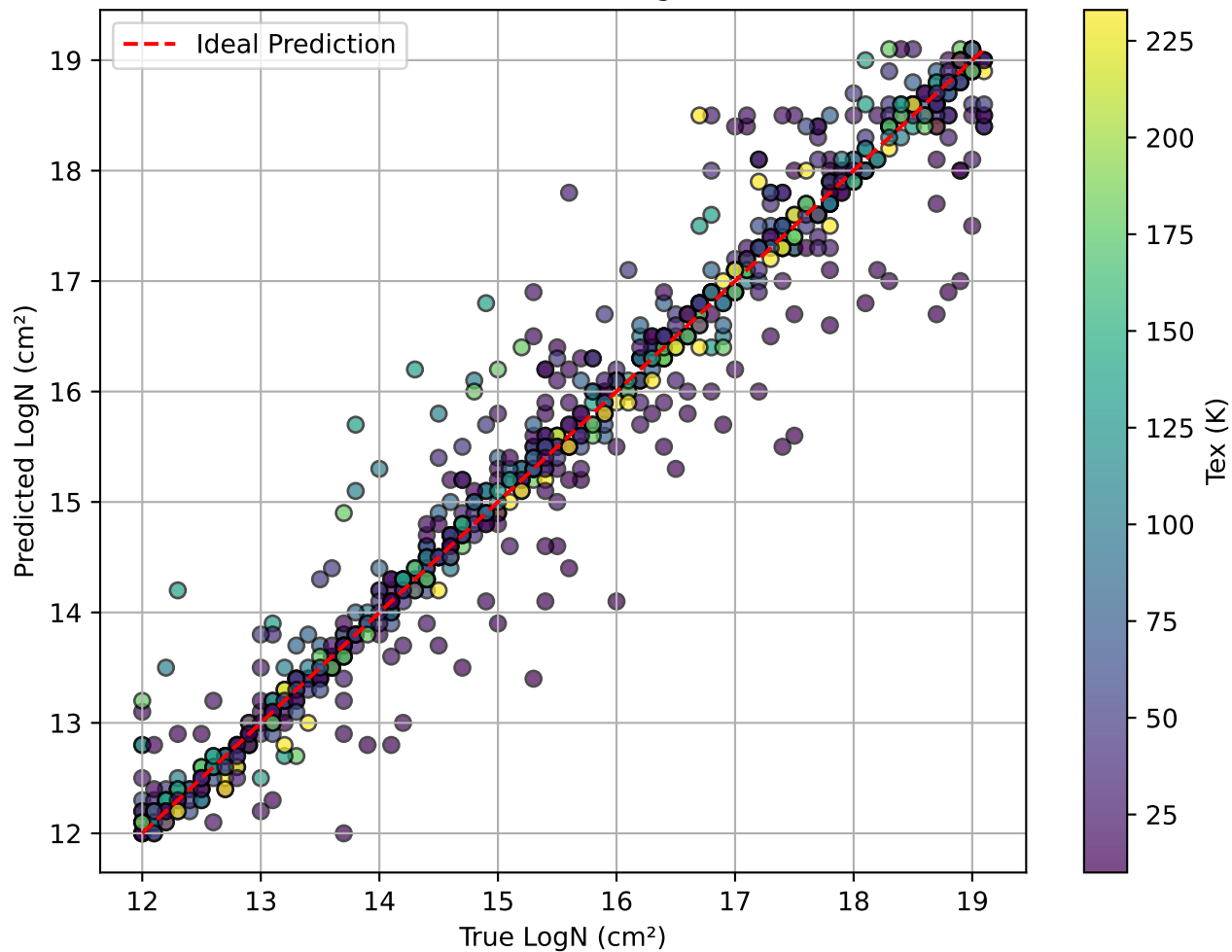
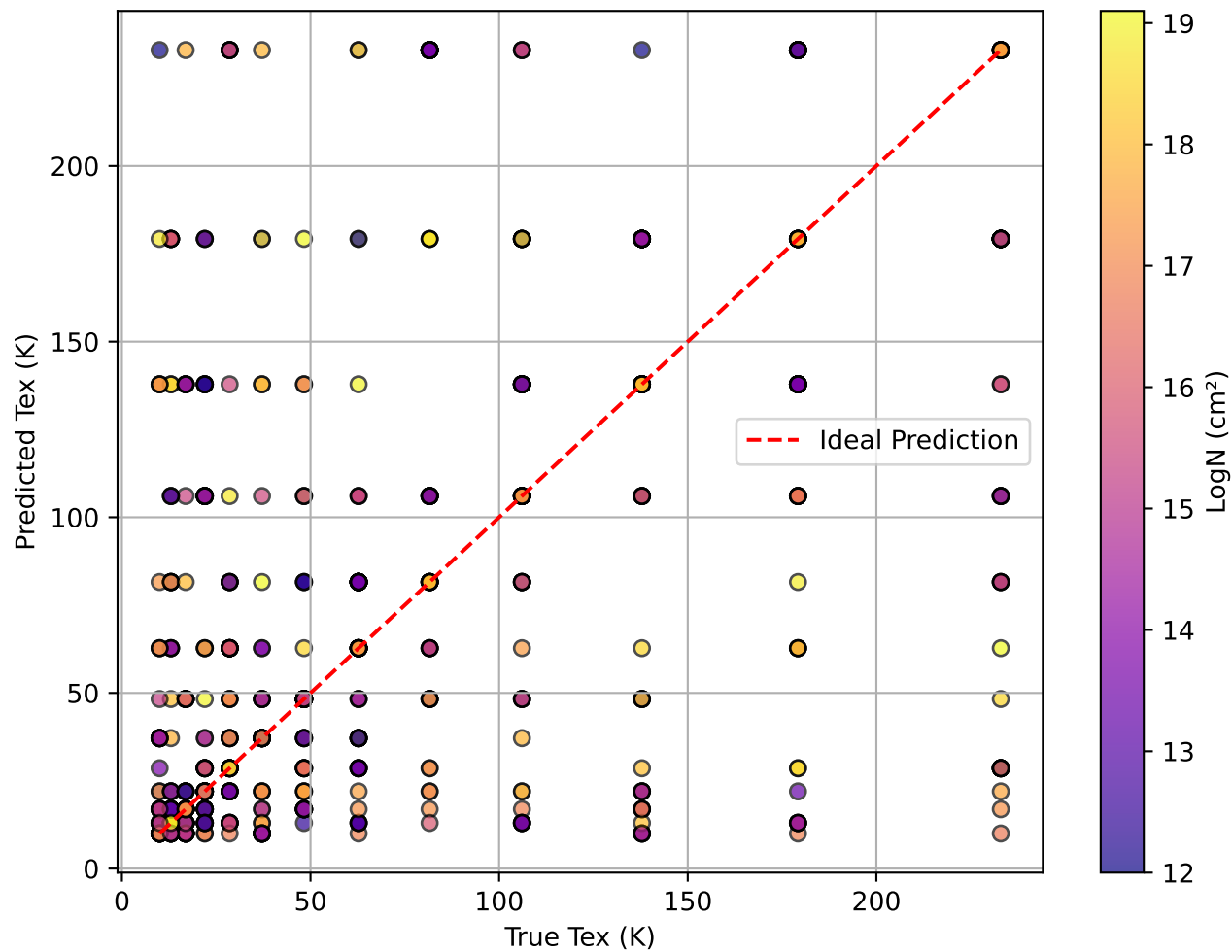


