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Produced by the NASA Center for Aerospace Information (CASI)
An Interactive Computer Code for Calculation of Gas-Phase Chemical Equilibrium (EQLBRM)

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Seattle, Washington

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PREFACE

This report describes EQLBRM, an interactive FORTRAN IV computer program for computing adiabatic flame temperatures and equilibrium product composition for the combustion of hydrocarbon fuels in air. The authors are grateful to Paul S. Cramer of the AVCO-Lycoming Corporation for extensively testing EQLBRM and suggesting many improvements.
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1. INTRODUCTION

1.1 Overview

EQLBRM is a user-friendly, menu-driven, interactive computer program which calculates the adiabatic equilibrium temperature and product composition resulting from the combustion of hydrocarbon fuels with air, at specified constant pressure and enthalpy. EQLBRM was developed primarily as an instructional tool to be run on small computers to allow the user to economically and efficiently explore the effects of varying fuel type, air-fuel ratio, inlet air and/or fuel temperature, and operating pressure on the performance of continuous combustion devices such as gas turbine combustors, Stirling engine burners, and power-generation furnaces.

1.2 Method of Solution

EQLBRM is based on the principle of minimization of the Gibbs function of a reactive mixture at constant pressure, subject to constraints of conservation of atomic species and enthalpy. The minimization algorithm used is the very efficient and robust Gordon and McBride method (ref. 1), as refined and simplified by Pratt (refs. 2 to 4), and recently developed independently by Reynolds (ref. 5). A detailed description of the solution algorithm is contained in reference 4.

1.3 Special Features

The many "user-friendly" features offered by EQLBRM include:

1. A resident default combustion problem specification allows immediate demonstration of program execution and output format.

2. Interactive, run-time selection of SI or English units.

3. Fuel compounds may be either chosen from a resident data file, EQLBRM.DAT, or entered interactively at run time.

4. Composition of the adiabatic equilibrium combustion products is output in mole numbers, mole fractions, mass fractions and emission index.

5. At the user's option, a summary of the inlet conditions together with the equilibrium solution may be written to a file, EQLOUT.LST, for subsequent hard copy output.

1.4 Limitations

Compared to the powerful NASA equilibrium code CEC71(1), EQLBRM has the following limitations:

1. To fit in a small computer, program size is limited to 80 species, compared to over 500 species in CEC71. (The limitation to 80 species includes any new species added during an interactive session.)

2. Condensed species are not considered.
3. Only the HP problem (assigned enthalpy and pressure) is considered by
EQLBRM, while CEC71 considers also the TP, UV, TV, SP and SV problems (ref. 1).

4. Multiple problem statements are not considered.

5. EQLBRM considers only the first NPROD species in the THERMO data file
(see Chapter 3) as combustion products; NPROD is presently hard-coded as 11.
NPROD can be modified only by altering a DATA statement in the BLOCK DATA sub-
program source code.

2. USER'S GUIDE

A well-written interactive program contains the equivalent of a User's
Guide in its interactive menu structure and prompts. As a practical matter,
it is necessary to actually use an interactive program to discover how it
works. One way to resolve this "Catch-22" dilemma is to obtain telephone-line
access to a demonstration run on a computer in which the interactive program
is installed. By this method a prospective user can "take a test drive" and
see whether or not the program is of further interest.

An alternative method is the presentation of example interactive sessions.
Two such example sessions are presented here. In both examples, everything
that would appear on the user's terminal (either a printing terminal or
cathode-ray tube screen) appears in UPPER CASE text, as it does in actual pro-
gram execution. The user response to prompts is boxed, and the lower case
text are comments provided to rationalize the hypothetical user's choices, or
to highlight some of the features of EQLBRM.

2.1 Example Run 1

In this example, it is assumed that the user is running EQLBRM for the
first time, and wishes only to become familiar with its operation, rather than
solve a specific problem. The user therefore types in "RUN EQLBRM" or an
equivalent system command to execute the compiled and linked load module named
"EQLBRM."

Immediately EQLBRM introduces itself by printing to the terminal:

*** PROGRAM EQLBRM ***

CALCULATES THE ADIABATIC EQUILIBRIUM TEMPERATURE AND PRODUCT COMPOSITION
RESULTING FROM COMBUSTION OF HYDROCARBON FUELS AND AIR AT CONSTANT PRES-
SURE AND ENTHALPY.

