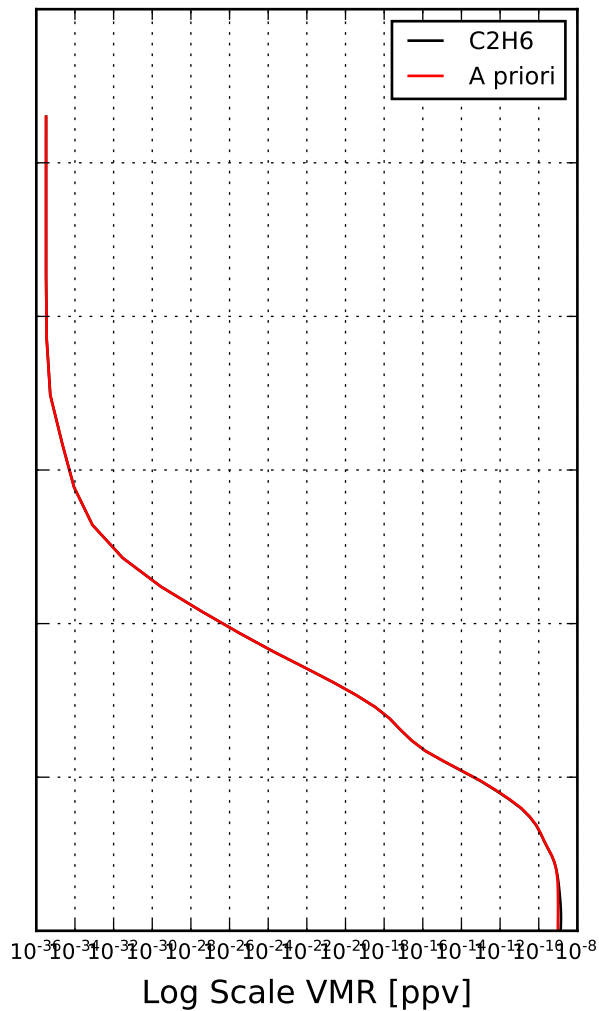
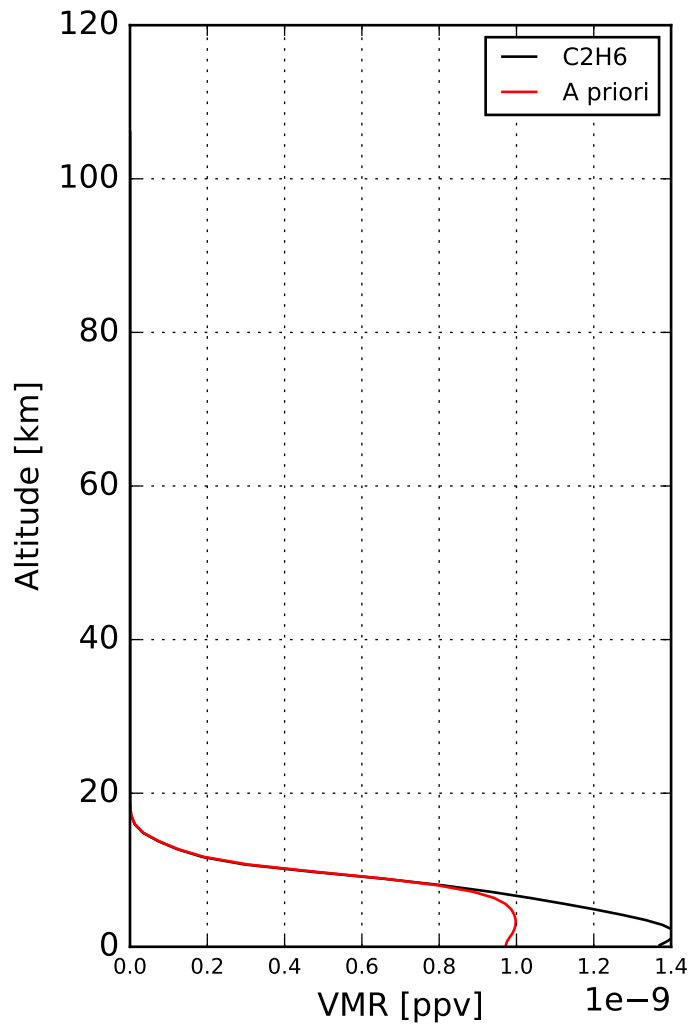
### Micro-Window 3



