Date: March 20, 2006

OUTPUT JACOBIAN/DERIVATIVE FILES (in the AJ directory):

```
RDderivDNW_xx_lll Layer downwelling Jacobian/derivative files
RDderivUPW_xx_lll Layer upwelling Jacobian/derivative files
(includes downwelling and surface components)
LEV_RDderivDNW_xx_lll Level downwelling Jacobian/derivative files
LEV_RDderivUPW_xx_lll Level upwelling Jacobian/derivative files
(includes downwelling and surface components)
```

Files related to surface derivatives

LEV_RDderivE-R1_000	TOA radiance Jacobian wrt surface emittance/reflectance (with E	+ R = 1)
LEV_RDderivEMI1_000	TOA radiance Jacobian wrt surface emittance	
LEV_RDderivRFL1_000	TOA radiance Jacobian wrt surface reflectance	
LEV_RDderivTSF1_000	TOA radiance Jacobian wrt surface temperature	

RECORD 1.1

CXID: 80 characters of user identification (80A1)

CXID(1) is the flag which determines program initialization and termination. The actual input data stream for LBLRTM commences with the record containing a '\$' in CXID(1). Any record that are read prior to a record containing a '\$' in CXID(1) are ignored. Since LBLRTM automatically recycles and reads Record 1.1 at the end of each run (this allows the stacking of multiple runs), each subsequent run of LBLRTM must also have a '\$' in CXID(1) on Record 1.1. In order to effect a normal termination of LBLRTM, CXID(1) must be set to '%', on the final record of

RECORD 1.2 (Continued)

IMRG (Continued)

IMRG OPTIONS FOR MERGED OUTPUT

RECORD 1.2 (Continued)

IMRG (Continued)

IMRG OPTIONS FOR SEQUENTIAL OUTPUT

 $\label{thm:monochromatic} \mbox{ Monochromatic Input, calculated in current run.}$

IMRG =
space to ground ground to space tangent

MERGE PATH: space to ground ground to space tangent Results
RADIAN9 MERGE nd ground to space tangent Results

RECORD 1.2 (Continued)

The actual results produced on MFILE or NFILE are determined by both IEMIT on Record 1.2 and JEMIT on either Record 6 (scanned results) or Record 7.1 (filtered results). The following table describes the options:

RECORD 1. or (required if ICNTNM = 6)

XSELF, XFRGN, XCO2C, XO3CN, XO2CN, XN2CN, XRAYL free format

XSELF H2O self broadened continuum absorption multiplicative factor

XFRGN H20 foreign broadened continuum absorption multiplicative factor

XO3CN O3 continuum absorption multiplicative factor

XO2CN O2 continuum absorption multiplicative factor

XN2CN N2 continuum absorption multiplicative factor

XRAYL Rayleigh extinction multiplicative factor

RECORD 1.2.1 (required if IEMIT = 2; otherwise omit)

NOTE: IEMIT = 2 requires binary solar radiance file SOLAR.RAD, containing extraterrestrial source spectra in units W/(m2 cm-1).

INFLAG, IOTFLG, JULDAT

1-5, 6-10, 13-15

I5, I5, 2X, I3

INFLAG (0,1,2,3) input flag for solar radiance calculation

- = 0 input previously calculated radiance and transmittance from TAPE12 (default)
- = 1 input previously calculated optical depth from TAPE12 = 2 input previously calculated upwelling radiance and transmittance from TAPE12,

RECORD 1.3 (required if IHIRAC > 0; IAERSL > 0; IEMIT = 1; IATM = 1; or ILAS > 0; otherwise omit)

V1,	V2,	SAMPLE,	DVSET,	ALFAL0,	AVMASS,	DPTMIN,	DPTFAC,	ILNFLG,	DVOUT,	NMOL_SCAL	
1-10,	11-20,	21-30,	31-40,	41-50,	51-60,	61-70,	71-80,	85,	90-100,	105	
E10.3,	E10.3,	E10.3,	E10.3,	E10.3,	E10.3,	E10.3,	E10.3,	4X(n0 0	8X,E10.3,	3x,I2)Tj ET	

```
RECORD 1.4 (required if IEMIT = 1, or both IEMIT=2 and IOTFLG=2; otherwise omit)
```

```
TBOUND, SREMIS(1), SREMIS(2), SREMIS(3), SRREFL(1), SRREFL(2), SRREFL(3), surf_refl
 1-10,
         11-20,
                   21-30,
                             31-40,
                                       41-50,
                                                 51-60,
                                                          61-70,
                                                                   75
E10.3,
         E10.3,
                  E10.3,
                             E10.3,
                                       E10.3,
                                                E10.3,
                                                         E10.3
                                                                  4X,1A
```

for down looking case, the effect of reflected atmospheric radiance above lower boundary included for SRREFL(I) > 0. (reflected atmosphere from H1 to H2 included)