USER MUST SPECIFY INITIAL FUEL COMPOSITION AND TEMPERATURE, INITIAL AIR
TEMPERATURE, AIR-FUEL MIXTURE RATIO, AND PRESSURE. A VARIETY OF UNITS ARE
AVAILABLE:

<table>
<thead>
<tr>
<th>TEMPERATURE</th>
<th>PRESSURE</th>
<th>AIR-FUEL RATIO</th>
</tr>
</thead>
<tbody>
<tr>
<td>KELVIN</td>
<td>KILO PASCALS</td>
<td>KG AIR/KG FUEL</td>
</tr>
<tr>
<td>CELSIUS</td>
<td>P.S.I.</td>
<td>KG FUEL/KG AIR</td>
</tr>
<tr>
<td>FAHRENHEIT</td>
<td>ATMOSPHERES</td>
<td>MOLS AIR/MOLS FUEL</td>
</tr>
<tr>
<td>RANKINE</td>
<td>BAR</td>
<td>MOLS FUEL/MOLS AIR</td>
</tr>
<tr>
<td></td>
<td></td>
<td>FUEL-AIR EQUIVALENCE RATIO</td>
</tr>
</tbody>
</table>
EQLBRM WILL RETURN THE EQUILIBRIUM TEMPERATURE AND FUEL-AIR EQUIVALENCE RATIO, AS WELL AS THE MOLE NUMBER, MOLE FRACTION, MASS FRACTION AND EMISSION INDEX OF THE PRODUCTS AT EQUILIBRIUM.

PRESS <RETURN> TO CONTINUE

Being an interactive program, EQLBRM always prompts the user for the proper response or range of responses. The user responds with:

<return>

IF THE USER SPECIFIES A FUEL FOR WHICH EQLBRM HAS NO THERMOCHEMICAL DATA, THE USER MAY EXPLICITLY SUPPLY THE FUEL CHEMICAL FORMULA AND STANDARD STATE ENTHALPY OF FORMATION OR LOWER HEATING VALUE. AGAIN, A CHOICE OF UNITS WILL BE AVAILABLE TO THE USER.

WOULD YOU LIKE TO SEE A SAMPLE DATA SET (Y/N)

Since this is just a familiarization run, the user elects to inspect the sample data set.

Y

***CURRENT DATA STATUS***

FUELS AND FUEL AMOUNTS:

<table>
<thead>
<tr>
<th>FUEL</th>
<th>RELATIVE MOLES AMONG FUELS</th>
</tr>
</thead>
<tbody>
<tr>
<td>CO</td>
<td>1.0000000</td>
</tr>
<tr>
<td>H2</td>
<td>2.0000000</td>
</tr>
</tbody>
</table>

FUEL TEMP = 2.9800E+02 DEG- KELVIN
AIR TEMP = 7.5000E+02 DEG- KELVIN
PRESSURE = 5.0000E+00 ATMOSPHERES
A-F RATIO = 8.5000E-01 FUEL-AIR EQUIVALENCE RATIO

SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

Again, since this is just a familiarization run, the user chooses to run with the sample data set, just to see what the output looks like.

0

RUN WITH THIS DATA: ARE YOU SURE? (Y/N)