Number of Obs Filtered = 0

# C2H6

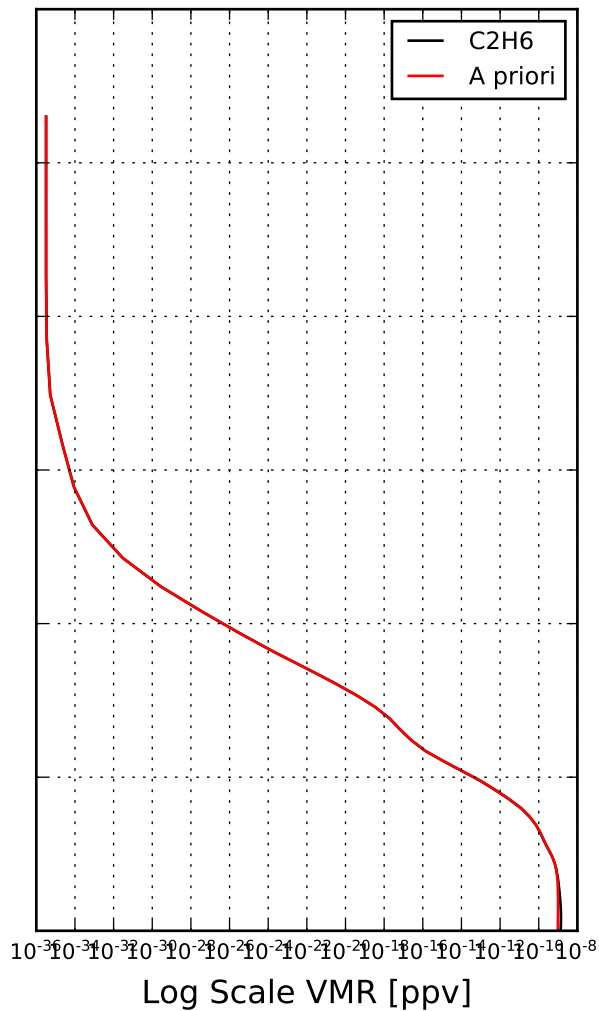
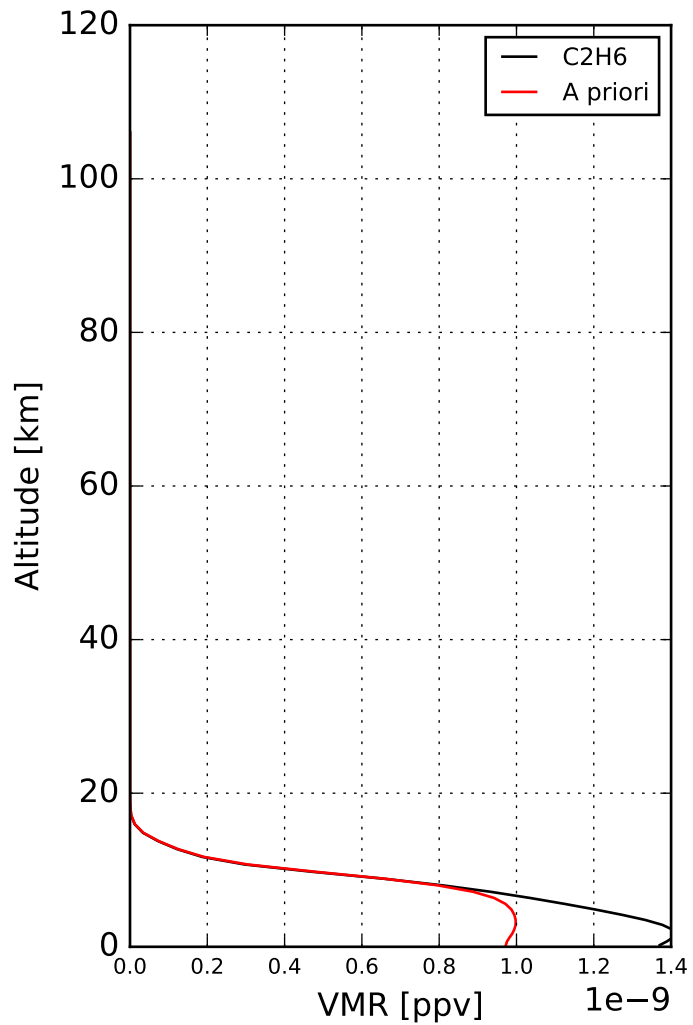
Number of Obs After Filtering = 1



Number of Obs Filtered = 0

# C2H6

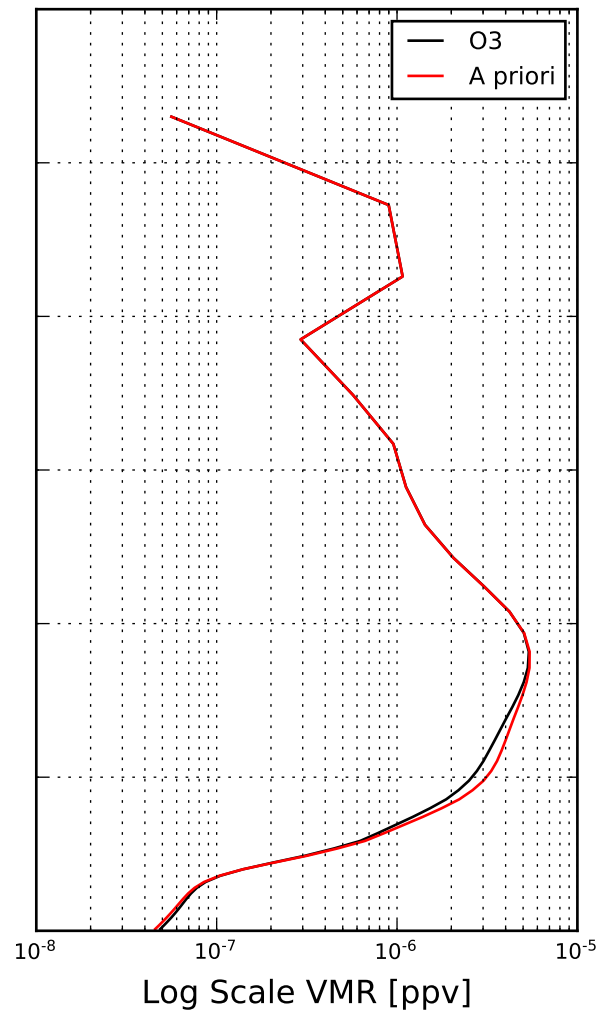
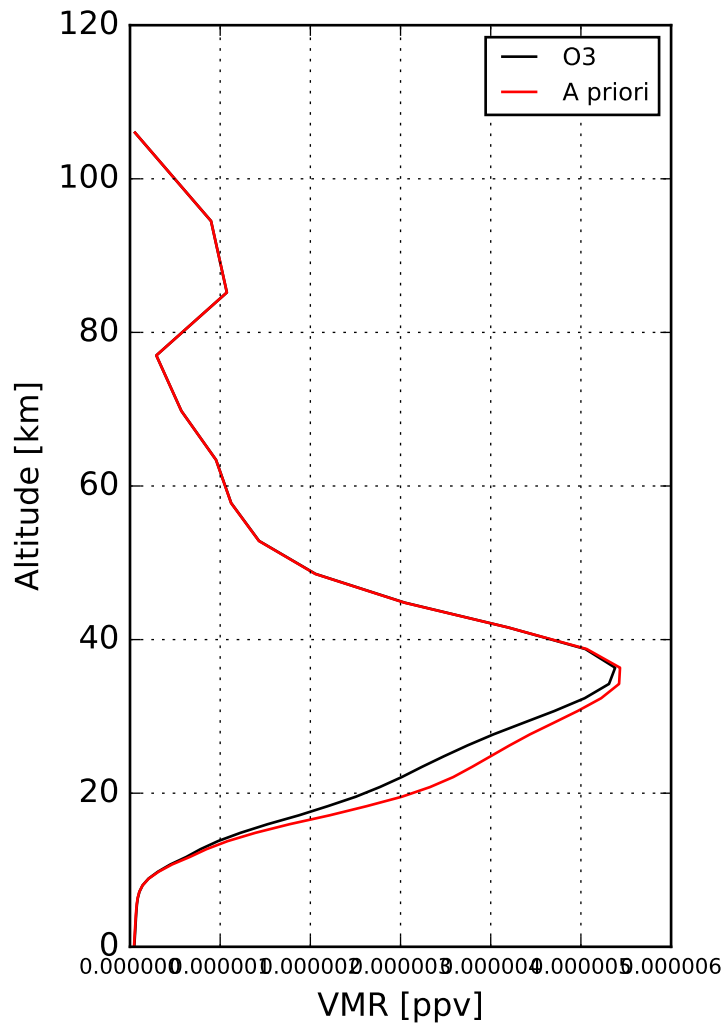
Number of Obs After Filtering = 1