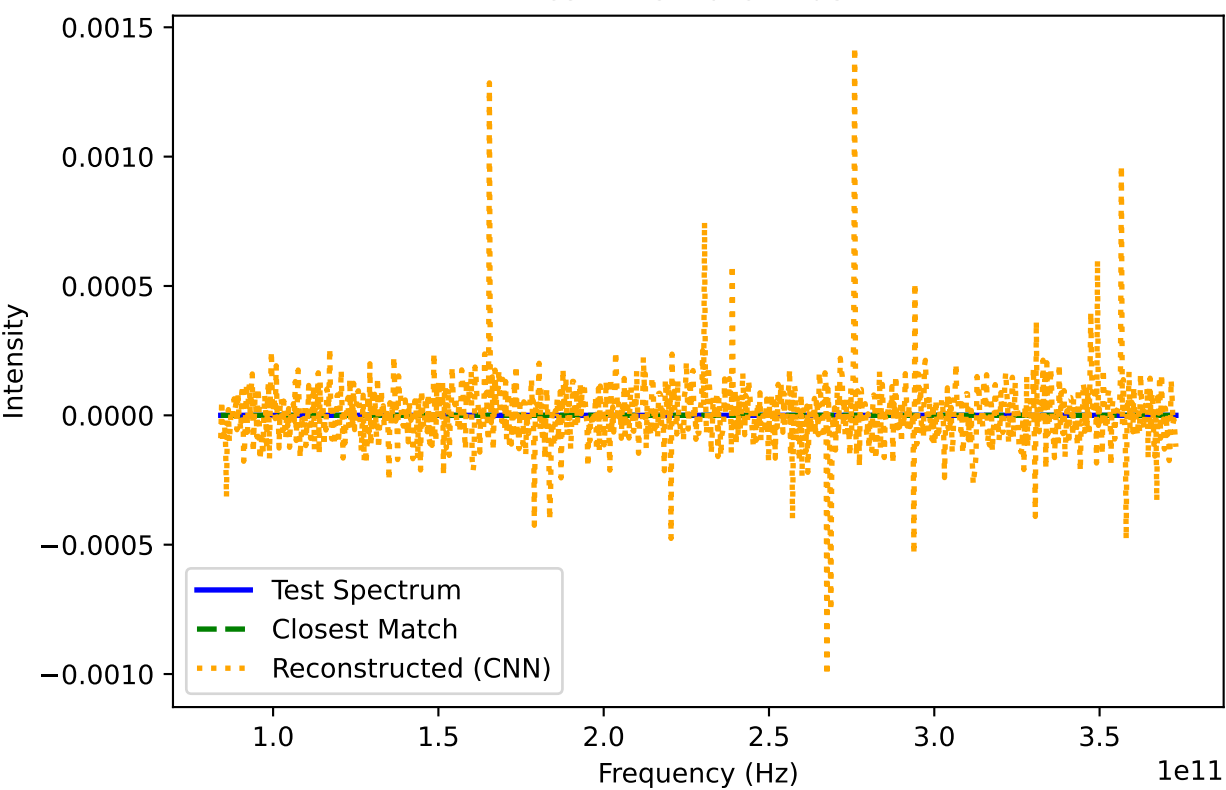
Predicted vs True LogN - ALL



Predicted vs True Tex - ALL



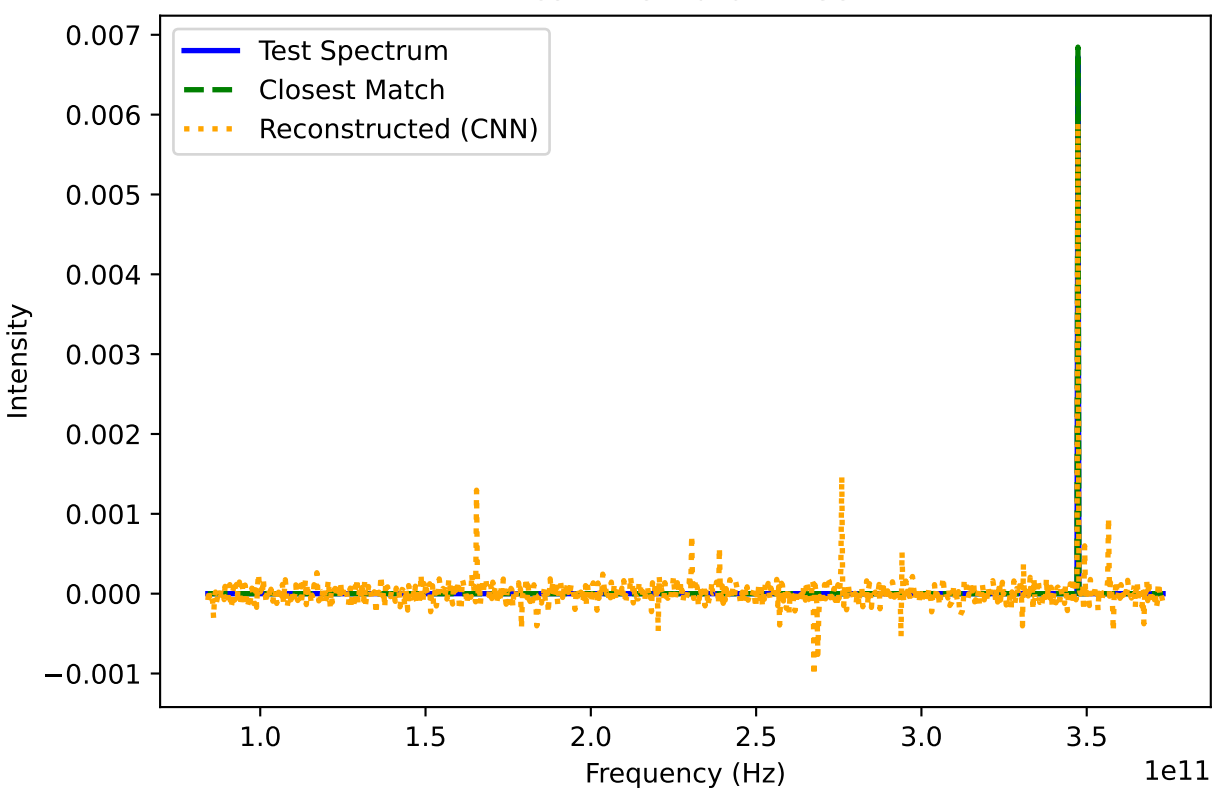
Test 1 vs Match 2087



Test File: file_JPL_CO_T16.900000000000002_N12.999999999999996_simulate_generate
Header: //molecules='CO[1]' logn=13.0 tex=16.9 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_CO_T10_N13.099999999999996_simulate_generate
