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- Pr = 0.707
- T0 = 300 K
- P0 = 101325 Pa
- Uniform molar fraction of H2O = 0.02

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+++++
GENERAL LAYOUT
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&Version File_Version="VERSION2.0"/
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FLUID PROPERTIES
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====
INCOMPRESSIBLE FLUID FLOW --> Constant Density
HEAT DRIVEN FLOW          --> Activation of Heat Transfer
BOUSSINESQ ASSUMPTION     --> Thermal Expansion Coefficient = 1/T0 (
here beta = 0 ==> beta = 1/T0 )

&Fluid_Properties   Variable_Density   = .false.   ,
Constant_Mass_Flow   = .true.   , Heat_Transfer_Flow = .true.   ,
                    Heat_Capacity_Ratio = 1.4   , Reference_Density= 1.225,
Reference_Dynamic_Viscosity= 1.852D-05,
                    Reference_Temperature= 300.0   , Prandtl = 0.707,
Reference_Heat_Capacity = 1004.D0 , Thermal_Expansion_coefficient =
0.0/

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====
INITIALIZATION OF THE VELOCITY COMPONENTS, THE TEMPERATURE AND
SPECIES
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====
START FROM FLOW AT REST
AND UNIFORM TEMPERATURE at T0 = 300 K

&Velocity_Initialization   I_Velocity_Reference_Value = 0.0   ,
J_Velocity_Reference_Value = 0.0   , K_Velocity_Reference_Value = 0.0   ,
                    Initial_Field_Option_For_Velocity_I = 0,
Initial_Field_Option_For_Velocity_J = 0 ,
Initial_Field_Option_For_Velocity_K = 0/

&Temperature_Initialization   Temperature_Reference_Value = 300.0,
Initial_Field_Option_For_Temperature = 0 /

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GRAVITY

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FORCE GRAVITY ALONG THE VERTICAL AXIS POINTING DOWNWARD ( i.e. gravity
= -g.\vec{z} )
CONSIDERING DIMENSIONAL PARAMETER g = 9.81 m/s^2

&Gravity Gravity_Enabled= .true. , Gravity_Angle_IJ= 90.0 ,
Gravity_Angle_IK= 0.0 , Reference_Gravity_Constant= 9.81 /

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                        RADIATION

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AS RADIATION IS CONSIDERED :
- ACTIVATE THE RADIATIVE SOLVER [default = .false.] ( ONLY FOR 3D
CARTESIAN PROBLEMS !! )
- SOLVE THE RADIATIVE PROBLEM EVERY 5 CONVECTIVE TIMESTEP ( LIMIT TIME
CONSUMPTION , KEEP THIS PARAMETER LOWER THAN 5~8 FOR STABILITY ... )
[default = 1]
- IF STARTED FROM SCRATCH, FORCE THE SOLVER TO ITERATE OVER
FirstIterations=200 LOCAL ITERATIONS FOR INCIDENT FLUXES CONVERGENCES
AT WALLS AND VOLUMIC RADIATIVE SOURCE TERM [default = 20]
- FOR EACH RADIATIVE PROBLEM SOLVING STEPS, ITERATE OVER
RadiativeLocalIterations=20 SUB-ITERATIONS OR UNTIL
RadiativeConvergenceTolerance=5.E-05 RESIDUAL
ERROR IS REACHED [default = 1.E-15]
- WallRadCoeff AND VolRadCoeff ARE FOR DEVELOPPEMENT ONLY ... [default
= 1]
- CONSIDER THE "LATHROP" SCHEME TO INTERPOLATE THE CELL-FACES
RADIATIVE INTENSITY [default = STEP]
- CONSIDER THE ANGULAR DISCRETISATION WITH S10 LEVEL SYMMETRIC
QUADRATURES Squad = 10 ( 120 DIRECTIONS IN VOLUMES, 60 DIRECTIONS ON
WALLS) [default = 8]
- CONSIDER BLACK WALLS ON DIRICHLET WALLS AND REFLECTIVE WALLS ON THE
OTHERS [default = 0.1]
- CONSIDER THE MEDIUM AS A REAL GAS MIXTURE :
+ ACTIVATE THE SLW MODEL ActivateGas=.true. [default = .false.]
+ SPLIT THE ABSORPTION COEFFICIENT DOMAIN IN 8 WEIGHTED SUM OF GRAY-
GASES NbGas = 8 [default = 1]
+ ka_min AND ka_max REPRESENTS THE MINIMUM AND MAXIMUM RANGE OF THE
ABSORPTION COEFFICIENT DOMAIN in m^{-1} [default = 0]
+ CONSIDERS THE MEDIUM AS AN AIR-H2O GAS MIXTURE WITH UNIFORM MOLAR
FRACTION x = 0.02 [default = 0.07]

&Radiative_Heat_Transfer_DOM activateRadiation=.true. ,
RadiativePeriod = 5, FirstIterations=200,
RadiativeLocalIterations=20,
RadiativeConvergenceTolerance = 5.E-05,
WallRadcoeff = 1.0 , VolRadCoeff = 1.0, RadiativeScheme
= "LATHROP",

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        ActivateGas=.true., NbGas = 8, ka_max=570.,
ka_min=6.3e-07,
        Pref=101325.0, Href = 1.0, spec='H2O',xaref=0.02,
xaUniform=0.02,
        Squad = 10, WallEmissivity = 1.0 1.0 0.0 0.0 0.0 0.0 /