Number of Obs Filtered = 0

Number of Obs After Filtering = 1

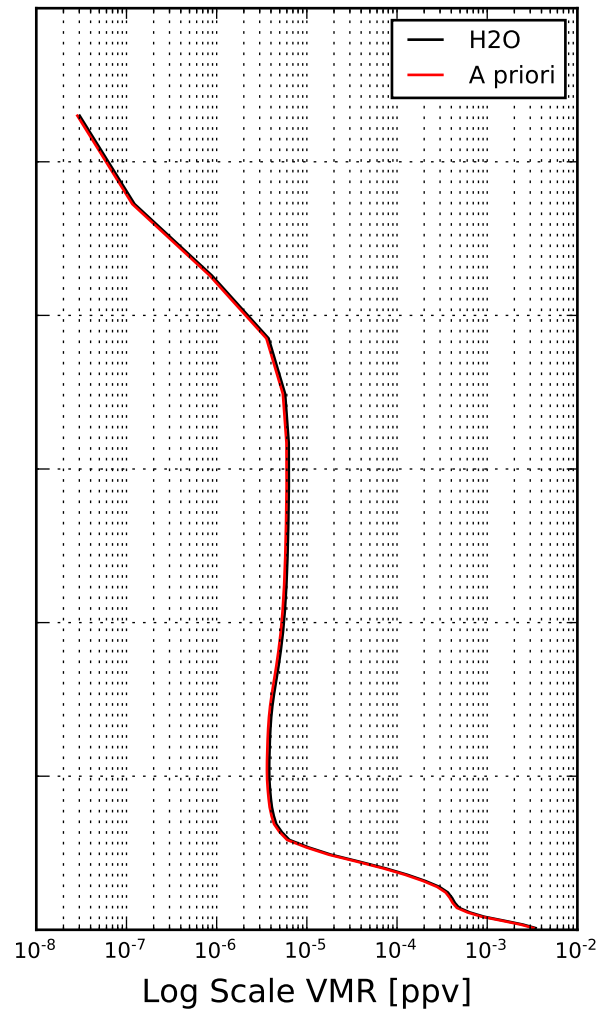
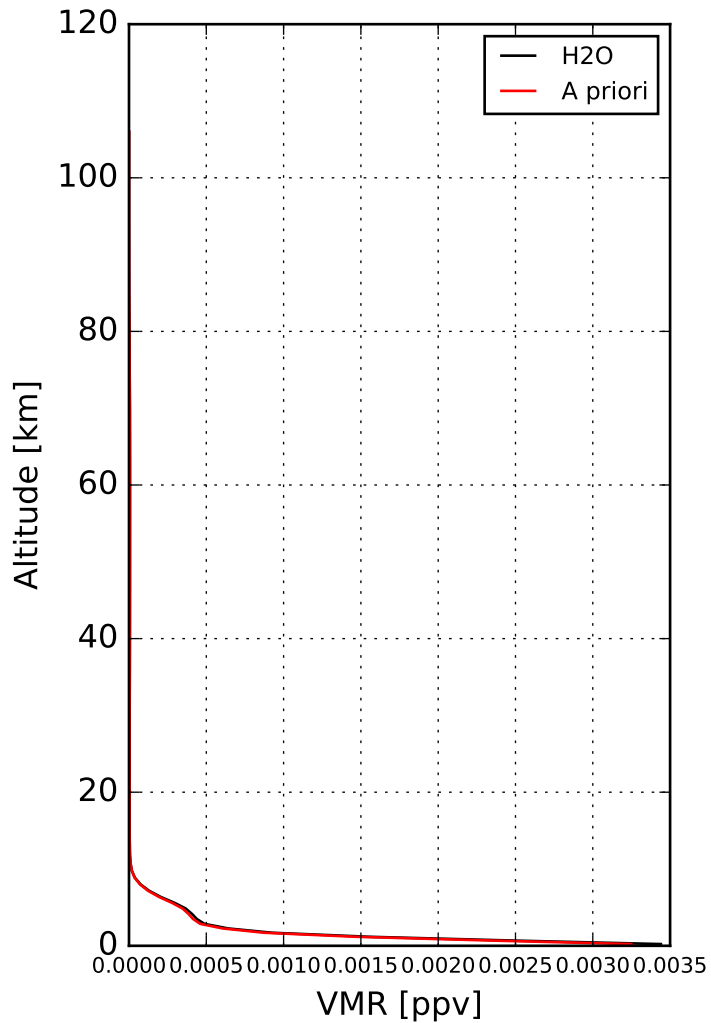
O3