Header: //molecules='CO[1]' logn=13.1 tex=10.0 velo=250.0 fwhm=50.0 sourcesize=10.0

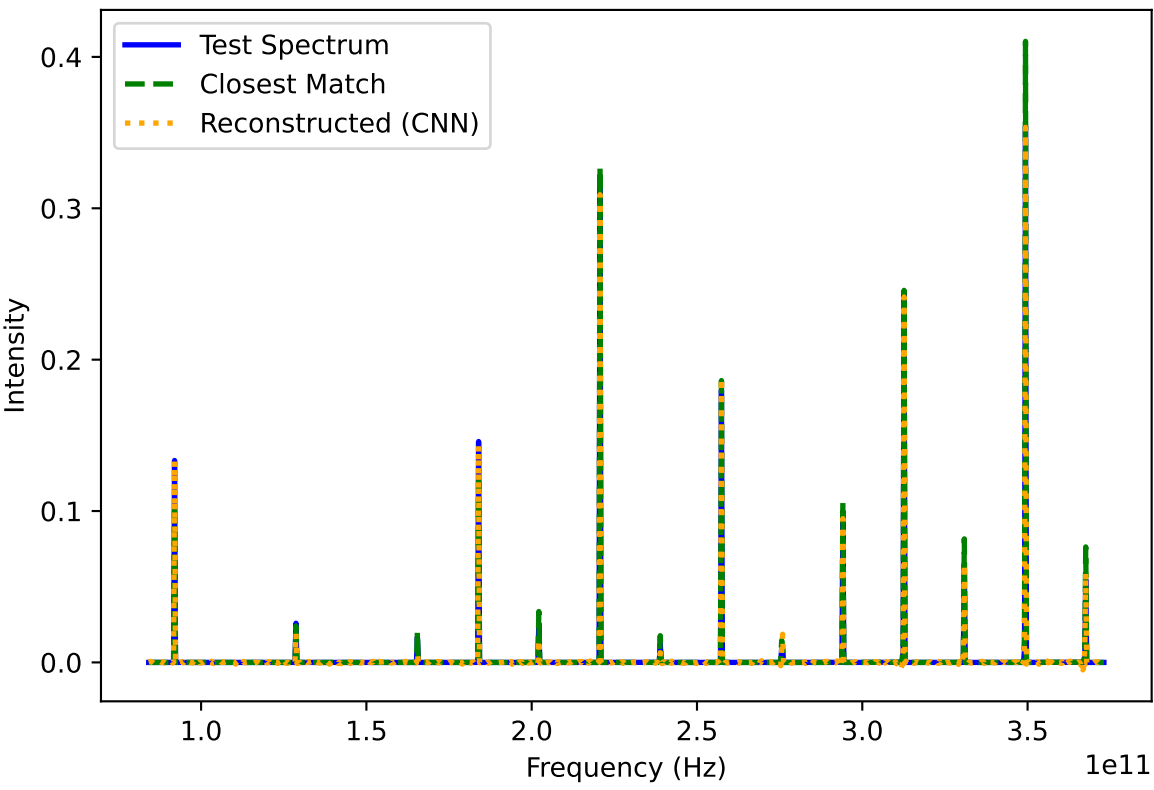
Test 2 vs Match 1735



Test File: file_JPL_SiO_T137.85849184900007_N17.1_simulate_generate
Header: //molecules='SiO[1]' logn=17.1 tex=137.85849 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_SiO_T48.268090000000015_N17.1_simulate_generate
Header: //molecules='SiO[1]' logn=17.1 tex=48.26809 velo=250.0 fwhm=50.0 sourcesize=10.0