TBOUND temperature of boundary (K)

SREMIS(I) frequency dependent boundary emissivity coefficients (I = 1,2,3)

EMISSIVITY = SREMIS(1) + SREMIS(2)*V + SREMIS(3)*(V**2)

*** NOTE: Entering a value for SREMIS(1) < 0 allows for direct input of boundary emissivities from file 'EMISSIVITY'

SRREFL(I) frequency dependent boundary reflectivity coefficients (I = 1,2,3)

REFLECTIVITY = SRREFL(1) + SRREFL(2)*V + SRREFL(3)*(V**2)

*** NOTE: Entering a value for SRREFL(1) < 0 allows for direct input of boundary reflectivities from file 'REFLECTIVITY'

surf_refl specifies the surface type used in computing the reflected downward radiance

```
RECORD 1.6b, 1.6c, and 1.6d (required if IMRG = 40,41 and IEMIT=3; otherwise omit)

Note: requires separate optical depth files for each layer (use IMRG = 10 prior to this run)

PTHODL

1-55

A55

PTHODTupw

1-55

A55

PTHODTdnw

1-55

A55

PTHODL = layer optical depth file (e.g., "ODint_")

PTHODL = layer optical depth file (e.g., "ODint_")

PTHODTupw = upwelling total optical depth file (e.g., "ODtoupw_)
```

```
******************
 ***** these records applicable only if LBLATM not selected (IATM=0) ******
                    LAYER INPUT (MOLECULES ONLY)
RECORD 2.1
         IFORM, NLAYRS, NMOL, SECNTO,
                                        ZH1,
                                                   ZH2,
                                                           ZANGLE
           2
                3-5,
                     6-10, 11-20,
                                      41-48, 53-60,
                                                          66-73
                     I5, F10.2, 20X, F8.2, 4X, F8.2, 5X, F8.3
         1X,I1
               I3,
             IFORM
                       (0,1) column amount format flag
                      = 0 read PAVE(L), WKL(M,L), WBROADL(L) in F10.4, E10.3, E10.3 formats (default)
                      = 1 read PAVE(L), WKL(M,L), WBROADL(L) in E15.7 format
            NLAYRS
                       number of layers (maximum of 200)
                       value of highest molecule number used (default = 7; maximum of 35)
              NMOL
                              See Table I for molecule numbers.
            SECNTO
                       user entered scale factor for the column amount for the layers defined by NLAYRS
                              if positive, looking up
                              if negative, looking down
                       normal value = 1.0
               ZH1
                       observer altitude
               ZH2
                       end point altitude
                       mean zenith angle for path calculation (degrees)
            ZANGLE
                       NOTE: With the surf_refl = 'l' option on Record 1.4
                             ZANGLE must be set. (ZANGLE .gt.90 and .le.180 )
```

RECORD 2.1.1

PAVE(L), TAVE(L), SECNTK(L), ITYL(L), IPATH, ALTZ(L-1), PZ(L-1), TZ(L-1), ATLZ(L), PZ(L), TZ(L)

1-10, 11-20, 21-30, 31-33, 34-35, 37-43, 44-51, 52-58, 59-65, 66-73, 74-80

ITYL(L) overrides the LBLRTM internal calulation of ITYPE for the layer L.
 ITYPE controls the DV ratio of the previous layer to the current layer.
 A blank defaults to the internal calulation.
 = 0 specifies 1/1

******	****	****	****	******	***	****	*****	***	****	*****	*****	*****	****	****	***	*****
****	Note:	all	the	quantites	on	thid	Record,	are	for	information	only,	LBLRTM	will	use	the	corresponding

```
RECORD 2.2.4
```

```
(XAMNT(I,L), I=1, 7), WBROADX(L)
(8E10.3)
```

XAMNT(I,L) column densities (molecules/cm**2) or mixing ratios (ppv) for 7 cross-section molecules

WBROADX(L) column density for broadening gases (molecules/cm**2), ** information only

should be the same as WBROADL, on Record 2.1.2

NOTE If IFRMX=1, then XAMNT(I,L) and WBROADX(L) are in 8E15.7 format

RECORD 2.2.5 only if (IXMOL .GT . 7) # records depends on IXMOL

(XAMNT(I,L), I=8, IXMOL)
(8E10.3)