Number of Obs Filtered = 0

# H2O

Number of Obs After Filtering = 1

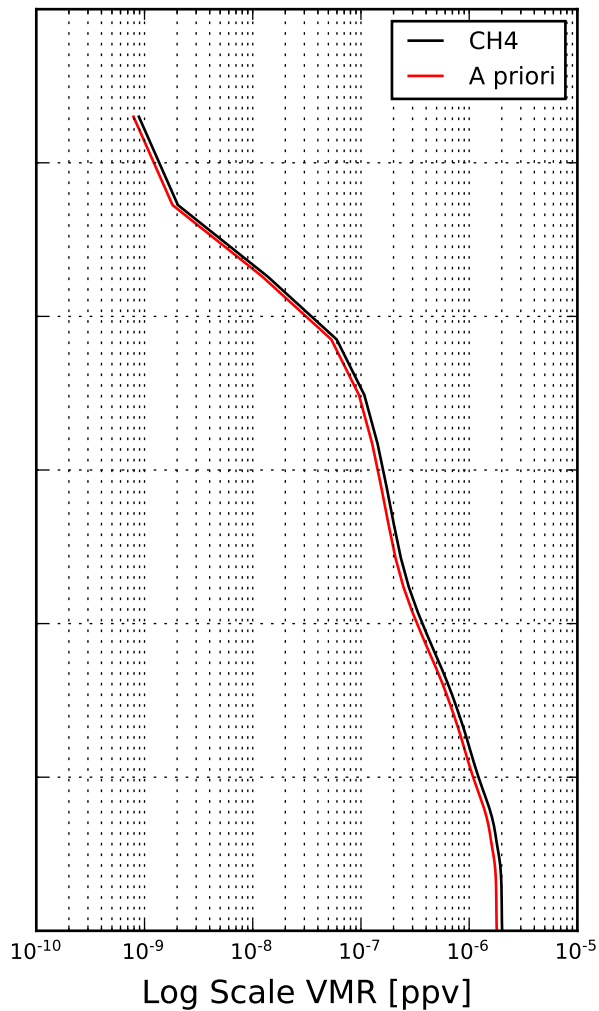
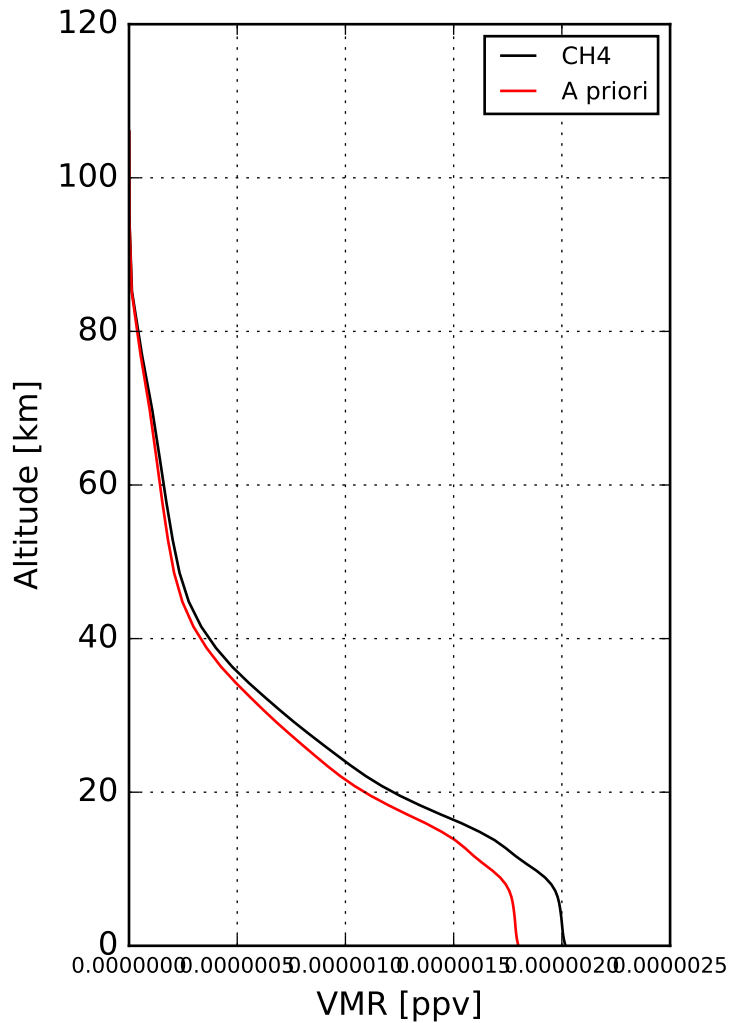




Number of Obs Filtered = 0

Number of Obs After Filtering = 1

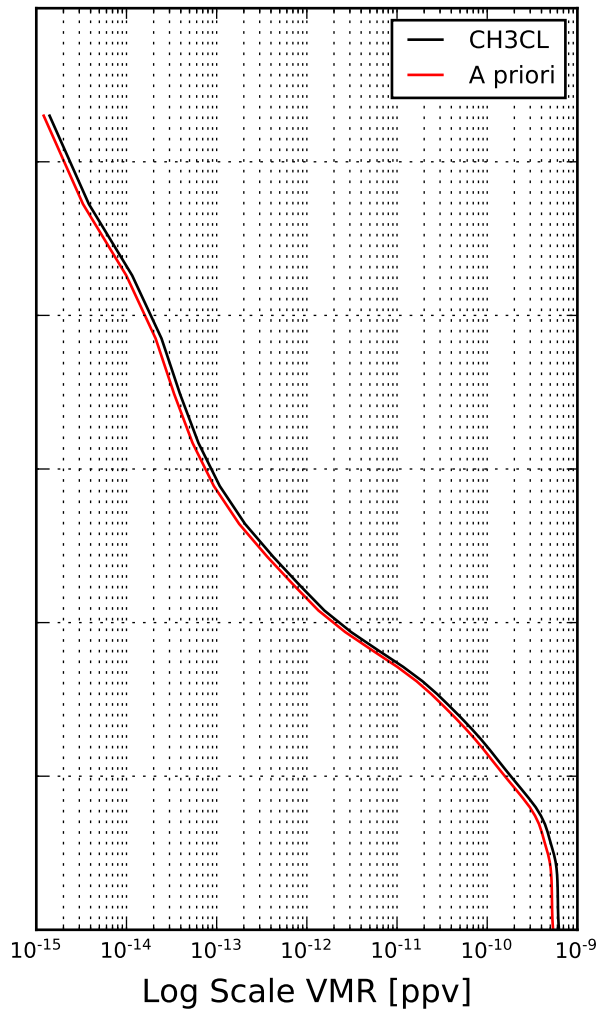
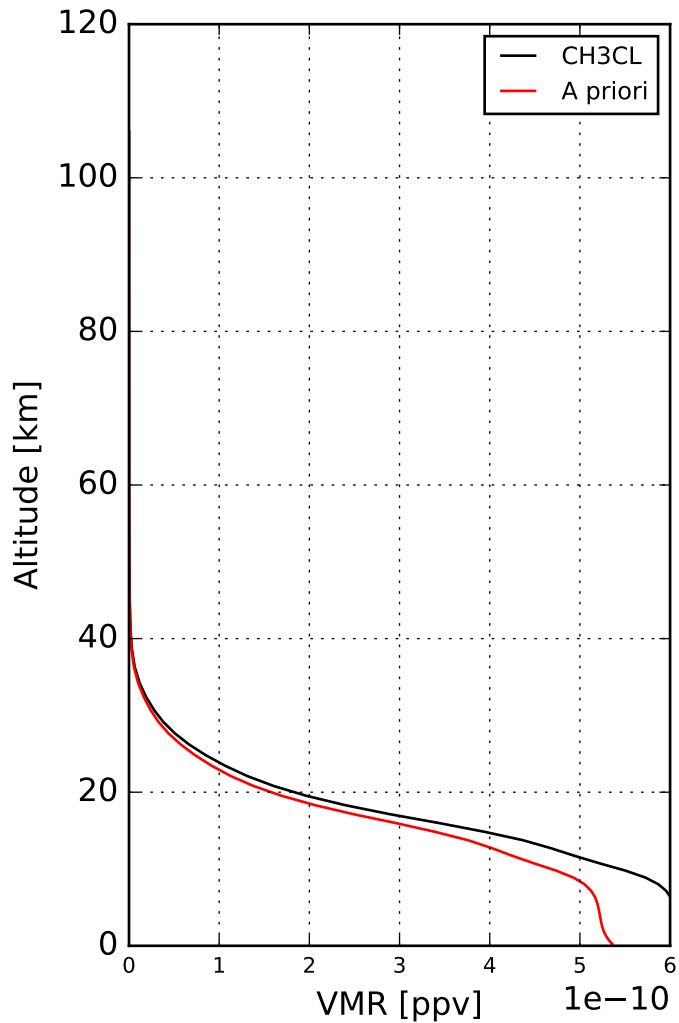
# CH4