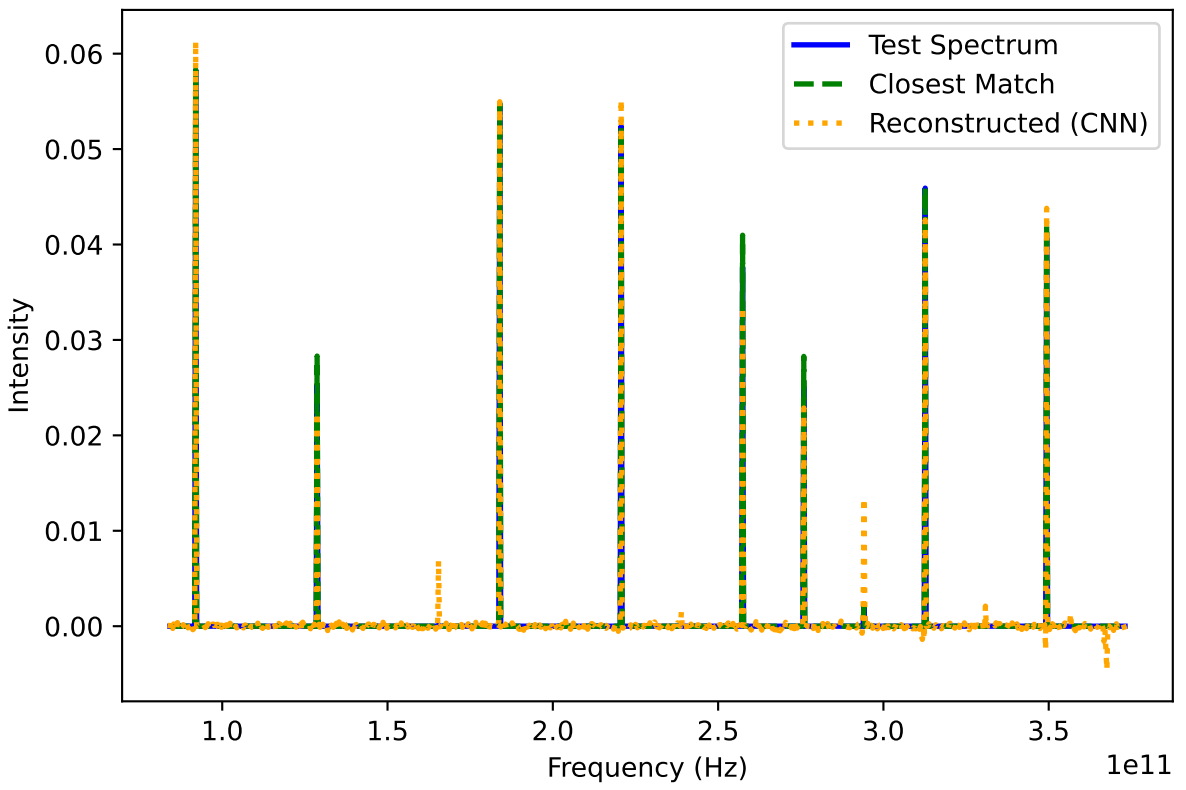
Test 3 vs Match 613



Test File: file_JPL_CH3CN_T106.04499373000004_N16.499999999999993_simulate_generate
Header: //molecules='CH3CN[1]' logn=16.5 tex=106.04499 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_CH3CN_T137.85849184900007_N16.499999999999993_simulate_generate
Header: //molecules='CH3CN[1]' logn=16.5 tex=137.85849 velo=250.0 fwhm=50.0 sourcesize=10.0