EQLBRM always double-checks before execution to avoid possible time-costly user error.
Now EQLBRM executes the sample data set and returns the solution to the screen:

```
OUTPUT

*** ADIABATIC EQUILIBRIUM AT CONSTANT PRESSURE AND ENTHALPY ***

PRESSURE = 5.000E+00 ATMOSPHERES
TEMPERATURE = 2444.76 DEGREES KELVIN
FUEL-AIR EQUIVALENCE RATIO = 0.850

<table>
<thead>
<tr>
<th>SPECIES NAME \ (&quot;I&quot;)</th>
<th>MOLE NUMBER \ (MOL/-G-X)</th>
<th>MOLE FRACTION \ (MOL/-MOL-X)</th>
<th>MASS FRACTION \ (G/-G-X)</th>
<th>EMISSION INDEX \ (G/-KG-FUEL)</th>
</tr>
</thead>
<tbody>
<tr>
<td>CO</td>
<td>2.204E-04</td>
<td>6.078E-03</td>
<td>6.173E-03</td>
<td>6.173E+00</td>
</tr>
<tr>
<td>CO2</td>
<td>3.270E-03</td>
<td>9.019E-02</td>
<td>1.439E-01</td>
<td>1.439E+02</td>
</tr>
<tr>
<td>H</td>
<td>1.458E-05</td>
<td>4.020E-04</td>
<td>1.469E-05</td>
<td>1.469E+02</td>
</tr>
<tr>
<td>H2</td>
<td>7.696E-05</td>
<td>2.123E-03</td>
<td>1.551E-03</td>
<td>1.551E+01</td>
</tr>
<tr>
<td>H2O</td>
<td>6.771E-03</td>
<td>1.868E-01</td>
<td>1.220E-01</td>
<td>1.220E+02</td>
</tr>
<tr>
<td>N</td>
<td>2.282E-09</td>
<td>6.292E-08</td>
<td>3.196E-08</td>
<td>3.196E-05</td>
</tr>
<tr>
<td>NO</td>
<td>2.848E-04</td>
<td>7.854E-03</td>
<td>8.545E-03</td>
<td>8.545E+00</td>
</tr>
<tr>
<td>N2</td>
<td>2.417E-02</td>
<td>6.667E-01</td>
<td>6.772E-01</td>
<td>6.772E+02</td>
</tr>
<tr>
<td>O</td>
<td>3.161E-05</td>
<td>8.719E-04</td>
<td>5.058E-04</td>
<td>5.058E-01</td>
</tr>
<tr>
<td>OH</td>
<td>2.503E-04</td>
<td>6.903E-03</td>
<td>4.257E-03</td>
<td>4.257E+00</td>
</tr>
<tr>
<td>O2</td>
<td>1.163E-03</td>
<td>3.209E-02</td>
<td>3.723E-02</td>
<td>3.723E+01</td>
</tr>
</tbody>
</table>

DO YOU WANT A HARD COPY OF YOUR INPUT AND OUTPUT? (Y/N)

The user elects to make a hardcopy of both the input and output datasets.

```

Y

DATA AND RESULTS ARE BEING APPENDED TO FILE "EQLOUT.LST" IN YOUR DIRECTORY.

The hard copy output is directed to logical unit 2. System-specific commands must be executed at run time (or added to the source code) to assign this unit to the file "EQLOUT.LST," or directly to a printer.

DO YOU WANT ANOTHER RUN? (Y/N)

A "Y" response would return the user to the main input menu, where the current input conditions can be modified interactively, as demonstrated in the next example run. However, the user in this case chooses to end the familiarization session.

N
2.2 Example Run 2

In this example, we assume the user wishes to solve the following problem: Propane at 125° C burns with air at 375° C, at constant pressure of 4.5 atmospheres, with a slightly rich fuel-air equivalence ratio of 1.1.

The user initiates EQLBRM, and the program introduces itself:

*** PROGRAM EQLBRM ***

CALCULATES THE ADIABATIC EQUILIBRIUM TEMPERATURE AND PRODUCT COMPOSITION RESULTING FROM COMBUSTION OF HYDROCARBON FUELS AND AIR AT CONSTANT PRESSURE AND ENTHALPY.

USER MUST SPECIFY INITIAL FUEL COMPOSITION AND TEMPERATURE, INITIAL AIR TEMPERATURE, AIR-FUEL MIXTURE RATIO, AND PRESSURE. A VARIETY OF UNITS ARE AVAILABLE:

| Temperature  | Pressure   | Air-Fuel Ratio
|--------------|------------|----------------|
| Kelvin       | KILO PASCALS | KG AIR/KG FUEL
| CELSIUS      | P.S.I.     | KG FUEL/KG AIR
| FAHRENHEIT   | ATMOSHERES | MOLS AIR/MOLS FUEL
| RANKINE      | BAR        | MOLS FUEL/MOLS AIR

EQLBRM WILL RETURN THE EQUILIBRIUM TEMPERATURE AND FUEL-AIR EQUIVALENCE RATIO, AS WELL AS THE MOLE NUMBER, MOLE FRACTION, MASS FRACTION AND EMISSION INDEX OF THE PRODUCTS AT EQUILIBRIUM.

PRESS <RETURN> TO CONTINUE.

IF THE USER SPECIFIES A FUEL FOR WHICH EQLBRM HAS NO THERMOCHEMICAL DATA, THE USER MAY EXPLICITLY SUPPLY THE FUEL CHEMICAL FORMULA AND STANDARD STATE ENTHALPY OF FORMATION OR LOWER HEATING VALUE. AGAIN, A CHOICE OF UNITS WILL BE AVAILABLE TO THE USER.