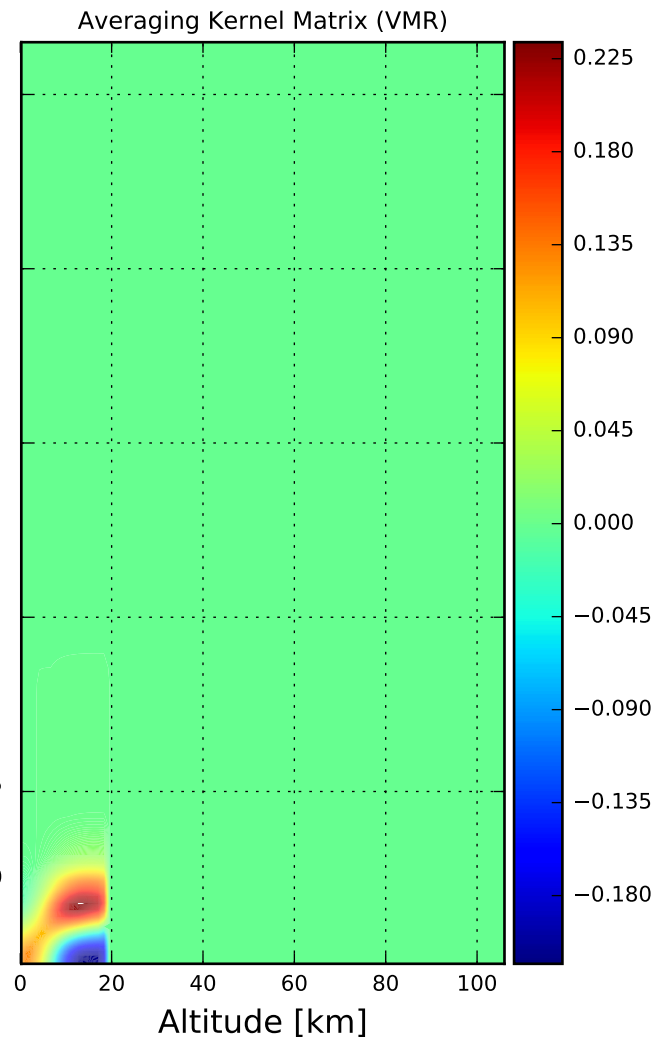
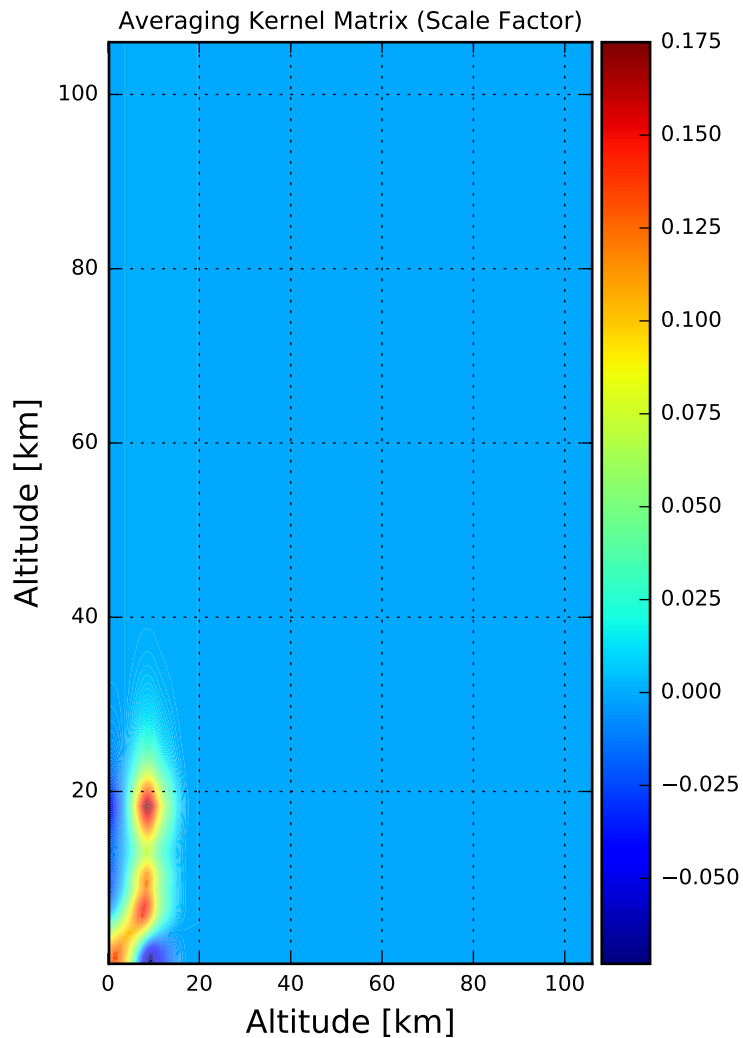