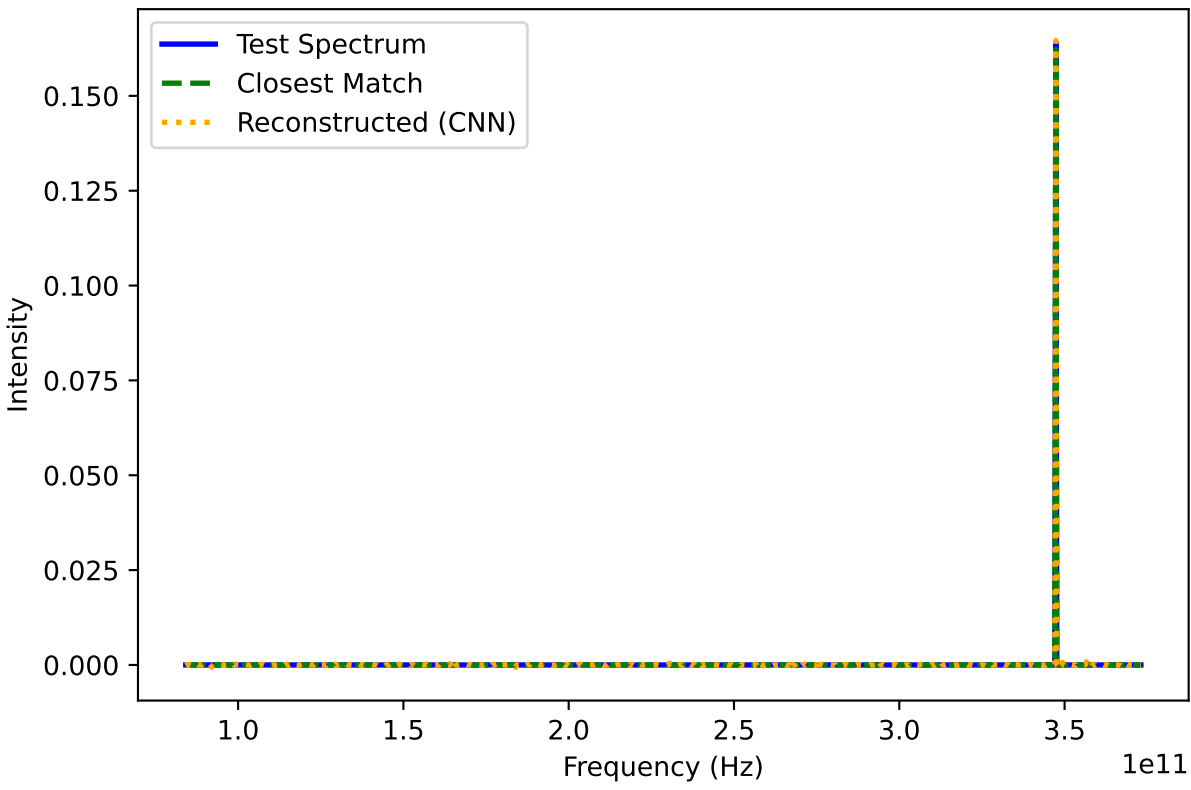
Test 4 vs Match 1114



Test File: file_JPL_CH3CN_T16.900000000000002_N18.300000000000002_simulate_generate
Header: //molecules='CH3CN[1]' logn=18.3 tex=16.9 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_CH3CN_T16.900000000000002_N18.400000000000002_simulate_generate
Header: //molecules='CH3CN[1]' logn=18.4 tex=16.9 velo=250.0 fwhm=50.0 sourcesize=10.0

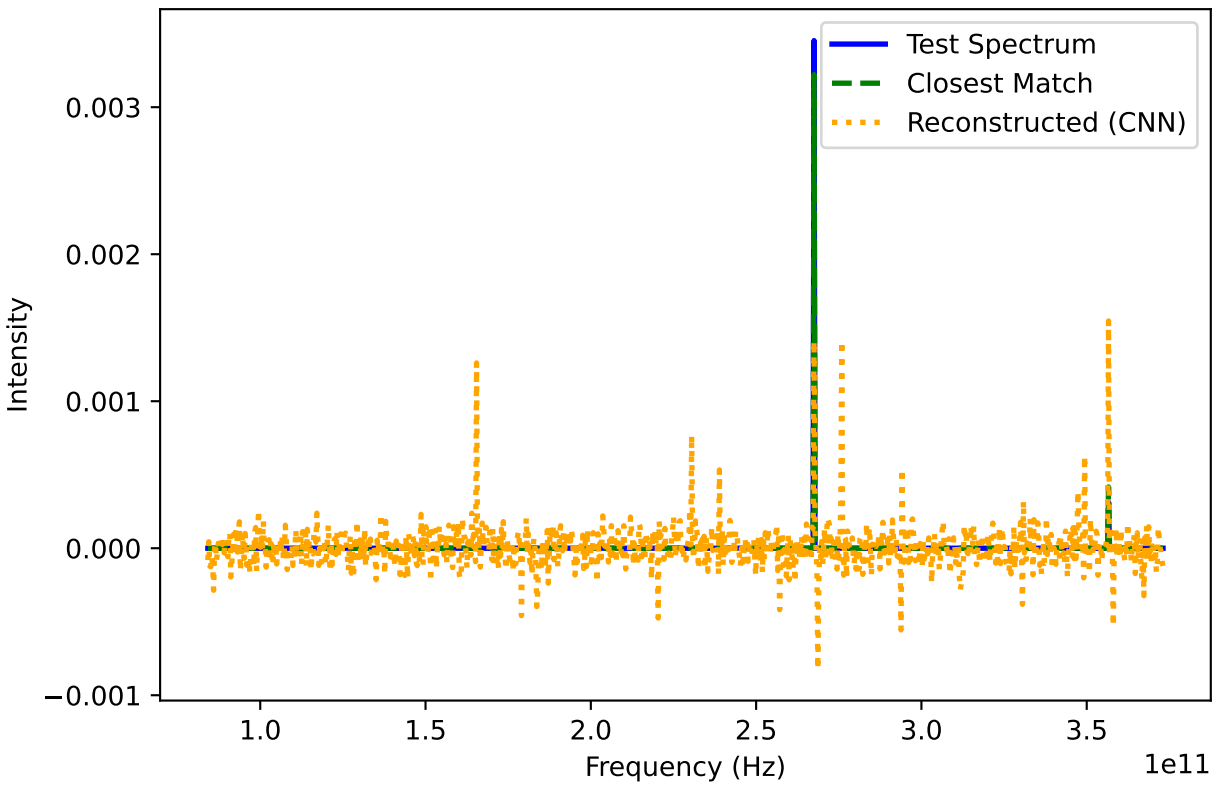
Test 5 vs Match 2625



Test File: file_JPL_SiO_T81.57307210000003_N18.600000000000023_simulate_generate
Header: //molecules='SiO[1]' logn=18.6 tex=81.573074 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_SiO_T179.2160394037001_N18.600000000000023_simulate_generate
Header: //molecules='SiO[1]' logn=18.6 tex=179.21603 velo=250.0 fwhm=50.0 sourcesize=10.0