NOTE: If IFRMX=1 then XAMNT(I,L) in 8E15.7 format

REPEAT RECORDS 2.2.3 through 2.2.5 for the remaining layers (up to NLAYXS)

```
RE
             radius of earth (km)
             defaults for RE=0:
             a) MODEL 0,2,3,6 RE = 6371.23 km
                                RE = 6378.39 km
RE = 6356.91 km
             b)
                      1
                        4,5
             c)
  HSPACE
             altitude definition for space (default = 100 km)
             internal models defined to 120 km
  VBAR
             frequency for refractive geometry calculation
                (default: VBAR = (V1+V2) / 2) (V1,V2 from Record 1.3)
REF_LAT
             latitude of location of calculation (degrees)
            defaults for REF_LAT = 0:
            a) MODEL 0,2,3,6 REF_LAT = 45.0 degrees
b) MODEL 1 REF_LAT = 15.0
c) MODEL 4,5 REF_LAT = 60.0
```

RECORD 3.3 options

RECORD 3.3A For IBMAX = 0 (from RECORD 3.1)

AVTRAT, TDIFF1, TDIFF2, ALTD1, ALTD2

1-10, 11-20, 21-30, 31-40, 41-50

F10.3, F10.3, F10.3, F10.3, F10.3

AVTRAT maximum Voigt Qidth ratio across a layer (if zero, default = 1.5)

TDIFF1 maximum layer temperature difference at ALTD1 (if zero, default = 5 K)

TDIFF2 maximum layer temperature difference at ALTD2 (if zero, default = 8 K)

ALTD1 altitude of TDIFF1 (if zero, default = 0 Km)

ALTD2 altitude of TDIFF2 (if zero, default = 100 Km)

RECORD 3.3B For IBMAX > 0 (from RECORD 3.1)

TABLE I

USER OPTIONS FOR PRESSURE, TEMPERATURE, AND MOLECULAR DENSITY

JCHARP PRESSURE 1-6 default to value for specified model atmosphere " ",A pressure in (mb) (JCHARP) " " (atm) " " (torr) В С JCHART TEMPERATURE default to value for specified model atmosphere 1-6 (JCHART) " ",A ambient temperature in deg (K) В JCHAR(M) (1) H2O (2) CO2 (3) (8) NO (9) SO2 (10) (M): AVAILABLE 03 (4) N2O (5) CO (6) CH4 (7) MOLECULAR SPECIES NO2 (11) NH3 (12) HNO3 (13) HF OH (14) (15) HCL (16) HBR (17) HI (18) CLO (19) OCS (20) H2CO (21) HOCL (22) N2 (23) HCN (24) CH3CL (25) H2O2 (26) C2H2 (27) C2H6 (28) (29) COF2 (30) SF6 (31) H2S (32) HCOOH (33) EMPTY (34) EMPTY (35) EMPTY potential choice of units for above species: JCHAR = 1-6 - default to value for specified model atmosphere = " ",A - volume mixing ratio (ppmv) = B - number density (cm-3) - mass mixing ratio (gm/kg) = C - mass density (gm m-3) = D = E - partial pressure (mb) - dew point temp (K) *H2O only* - dew point temp (C) *H2O only* = F = G

- relative humidity (percent) *H2O only*

- available for user definition

JCHAR must be less than "J"

= H

= I

TABLE II. Structure of file FSCDXS. (Continued)

V1X Beginning wavenumber for the cross-section data (cm-1) V2X Ending wavenumber for the cross-section data (cm-1) ** for information only ** DVX Wavenumber spacing for the cross-section data (cm-1) number of temperature dependent cross-section files to be read in. '1' indicates no temperature dependence. Maximum of '6' files (i.e. 6 temperatures) is allowed. ***** no default ***** NTEMP is the format specifier IFRM = 86 - HITRAN 86 format = 90 - HITRAN 90 format ** default ** Note: the headers are different between HITRAN 86 and HITRAN 90 CFRM is the blocking specifier

```
User Defined Atmospheric Profile
----- (IPRFL = 0) -----
RECORD 3.8
        LAYX, IZORP, XTITLE
              10, 11-60
              I5 A50
         I5,
       LAYX
             number of atmospheric profile boundaries
       IZORP (0,1) flag which determines value of ZORP on Record 3.8.1
                 = 1 ZORP is a pressure in millibars
       XTITLE
                 50 character descmiption of profile
RECORD 3.8.1
   ZORP, (JCHAR(K), K = 1, 28)
  1-10, 16 through 50
  F10.3, 5X,
                 35A1
```

flag for units and input options for

boundary altitude (km) or pressure (millibars) as determined by IZORP on Record 3.8

ZORP

JCHAR(K)