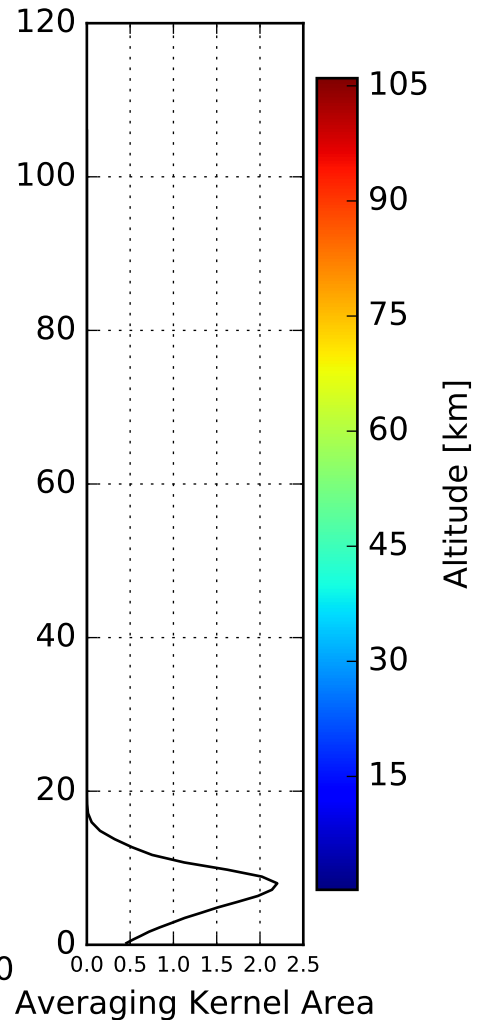
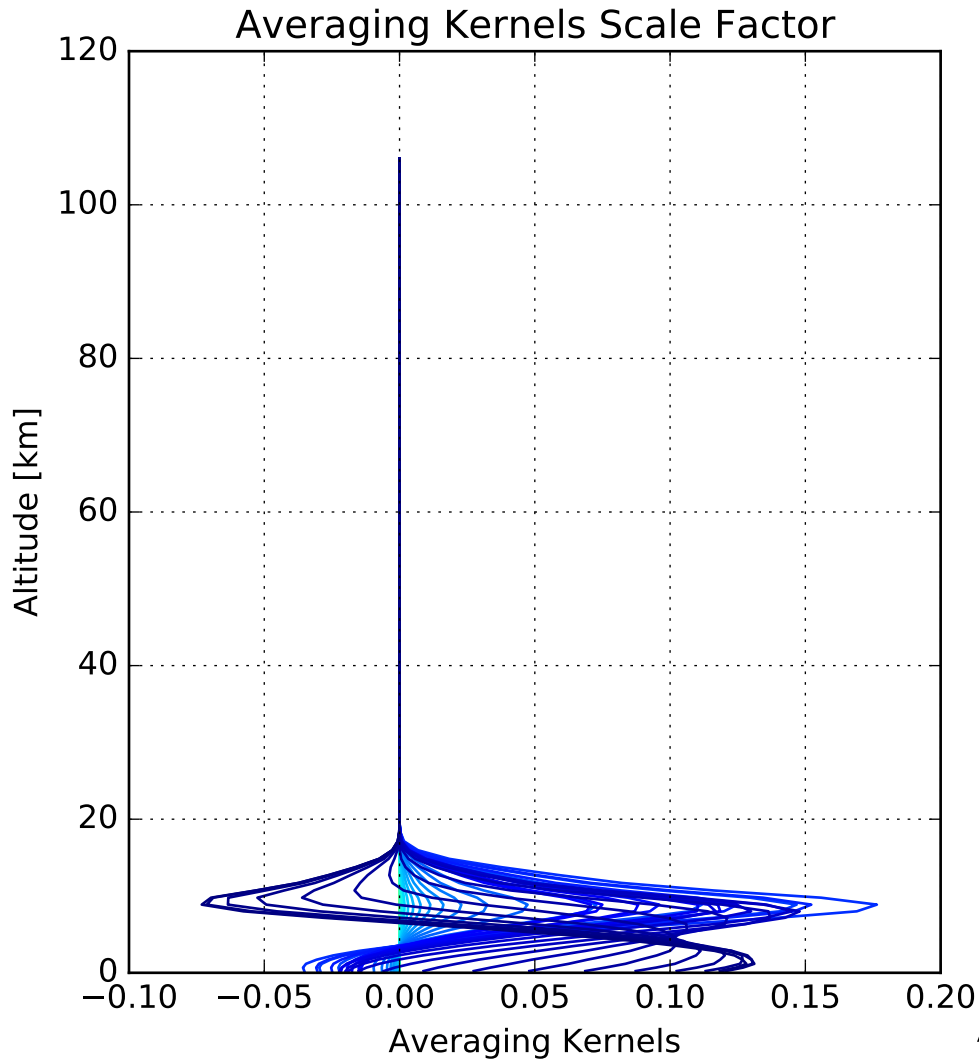
Number of Obs Filtered = 0