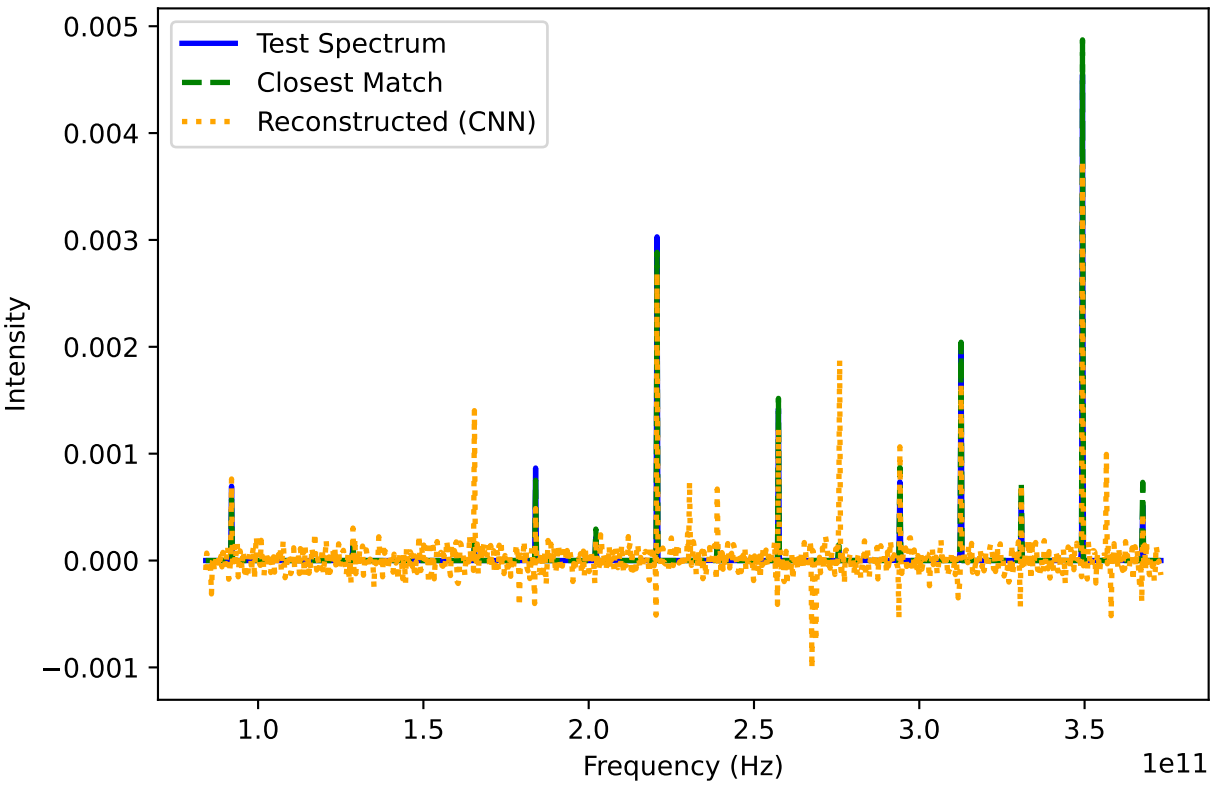
Test 6 vs Match 573



Test File: file_JPL_HCO+v=0,1,2_T10_N13.999999999999993_simulate_generate
Header: //molecules='HCO+V=0,1,2[1]' logn=14.0 tex=21.97 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_HCO+v=0,1,2_T13.0_N13.799999999999994_simulate_generate
Header: //molecules='HCO+V=0,1,2[1]' logn=13.8 tex=13.0 velo=250.0 fwhm=50.0 sourcesize=10.0

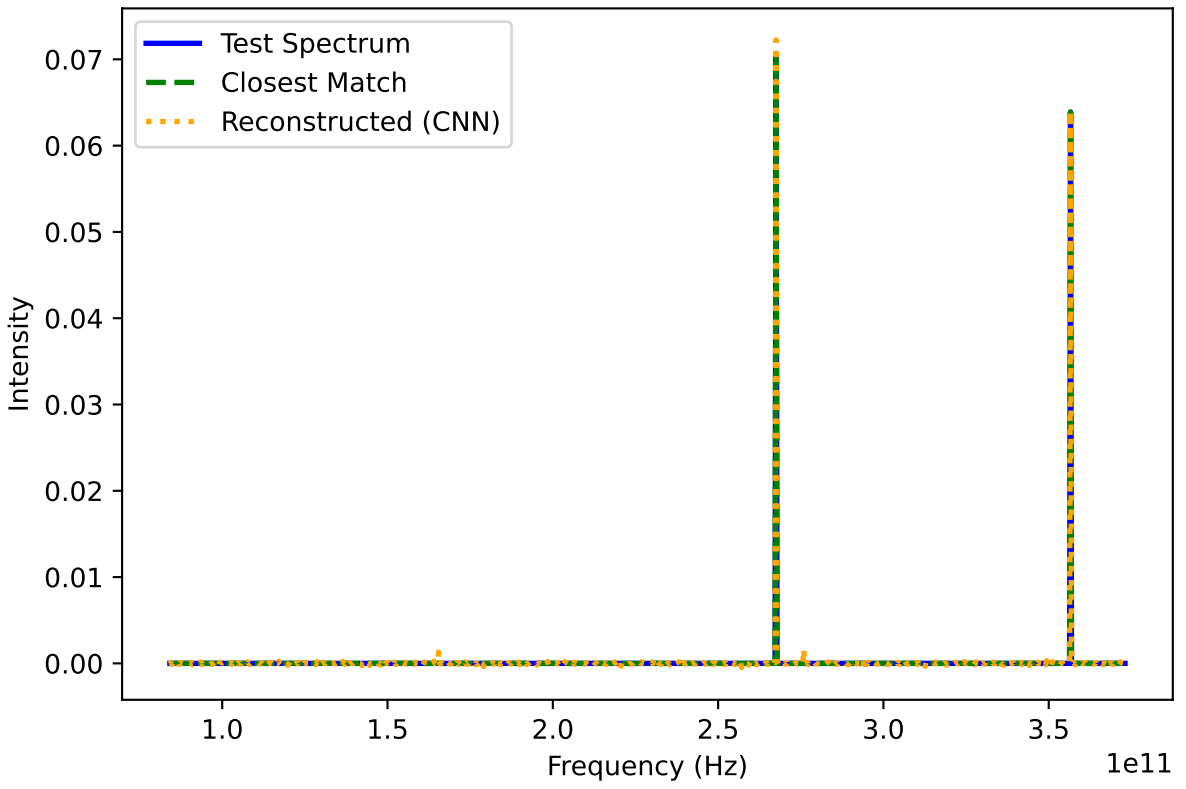
Test 7 vs Match 2150



Test File: file_JPL_CH3CN_T137.85849184900007_N14.299999999999992_simulate_generate
Header: //molecules='CH3CN[1]' logn=14.3 tex=137.85849 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_CH3CN_T179.2160394037001_N14.399999999999991_simulate_generate
Header: //molecules='CH3CN[1]' logn=14.4 tex=179.21603 velo=250.0 fwhm=50.0 sourcesize=10.0