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=====
                        DOMAIN FEATURES
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=====
- CONSIDER HERE A CUBICAL CAVITY WITH WALL REFINED CELLS GIVEN IN
SEPARATE MESH FILES
- WE CONSIDERS AN MPI DOMAIN DECOMPOSITION PROBLEM ON 2x2x3 MPI
PROCESSES

&Domain_Features Start_Coordinate_I_Direction= 0.00 ,
End_Coordinate_I_Direction= 1.00,
                Start_Coordinate_J_Direction= 0.00 ,
End_Coordinate_J_Direction= 1.00,
                Start_Coordinate_K_Direction= 0.00 ,
End_Coordinate_K_Direction= 1.00,
                Cells_Number_I_Direction= 40
,Cells_Number_J_Direction= 40 ,Cells_Number_K_Direction= 30,
                Number_OMP_Threads= 1,
                MPI_Cartesian_Topology= .true. ,
                Total_Number_MPI_Processes= 12,
                Max_Number_MPI_Proc_I_Direction= 2 ,
Max_Number_MPI_Proc_J_Direction= 2, Max_Number_MPI_Proc_K_Direction= 3,
                Regular_Mesh= .false. /

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+++++
                        DEFINITION OF BOUNDARY CONDITIONS
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=====
                        WALL BOUNDARY CONDITION SETUP
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=====
- WE CONSIDER DIRICHLET TEMPERATURE CONDITION ON HOT AND COLD WALLS
(Heat_BC_Option = 0)
- AND WALL CONVECTION-RADIATION COUPLING AT THE OTHER WALLS
(Heat_BC_Option = 5)

&Heat_Wall_Boundary_Condition_Setup
        Wall_BC_DataSetName ="Set1",
        West_Heat_BC_Option = 0 , East_Heat_BC_Option = 0 ,
Back_Heat_BC_Option = 5 , Front_Heat_BC_Option = 5 ,

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South_Heat_BC_Option = 5 , North_Heat_BC_Option = 5,
      West_Wall_BC_Value= 300.005 , East_Wall_BC_Value= 299.995 ,
Back_Wall_BC_Value= 0.0 , Front_Wall_BC_Value= 0.0 ,
South_Wall_BC_Value= 0.0 , North_Wall_BC_Value= 0.0 /

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=====
                                BORDER BOUNDARY CONDITIONS
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!--- No new boundary conditions are defined at the ends of the domain
: walls by default are preserved, the inlet and outlet previously are
defined above)
!--- As "None" is the default setting for this namelist, it can be
removed

&Border_Domain_Boundary_Conditions West_BC_Name= "None" , East_BC_Name=
"None" , Back_BC_Name= "None" , Front_BC_Name= "None" , North_BC_Name=
"None" , South_BC_Name= "None" /
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+++++
                                NUMERICAL METHODS
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+++++
PARTIAL DIAGONALISATION TECHNIQUE IS EMPLOYED FOR THE POISSON PROBLEM
==> Numerical_Method_Poisson_Equation   = 3

&Numerical_Methods  NS_NumericalMethod= "BDF2-Scheme02"
,      !--- BDF2 + 2nd order centered scheme
      MomentumConvection_Scheme="Centered-02-
Conservative" ,      !--- conservative form for solving the velocity
(momentum) equation
      TemperatureAdvection_Scheme="Centered-02-
Conservative",      !--- conservative form for solving the temperature
(enthalpy) equation
      Poisson_NumericalMethod="Home-
PartialDiagonalization" /      !--- Partial Diagonalization for
Poisson's equation
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+++++
                                SIMULATION MANAGEMENT
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+++++
- START FROM SCRATCH IF Restart_Parameter= 0 OR FROM EXISTING FILES IF
Restart_Parameter= 3
- WE CONSIDERS THAT THE PROBLEM WILL REACH A STEADY STATE AND WILL
EVOLVE IN TIME WITH FIXED CFL PARAMETER

&Simulation_Management      Restart_Parameter= 3 ,
      Steady_Flow_Stopping_Criterion_Enabled =
.true. , Steady_Flow_Stopping_Criterion = 1.D-14,

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Temporal_Iterations_Number = 100 , Final_Time
= 3.D+04 ,
TimeStep_Type = 1 ,
CFL_Min      = 0.3 , CFL_Max      = 0.3 ,
Timestep_Min = 1.D-03 , Timestep_Max =
1.D+01 ,
Iterations_For_Timestep_Linear_Progress= 1,
Probe_Recording_Rate      = 1000
,
Simulation_Backup_Rate    = 5000
, Simulation_Checking_Rate = 20 /

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PROBES MANAGEMENT
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FIELDS RECORDING DECLARATION
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=====
&Field_Recording_Setup      Check_Special_Features=
"NOHeat_Driven_Cavity_Flow" /

&Simulation_Management
    InstantaneousFields_RecordingReset=.false.      ,
    InstantaneousFields_TimeRecordingRate= 5.0E+01 ,
    InstantaneousFields_RecordingStartTime= 0.D-00 /
&Instantaneous_Fields_Listing Name_of_Field = "U"      " /      First
velocity component
&Instantaneous_Fields_Listing Name_of_Field = "V"      " /      Second
velocity component
&Instantaneous_Fields_Listing Name_of_Field = "W"      " /      Third
velocity component
&Instantaneous_Fields_Listing Name_of_Field = "T"      " /
Temperature
&Instantaneous_Fields_Listing Name_of_Field = "P"      " /      Pressure
&Instantaneous_Fields_Listing Name_of_Field = "divU"   " /      Momentum
divergence

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