WOULD YOU LIKE TO SEE A SAMPLE DATA SET? (Y/N)

Since the user has a data set in mind already, there is no interest in seeing the sample dataset.

N

***CURRENT DATA STATUS***

FUELS AND FUEL AMOUNTS:
FUEL RELATIVE MOLES AMONG FUELS
- NO FUEL SPECIFIED -
FUEL TEMP = 0.0000E-01 DEG- KELVIN
AIR TEMP = 0.0000E-01 DEG- KELVIN
PRESSURE = 0.0000E-01 ATMOSHERES
A-F RATIO = 0.0000E-01 FUEL-AIR EQUIVALENCE RATIO
SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE A/F TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

Note that the main input menu comes up blank. Data must be entered by the user interactively. The order in which this is done is not important, only it is important that no null data be left when the dataset is executed. For the sake of convenience, the user decides to begin with option 1, and work down:

1

*CURRENT FUELS DATA*

FUEL	 RELATIVE MOLES AMONG FUELS
- NO FUEL SPECIFIED -

SELECT AN OPTION:
0- NO MORE CHANGES HERE
1- CHANGE RELATIVE MOLES OF AN EXISTING FUEL
2- ADD A NEW FUEL
3- REMOVE AN EXISTING FUEL
4- LIST ALL COMPOUNDS AVAILABLE IN THE THERMO LIBRARY

OPTION? (0-4)

The user is now in the fuels sub-menu, and exercises the 2nd option - logically enough.

2

ENTER NEW FUEL NAME.

The user could enter any alphanumeric string here - but EQLBRM is looking for a chemical formula. We assume the user carelessly enters "C7H8" as propane:

C7H8

ENTER RELATIVE MOLES OF C7H8

If there is to be a mixture of fuels, it is necessary to specify the relative amount of each fuel. For the case of a single fuel, however, any nonzero value will do. The user chooses "1.0."

1.00

*CURRENT FUELS DATA*
The user notices the error that propane is C3H8, not C7H8 as entered. Option 3 must be exercised to remove the erroneous fuel selection.

3

DELETE C7H8

-- ARE YOU SURE? (Y/N)

EQLBRM double checks with the user in order to avoid time-consuming deletion errors.

Y

*CURRENT FUELS DATA*

FUEL RELATIVE MOLES AMONG FUELS
- NO FUEL SPECIFIED -

SELECT AN OPTION:
0- NO MORE CHANGES HERE
1- CHANGE RELATIVE MOLES OF AN EXISTING FUEL
2- ADD A NEW FUEL
3- REMOVE AN EXISTING FUEL
4- LIST ALL COMPOUNDS AVAILABLE IN THE THERMO LIBRARY

OPTION? (0-4)

Option 2 is used again to enter the correct fuel name.

2

ENTER NEW FUEL NAME.

C3H8

-- SORRY -- C3H8
IS NOT IN THE THERMOCHEMICAL LIBRARY. IF YOU WISH TO USE THIS COMPOUND
YOU WILL HAVE TO SPECIFY THE ENTHALPY OF FORMATION OR LOWER HEAT LUE.
DO YOU HAVE THIS INFORMATION (Y/N)
Thermochemical data for C3H8 was not included in the input data deck. We here assume that the user has in mind a reasonable figure for the LHV of typical hydrocarbon.

EQLBRM interrogates the user for the molecular composition of C3H8:

SPECIFY CHEMICAL COMPOSITION:

ENTER KG-ATOMS C PER KG-MOL C3H8

3.00

ENTER KG-ATOMS H PER KG-MOL C3H8

8.00

ENTER KG-ATOMS O PER KG-MOL C3H8

There is no oxygen in the C3H8 molecule, so a null response is required.

ENTER KG-ATOMS N PER KG-MOL C3H8

0.00

YOU MAY SPECIFY EITHER:

1- STANDARD STATE ENTHALPY OF FORMATION
2- LOWER HEATING VALUE

WHICH VALUE? (1/2)

The user recalls that a typical LHV for hydrocarbon fuels is about 18,500 Btu/lbm, and decides to use this value.