# CH3CL

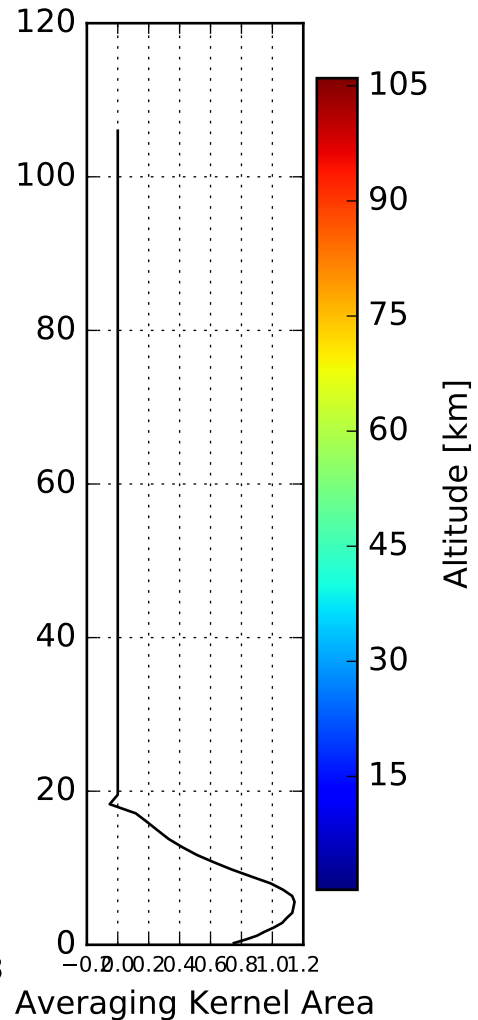
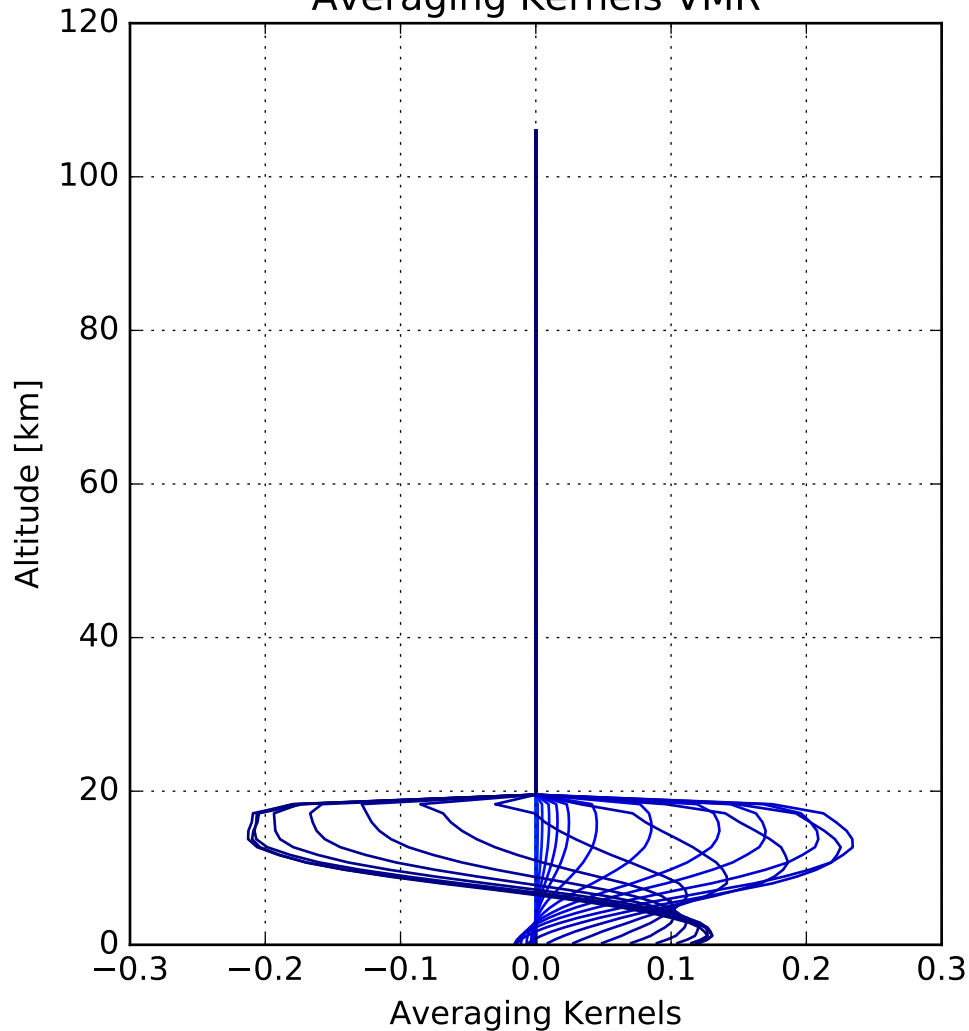
Number of Obs After Filtering = 1



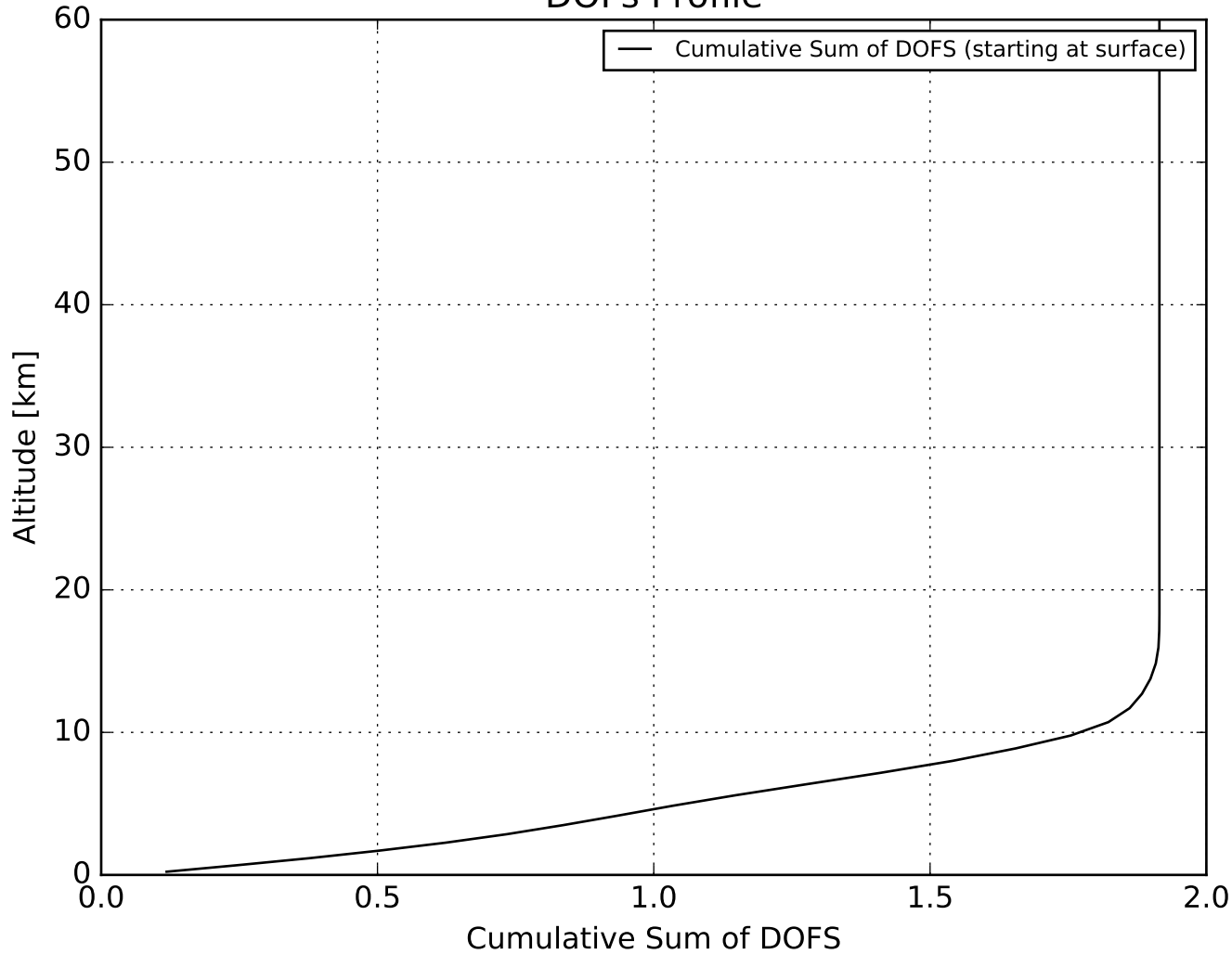




### Averaging Kernels VMR



# DOFs Profile



SFIT4: Pre-Rel:Pre-Release -- Sept 9 2014 RUNTIME:20190116-14:31:56 RETRIEVAL SUMMARY