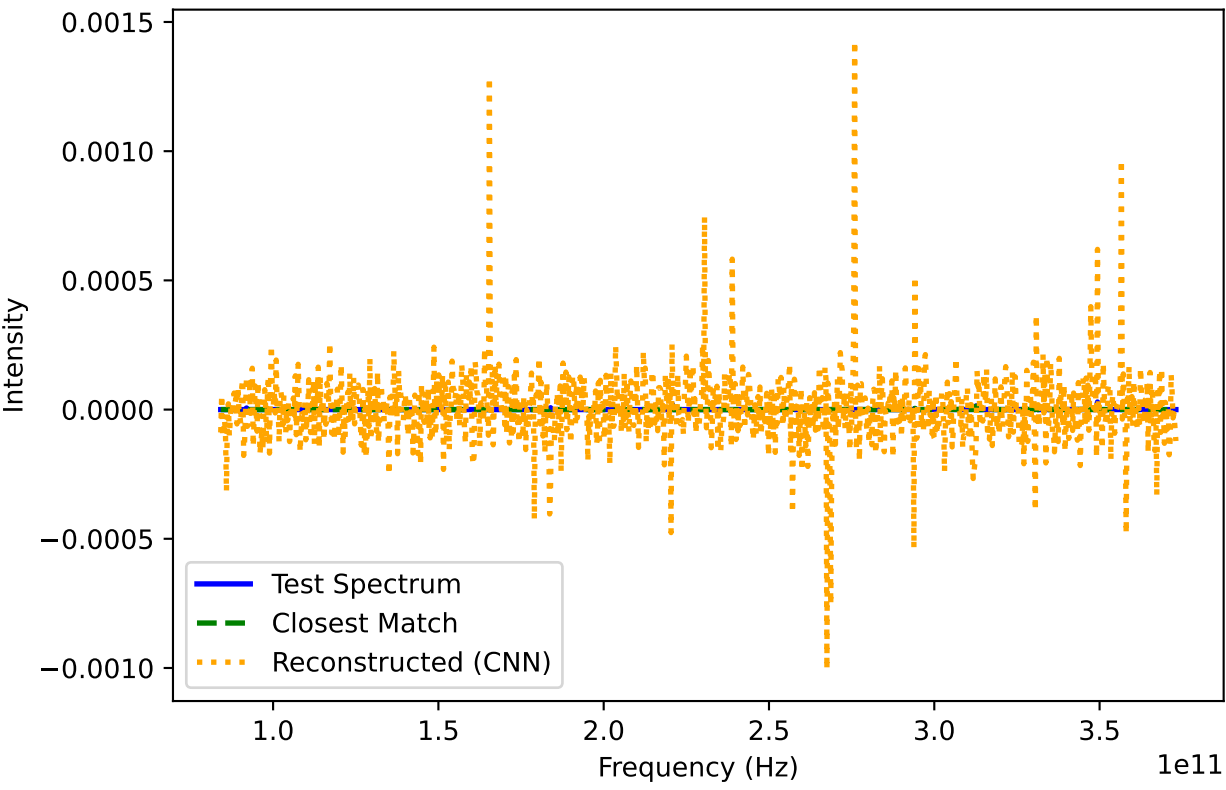
Test 8 vs Match 125



Test File: file_JPL_HCO+v=0,1,2_T21.970000000000002_N19.000000000000003_simulate_generate
Header: //molecules='HCO+V=0,1,2[1]' logn=19.0 tex=21.97 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_HCO+v=0,1,2_T21.970000000000002_N18.500000000000002_simulate_generate
Header: //molecules='HCO+V=0,1,2[1]' logn=18.5 tex=21.97 velo=250.0 fwhm=50.0 sourcesize=10.0