SELECT UNITS:

1- KJ/KMOL
2- KJ/KG
3- BTU/LB
4- CAL/G

WHICH UNITS? (1-4)

ENTER LOWER HEATING VALUE

18500.

ENTER RELATIVE MOLES OF C3H8

8
Again, since this is the only fuel to be included, any nonzero value will do.

1.00

*CURRENT FUELS DATA*

FUEL RELATIVE MOLES AMONG FUELS
C3H8 1.0000E+00

SELECT AN OPTION:
0- NO MORE CHANGES HERE
1- CHANGE RELATIVE MOLES OF AN EXISTING FUEL
2- ADD A NEW FUEL
3- REMOVE AN EXISTING FUEL
4- LIST ALL COMPOUNDS AVAILABLE IN THE THERMO LIBRARY

OPTION? (0-4)

No more changes are required, but the user decides to see what compounds are
in the properties library:

4

*COMPOUNDS AVAILABLE IN THERMO LIBRARY*

C0  C02  H  H2
H2O  N  NO  N2
O  OH  O2  NO2
CH4  C6H6  C6H6(L)  C7H8
C8H16  C8H18  C8H18(L)  C12H26
CH40  C2H50H  C3H80  CH2
CH20  C2H2  C2H4  C2H6
C3H8

PRESS <RETURN> TO CONTINUE.

Note that C3H8 has been added to the list of recognised compounds. It will,
however, be lost when EQLBRM is exited. The user must add thermochemical data
for C3H8 to the input deck if C3H8 is to be used repeatedly. EQLBRM can ac-
commodate up to 80 species, inclusive of those added interactively in the
manner just demonstrated.

*CURRENT FUELS DATA*

FUEL RELATIVE MOLES AMONG FUELS
C3H8 1.0000E+00

SELECT AN OPTION:
0- NO MORE CHANGES HERE
1- CHANGE RELATIVE MOLES OF AN EXISTING FUEL
2- ADD A NEW FUEL
3- REMOVE AN EXISTING FUEL
4- LIST ALL COMPOUNDS AVAILABLE IN THE THERMO LIBRARY
OPTION? (0-4)

No more changes are required.

0

***CURRENT DATA STATUS***

FUELS AND FUEL AMOUNTS:

<table>
<thead>
<tr>
<th>FUEL</th>
<th>RELATIVE MOLES AMONG FUELS</th>
</tr>
</thead>
<tbody>
<tr>
<td>C3H8</td>
<td>1.0000000</td>
</tr>
</tbody>
</table>

FUEL TEMP = 0.0000E-01 DEG- KELVIN
AIR TEMP = 0.0000E-01 DEG- KELVIN
PRESSURE = 0.0000E-01 ATMOSPHERES
A-F RATIO = 0.0000E-01 FUEL-AIR EQUIVALENCE RATIO

SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

The user wishes to work with Celsius as temperature units.

4

SELECT AN OPTION:
1- CHANGE UNITS TO KELVIN
2- CHANGE UNITS TO CELSIUS
3- CHANGE UNITS TO FAHRENHEIT
4- CHANGE UNITS TO RANKINE

OPTION? (1-4)

2

***CURRENT DATA STATUS***

FUELS AND FUEL AMOUNTS:

<table>
<thead>
<tr>
<th>FUEL</th>
<th>RELATIVE MOLES AMONG FUELS</th>
</tr>
</thead>
<tbody>
<tr>
<td>C3H8</td>
<td>1.0000000</td>
</tr>
</tbody>
</table>

FUEL TEMP = 0.0000E-01 DEG- CELSIUS
AIR TEMP = 0.0000E-01 DEG- CELSIUS
PRESSURE = 0.0000E-01 ATMOSPHERES
A-F RATIO = 0.0000E-01 FUEL-AIR EQUIVALENCE RATIO
SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

The user notes that fuel and air use the same units for temperature. Again, order is unimportant, but the user chooses to work top-down through the menu. Next in the menu is fuel temperature selection.

**CURRENT DATA STATUS**

FUELS AND FUEL AMOUNTS:
FUEL	 RELATIVE MOLES AMONG FUELS
C3H8	 1.0000000
FUEL TEMP = 1.250E+02 DEG- CELSIUS
AIR TEMP = 0.000E-01 DEG- CELSIUS
PRESSURE = 0.000E-01 ATMOSPHERES
A-F RATIO = 0.000E-01 FUEL-AIR EQUIVALENCE RATIO

SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

The desired fuel temperature is specified.