3

20170520 13:56:01UT Z:059.701 A:316.91 D:0204.7 R:0.0035 P:BX V:02.3923 E:6397□□  
20170520 13:56:01UT Z:059.701 A:316.91 D:0204.7 R:0.0035 P:BX V:02.3923 E:6397□□  
20170520 13:56:01UT Z:059.701 A:316.91 D:0204.7 R:0.0035 P:BX V:02.3923 E:6397□□

5

IRET	GAS_NAME	IFPRF	APR_COLUMN	RET_COLUMN
1	C2H6	T	4.66631E+15	1.96177E+16
2	O3	T	1.03827E+19	9.37520E+18
3	H2O	F	1.29288E+22	1.36926E+22
4	CH4	F	3.48390E+19	3.90509E+19
5	CH3CL	F	9.76489E+15	1.13483E+16

3

IBAND	NUSTART	NUSTOP	SPACE	NPTSB	PMAX	FOVDIA	MEAN_FIT_SNR	NSCAN	JSCAN	INIT_SNR	FIT
1	2976.66000	2977.05900	0.001945525	205	257.00	2.392300	977.444062	1			
						1	1251.597	977.444			
2	2983.20000	2983.50000	0.001945525	154	257.00	2.392300	996.888659	1			
						1	1481.987	996.889			
3	2986.45000	2986.85000	0.001945525	205	257.00	2.392300	673.529970	1			
						1	1495.984	673.530			

FITRMS	CHI_2_Y	DOFS_ALL	DOFS_TRG	DOFS_TPR	ITER	MAX_ITER	CONVERGED	DIVWARN
0.120370	0.986385	15.676	1.915	0.000	5	25	T F	

Primary gas	=	C2H6	
Total column amount	=	1.96177E+16 [molecules cm <sup>-2</sup> ]	
DOFs (total column)	=	1.915	
Smoothing error (Ss, using sa+sb.ctf)	=	0.353 [%]	
Measurement error (Sm)	=	1.657 [%]	
Interference error (retrieved params)	=	0.316 [%]	
Interference error (interfering spcs)	=	0.132 [%]	
Temperature (Random)	=	1.063 [%]	
Temperature (Systematic)	=	3.261 [%]	
Total random error	=	2.490 [%]	
Total systematic error	=	14.430 [%]	
Total random uncertainty	=	4.885E+14 [molecules cm <sup>-2</sup> ]	
Total systematic uncertainty	=	2.831E+15 [molecules cm <sup>-2</sup> ]	
Total random uncertainty retrieval_parameters	=	6.192E+13 [molecules cm <sup>-2</sup> ] □	0.316 [%]
Total random uncertainty temperature	=	2.086E+14 [molecules cm <sup>-2</sup> ] □	1.063 [%]
Total random uncertainty apod_fcn	=	1.382E+12 [molecules cm <sup>-2</sup> ] □	0.007 [%]
Total random uncertainty sza	=	2.911E+14 [molecules cm <sup>-2</sup> ] □	1.484 [%]
Total random uncertainty curvature	=	1.616E+13 [molecules cm <sup>-2</sup> ] □	0.082 [%]
Total random uncertainty smoothing	=	6.927E+13 [molecules cm <sup>-2</sup> ] □	0.353 [%]
Total random uncertainty measurement	=	3.250E+14 [molecules cm <sup>-2</sup> ] □	1.657 [%]
Total random uncertainty omega	=	3.023E+10 [molecules cm <sup>-2</sup> ] □	0.000 [%]
Total random uncertainty interfering_species	=	2.593E+13 [molecules cm <sup>-2</sup> ] □	0.132 [%]
Total systematic uncertainty linepair_ch3cl	=	1.776E+13 [molecules cm <sup>-2</sup> ] □	0.091 [%]
Total systematic uncertainty linetair_c2h6	=	2.223E+13 [molecules cm <sup>-2</sup> ] □	0.113 [%]
Total systematic uncertainty linetair_ch3cl	=	3.794E+12 [molecules cm <sup>-2</sup> ] □	0.019 [%]
Total systematic uncertainty temperature	=	6.398E+14 [molecules cm <sup>-2</sup> ] □	3.261 [%]
Total systematic uncertainty linepair_h2o	=	9.921E+14 [molecules cm <sup>-2</sup> ] □	5.057 [%]
Total systematic uncertainty linepair_o3	=	2.168E+13 [molecules cm <sup>-2</sup> ] □	0.111 [%]
Total systematic uncertainty omega	=	3.023E+10 [molecules cm <sup>-2</sup> ] □	0.000 [%]
Total systematic uncertainty sza	=	2.911E+14 [molecules cm <sup>-2</sup> ] □	1.484 [%]
Total systematic uncertainty linepair_ch4	=	1.977E+14 [molecules cm <sup>-2</sup> ] □	1.008 [%]
Total systematic uncertainty lineint_c2h6	=	2.458E+15 [molecules cm <sup>-2</sup> ] □	12.531 [%]
Total systematic uncertainty curvature	=	1.616E+13 [molecules cm <sup>-2</sup> ] □	0.082 [%]
Total systematic uncertainty smoothing	=	2.714E+13 [molecules cm <sup>-2</sup> ] □	0.138 [%]
Total systematic uncertainty lineint_ch4	=	1.390E+13 [molecules cm <sup>-2</sup> ] □	0.071 [%]
Total systematic uncertainty linepair_c2h6	=	6.658E+14 [molecules cm <sup>-2</sup> ] □	3.394 [%]
Total systematic uncertainty lineint_o3	=	1.677E+13 [molecules cm <sup>-2</sup> ] □	0.086 [%]
Total systematic uncertainty linetair_ch4	=	2.411E+13 [molecules cm <sup>-2</sup> ] □	0.123 [%]
Total systematic uncertainty linetair_h2o	=	8.542E+13 [molecules cm <sup>-2</sup> ] □	0.435 [%]
Total systematic uncertainty lineint_h2o	=	7.713E+10 [molecules cm <sup>-2</sup> ] □	0.000 [%]
Total systematic uncertainty linetair_o3	=	1.402E+13 [molecules cm <sup>-2</sup> ] □	0.071 [%]
Total systematic uncertainty apod_fcn	=	1.382E+12 [molecules cm <sup>-2</sup> ] □	0.007 [%]