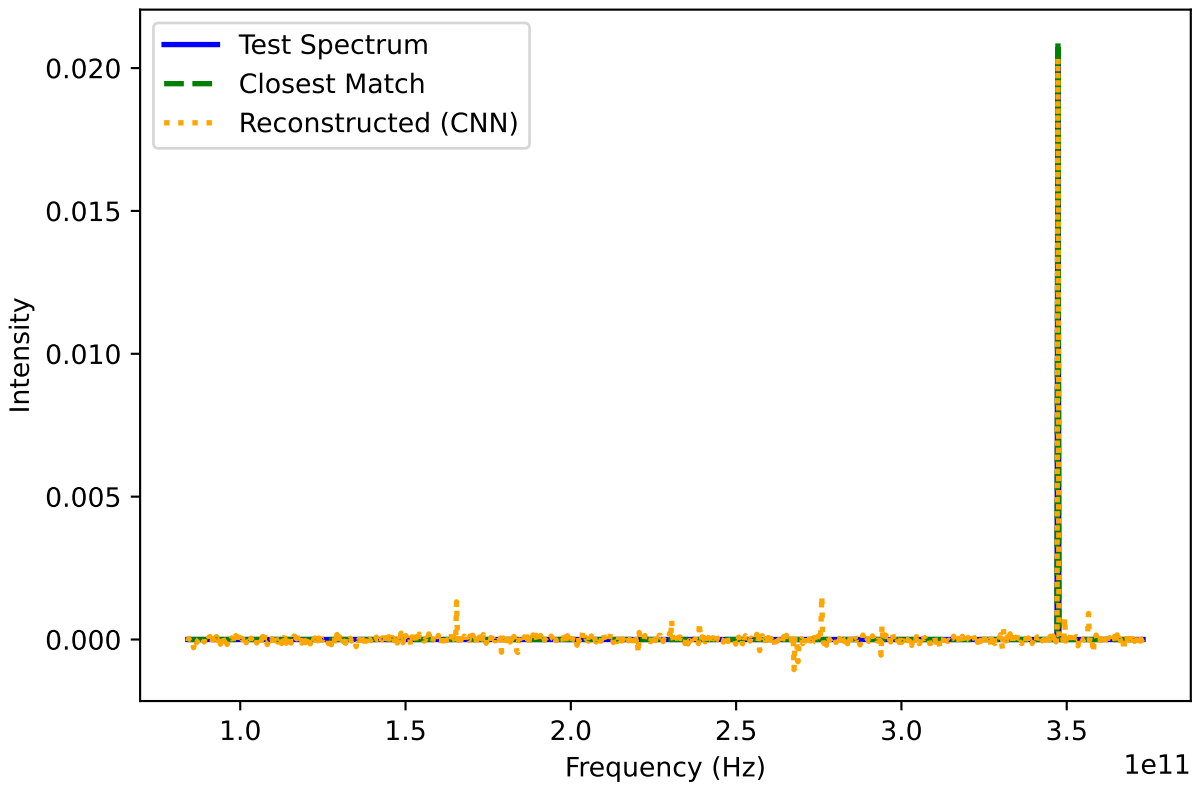
Test 9 vs Match 2269



Test File: file_JPL_CH3CN_T106.04499373000004_N12.1_simulate_generate
Header: //molecules='CH3CN[1]' logn=12.1 tex=106.04499 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_CH3CN_T137.85849184900007_N12.1_simulate_generate
Header: //molecules='CH3CN[1]' logn=12.1 tex=137.85849 velo=250.0 fwhm=50.0 sourcesize=10.0

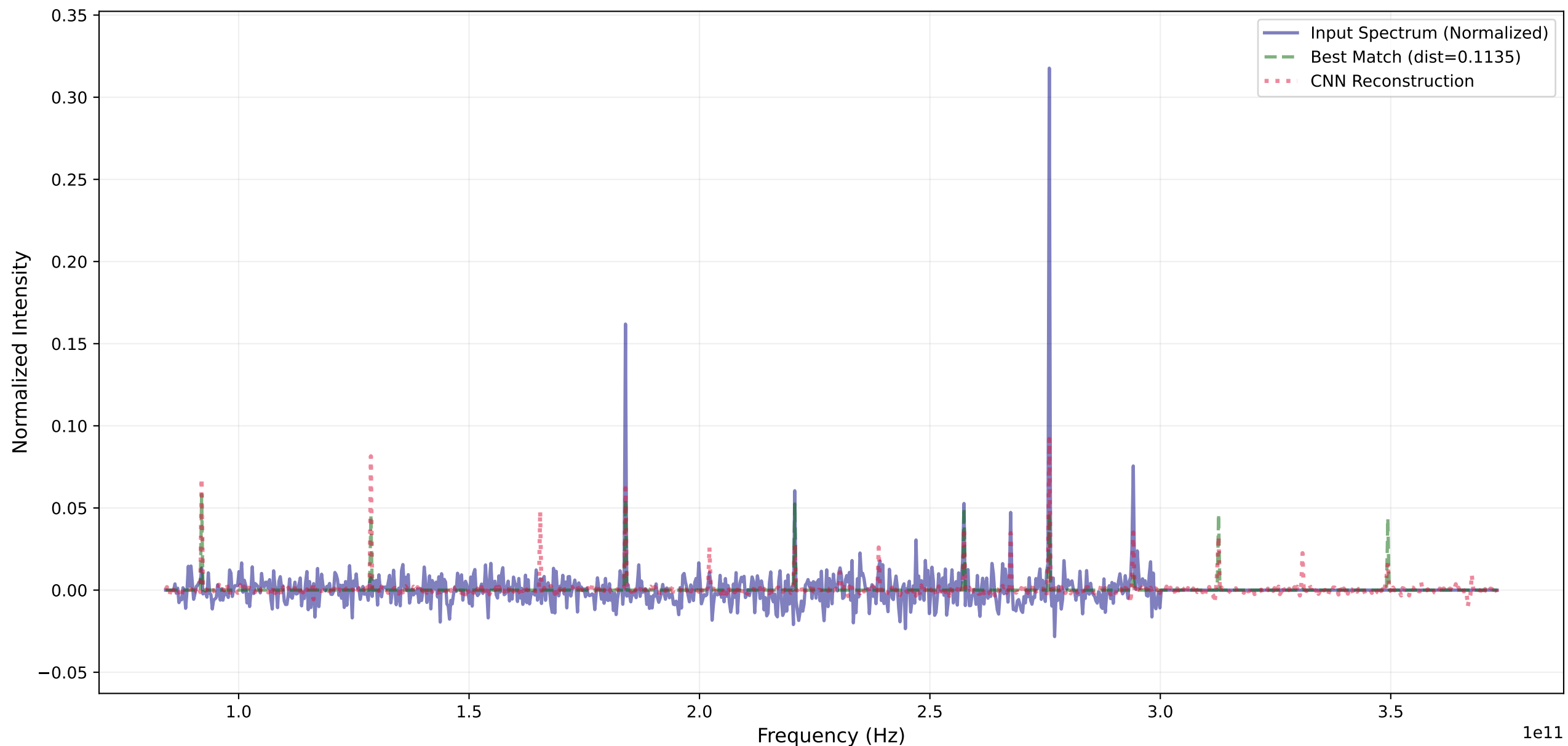
Test 10 vs Match 1243



Test File: file_JPL_SiO_T48.268090000000015_N17.600000000000001_simulate_generate
Header: //molecules='SiO[1]' logn=17.6 tex=48.26809 velo=250.0 fwhm=50.0 sourcesize=10.0

Match: file_JPL_SiO_T137.85849184900007_N17.600000000000001_simulate_generate
Header: //molecules='SiO[1]' logn=17.6 tex=137.85849 velo=250.0 fwhm=50.0 sourcesize=10.0

Input Spectrum Analysis: example_4mols_format



INPUT FILE:

=====

Filename: example_4mols_format
Header: //!xValues(GHz)[]yValues(K)

BEST MATCH:

- =====
- File: file_JPL_CH3CN_T16.9000000000000002_N19.1000000000000003_simulate_generate
 - Distance: 0.113487
 - Parameters: //molecules='CH3CN|1' logn=19.1 tex=16.9 velo=250.0 fwhm=50.0 sourcesize=10.0