**CURRENT DATA STATUS**

FUELS AND FUEL AMOUNTS:
FUEL	 RELATIVE MOLES AMONG FUELS
C3H8	 1.0000000

ENTER TEMPERATURE OF AIR, DEG- CELSIUS

Now the desired input air temperature is specified.

**CURRENT DATA STATUS**

FUELS AND FUEL AMOUNTS:
FUEL	 RELATIVE MOLES AMONG FUELS
C3H8	 1.0000000
FUEL TEMP = 1.2500E+02 DEG- CELSIUS
AIR TEMP = 3.7500E+02 DEG- CELSIUS
PRESSURE = 0.0000E-01 ATMOSPHERES
A-F RATIO = 0.0000E-01 FUEL-AIR EQUIVALENCE RATIO

SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

Still working arbitrarily top-down, only options 5 and 6 remain to be exercised before the dataset is complete.

5

SELECT AN OPTION:
1- SAME UNITS, NEW PRESSURE
2- CHANGE UNITS TO KILO PASCALS
3- CHANGE UNITS TO P.S.I.
4- CHANGE UNITS TO ATMOSPHERES
5- CHANGE UNITS TO BAR

OPTION? (1-5)

Note that entering either a "1" or a "4" would have the same effect here.

1

ENTER PRESSURE, IN ATMOSPHERES

4.50

***CURRENT DATA STATUS***

FUELS AND FUEL AMOUNTS:

<table>
<thead>
<tr>
<th>FUEL</th>
<th>RELATIVE MOLES AMONG FUELS</th>
</tr>
</thead>
<tbody>
<tr>
<td>C3H8</td>
<td>1.0000000</td>
</tr>
</tbody>
</table>

FUEL TEMP = 1.2500E+02 DEG- CELSIUS
AIR TEMP = 3.7500E+02 DEG- CELSIUS
PRESSURE = 4.5000E+00 ATMOSPHERES
A-F RATIO = 0.0000E-01 FUEL-AIR EQUIVALENCE RATIO

SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS
12
OPTION NUMBER? (0-6)

The only null data left is the air-fuel ratio.

6

SELECT AN OPTION:
1- SAME UNITS, NEW RATIO
2- CHANGE UNITS TO KG AIR/KG FUEL
3- CHANGE UNITS TO KG FUEL/KG AIR
4- CHANGE UNITS TO MOLS AIR/MOLS FUEL
5- CHANGE UNITS TO MOLS FUEL/MOLS AIR
6- CHANGE UNITS TO FUEL-AIR EQUIVALENCE RATIO

OPTION? (1-6)

Note that either a "1" or a "6" will have the same effect here.

1

ENTER FUEL-AIR EQUIVALENCE RATIO

1.10

***CURRENT DATA STATUS***

FUELS AND FUEL AMOUNTS:
FUEL RELATIVE MOLES AMONG FUELS
C3H8 1.0000000

FUEL TEMP = 1.2500E+02 DEG- CELSIUS
AIR TEMP = 3.7500E+02 DEG- CELSIUS
PRESSURE = 4.5000E+00 ATMOSPHERES
A-F RATIO = 1.1000E+00 FUEL-AIR EQUIVALENCE RATIO

SELECT AN OPTION:
0- RUN WITH THIS DATA
1- CHANGE FUELS OR FUEL AMOUNTS
2- CHANGE FUEL TEMPERATURE
3- CHANGE AIR TEMPERATURE
4- CHANGE TEMPERATURE UNITS
5- CHANGE PRESSURE OR PRESSURE UNITS
6- CHANGE AIR-FUEL RATIO OR RATIO UNITS

OPTION NUMBER? (0-6)

The user has now specified all input conditions of the problem.

0

RUN WITH THIS DATA: ARE YOU SURE? (Y/N)

The user has double-checked all the items on the menu, so as to avoid wasting computer time on bad input data.

Y
EQLBRM executes the dataset and presents the solution:

******
OUTPUT
******

*** ADIABATIC EQUILIBRIUM AT CONSTANT PRESSURE AND ENTHALPY ***

PRESSURE = 4.500E+00 ATMOSPHERES
TEMPERATURE = 2007.09 DEGREES CELSIUS
FUEL-AIR EQUIVALENCE RATIO = 1.100

<table>
<thead>
<tr>
<th>SPECIES NAME</th>
<th>MOLE NUMBER</th>
<th>MOLE FRACTION</th>
<th>MOLE MASS FRACTION</th>
<th>EMISION INDEX</th>
</tr>
</thead>
<tbody>
<tr>
<td>CO</td>
<td>1.080E-03</td>
<td>2.993E-02</td>
<td>3.026E-02</td>
<td>3.026E+01</td>
</tr>
<tr>
<td>CO2</td>
<td>3.999E-03</td>
<td>9.444E-02</td>
<td>1.500E-01</td>
<td>1.500E+02</td>
</tr>
<tr>
<td>H</td>
<td>1.407E-05</td>
<td>3.899E-04</td>
<td>1.418E-05</td>
<td>1.418E-02</td>
</tr>
<tr>
<td>H2</td>
<td>3.274E-04</td>
<td>9.071E-03</td>
<td>6.600E-04</td>
<td>6.600E-01</td>
</tr>
<tr>
<td>H2O</td>
<td>5.632E-03</td>
<td>1.560E-01</td>
<td>1.015E-01</td>
<td>1.015E+02</td>
</tr>
<tr>
<td>N</td>
<td>4.476E-10</td>
<td>1.240E-08</td>
<td>6.269E-09</td>
<td>6.269E-06</td>
</tr>
<tr>
<td>NO</td>
<td>1.784E-05</td>
<td>4.944E-04</td>
<td>5.354E-04</td>
<td>5.354E-01</td>
</tr>
<tr>
<td>N2</td>
<td>2.556E-02</td>
<td>7.083E-01</td>
<td>7.161E-01</td>
<td>7.161E+02</td>
</tr>
<tr>
<td>O</td>
<td>1.125E-06</td>
<td>3.118E-05</td>
<td>1.800E-05</td>
<td>1.800E-02</td>
</tr>
<tr>
<td>OH</td>
<td>3.805E-05</td>
<td>1.054E-03</td>
<td>6.472E-04</td>
<td>6.472E-01</td>
</tr>
<tr>
<td>O2</td>
<td>8.201E-06</td>
<td>2.272E-04</td>
<td>2.624E-04</td>
<td>2.624E-01</td>
</tr>
</tbody>
</table>

DO YOU WANT A HARD COPY OF YOUR INPUT AND OUTPUT? (Y/N)

The user elects to make a hardcopy of both the input and output datasets.

Y

DATA AND RESULTS ARE BEING APPENDED TO FILE "EQLOUT.LST" IN YOUR DIRECTORY.

The hard copy output is directed to logical unit 2. System-specific commands must be executed at run time (or added to the source code) to assign this unit to the file "EQLOUT.LST," or directly to an output device.

2.3 Debugging

If improper operation of EQLBRM is suspected, a debugging feature can be brought up by "un-commenting" the appropriate source code statements near the top of the main routine. With this feature restored, an interactive choice of debugging level (1-5) can be selected, which can be used (together with the source code listing) to discover implementation or source code transcription errors.

2.4 FORTRAN Language and Availability

Program EQLBRM was developed on a Digital Equipment Corporation PDP 11/23 computer. The source code is in 1966 ANSI-standard FORTRAN IV, and runs
without modification (except for system-specific I/O statements in the main routine KEYBRD) on VAX and IBM computers.

Source codes and the required data file EQLBRM.DA1 are available from COSMIC, University of Georgia, Athens, Georgia 30602.

3. THERMOCHEMICAL DATA

The thermochemical data for elements and for nonelemental molecular species are taken directly from the NASA element and thermochemical data files (ref. 1).

In the EQLBRM.DA1 data file, the ELEMENTS code word and elemental data must precede the THERMO code word and thermochemical data. The format for input data are summarized in tables I and II:
### TABLE I. - ELEMENTS INPUT RECORDS

<table>
<thead>
<tr>
<th>Order</th>
<th>Contents</th>
<th>Format</th>
<th>Record Columns</th>
</tr>
</thead>
<tbody>
<tr>
<td>First</td>
<td>ELEMENTS</td>
<td>3A4</td>
<td>1 to 8</td>
</tr>
<tr>
<td>Any</td>
<td>One record each distinct element present in the chemical system. Each record contains:</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>1) Atomic symbol of element—must agree with that used in THERMO data.</td>
<td>A2</td>
<td>1 to 2</td>
</tr>
<tr>
<td></td>
<td>2) Atomic weight of the element</td>
<td>F10.6</td>
<td>10 to 19</td>
</tr>
<tr>
<td></td>
<td>3) Valence or oxidation state of the element (positive, negative or zero)</td>
<td>F10.6</td>
<td>20 to 29</td>
</tr>
<tr>
<td>Last</td>
<td>Blank record card</td>
<td>------</td>
<td>all</td>
</tr>
</tbody>
</table>

**Note A:** EQLBRM is limited to ten distinct elements.
### TABLE II. - THERMO INPUT RECORDS

<table>
<thead>
<tr>
<th>Order</th>
<th>Contents</th>
<th>Format</th>
<th>Card Columns</th>
</tr>
</thead>
<tbody>
<tr>
<td>First</td>
<td>THERMO</td>
<td>3A4</td>
<td>1 to 6</td>
</tr>
<tr>
<td>See Note B</td>
<td>Sets of four records in sequence for each species in the chemical system. The record formats for each set are, in order:</td>
<td></td>
<td></td>
</tr>
<tr>
<td>First in set</td>
<td>1) Molecular symbol or name of species 3A4</td>
<td>1 to 12</td>
<td></td>
</tr>
<tr>
<td></td>
<td>2) Date 2A3</td>
<td>19 to 24</td>
<td></td>
</tr>
<tr>
<td></td>
<td>3) Atomic symbols and formula 4(A2,F3.0)</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>4) Phase (gas only, letter G) A1</td>
<td>45</td>
<td></td>
</tr>
<tr>
<td></td>
<td>5) Temperature 2F10.3</td>
<td>46 to 65</td>
<td></td>
</tr>
<tr>
<td></td>
<td>6) Integer 1 I15</td>
<td>80</td>
<td></td>
</tr>
<tr>
<td>Second in set</td>
<td>1) Coefficients ($z_i$, $i = 1,5$) for upper temperature range. See Note A. 5E15.8</td>
<td>1 to 75</td>
<td></td>
</tr>
<tr>
<td></td>
<td>2) Integer 2 I5</td>
<td>80</td>
<td></td>
</tr>
<tr>
<td>Third in set</td>
<td>1) Coefficients $z_6$ and $z_7$ for upper temperature range, and $z_1$, $z_2$ and $z_3$ for lower. See Note A. 5E15.8</td>
<td>1 to 75</td>
<td></td>
</tr>
<tr>
<td></td>
<td>2) Integer 3 I5</td>
<td>80</td>
<td></td>
</tr>
<tr>
<td>Fourth in set</td>
<td>1) Coefficients ($z_i$, $i = 4,7$) for low temperature interval. See Note A. 4E15.8</td>
<td>1 to 60</td>
<td></td>
</tr>
<tr>
<td></td>
<td>2) Integer 4 I20</td>
<td>80</td>
<td></td>
</tr>
<tr>
<td>Last</td>
<td>Blank record</td>
<td></td>
<td></td>
</tr>
</tbody>
</table>

**Note A:** Coefficients ($z_i$, $i = 1,7$) correspond to the following equations for constant-pressure specific heat capacity, enthalpy, and entropy:

\[
\frac{C_P}{R} = z_1 + \frac{z_2}{2} + \frac{z_3}{3} + \frac{z_4}{4} + \frac{z_5}{5} + z_6
\]

\[
\frac{h}{RT} = z_1 + \frac{z_2}{2} + \frac{z_3^2}{3} + \frac{z_4^3}{4} + \frac{z_5^4}{5} + \frac{z_6}{T}
\]

\[
\frac{s^o}{R} = z_1 \log T + \frac{z_2^2}{2} + \frac{z_3^3}{3} + \frac{z_4^4}{4} + z_7
\]

The individual species enthalpy $h$ include the standard-state enthalpy of formation and the sensible enthalpy. The individual species entropy $s^o$ is for a pressure of one atmosphere.

**Note B:** The first two sets must be molecular nitrogen $N_2$ and molecular oxygen $O_2$, the only constituents of air assumed in EOLBRM. All polyatomic fuel data-sets must reside at the end of the THERMO data file.

**Note C:** EQLBRM is limited to 80 distinct atomic and/or molecular species, including those entered interactively at run time.
4. REFERENCES