************************** AEROSOLS (IAERSL = 1,7) Subroutine LOWTRN Stripped down version of LOWTRAN 7 run as a subroutine to supply LBLRTM with the attenuation due to aerosols, clouds, fogs, and rain. Further details of most of the aerosol parameters are contained in the LOWTRAN7 User Guide (AFGL-TR-88-0177), the LOWTRAN6 Report (AFGL-TR-83-0187) the LOWTRAN5 Report (AFGL-TR-80-0167) and the Millimeter Aerosol Report (AFGL-TR-79-0253) Activated by IAERSL = 1 or 7 (on record 1.2) ** Note:(ip)ICERSK7and)CALT should berpCoRD4on specific@tITLE0) model levels (ZBND or PBND from Record 3.3B) ** * * 2) MODEL is read in LBLATM 3) M1, M2, and M3 are set depending on MODEL RECORD sequence as follows IHAZE, ISEASN, IVULCN, ICSTL, ICLD, IVSA, VIS, WSS, WHH, RAINRT, GNDALT 4.1 Format (6I5,5F10.3) ***Optional RECORDS*** (if ICLD = 18,19, or 20) RECORD 4.2 CTHIK, CALT, CEXT, ISEED Format (3F10.3,I10)

RECORD 4.3

ZCVSA, ZTVSA, ZINVSA

(if IVSA = 1)

RECORD 4.1 (Continued)

ICSTL is the air mass character (1 to 10); only used with Navy maritime model (IHAZE=3) (default value = 3)

ICSTL = 1 open ocean

.

10 strong continental influence

ICLD determines the inclusion of cirrus cloud attenuation or gives a choice of five cloud models and 5 rain models $\,$

ICLD for cloud and or rain

ICLD = 0 no clouds or rain

= 1 Cumulus cloud; base .66 km; top 3.0 km

RECORD 4.1 (Continued)

Optional input records after Record 4.1 Selected by parameters ICLD, IVSA, and IHAZE on Record 4.1 $\,$

RECORD 4.2 CTHIK, CALT, CEXT, ISEED

(if ICLD = 18, 19, or 20)

FORMAT (3F10.3,I1

Input record for cirrus altitude profile

RECORD 4.4 ML, TITLE

(if IAERSL = 7)

FORMAT (I5,18A4)

Additional aerosol profile

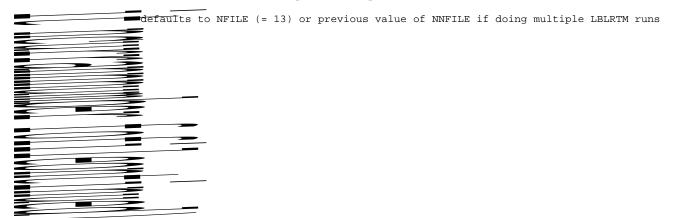
SCNMRG (Continued)

RECORD 6 (Continued)

SAMPL number of sample points per half width

- = 0 gives default value for each function
- < 0 this variable specifies the output spectral spacing (DELVO cm-1)</p>
 The value of SAMPL is calculated internally as SAMPL = HWHM/DELVO

NNFILE unit number for scanned sequential output



RECORD 7.1

V1F, DVF, NPTS, JEMIT, NNFILE, HEDDR 1-10, 11-20, 21-25, 26-30, 31-35, 46-80 F10.3, F10.4, I5, I5, I5, 10X, 8A4,A3

V1F wavenumber of initial filter value

DVF wavenumber increment between filter values

NPTS number of filter values

(if NPTS.LT.0, previous values of XF utilized XF 7tilized XF 7tilized XF7318 774 cm BT 33 0 0 -33 0

SCANFN

(ISCAN = 1)

RECORD 8.1

HWHM, V1, V2, JEMIT, JFN, JVAR, SAMPL, IUNIT, IFILST, NIFILS, JUNIT, NPTS 1-10, 11-20, 21-30, 34-35, 39-40, 44-45, 46-55, 59-60, 64-65, 69-70, 74-75, 76-80 F10.3, F10.3, F10.3, 3X,12, 3X,12, 3X,12, F10.4, 3X,12, 3X,12, 3X,12, I5

HWHM (Half Width Half Maximum)

negative value terminates SCANFN option

- Notes: 1. HWHM is first zero crossing of periodic functions for JV < 0. HWHM is redefined as $HWHM=(FIRST\ ZERO)/(PI/SCALE)$
 - 2. HWHM is instrument field of view half angle (in degrees) for JV=5,6 (e.g., for FOV of 10.0 degrees, HWHM=5.0)
- V1 beginning wavenumber value for performing SCAN
- V2 ending wavenumber value for performing SCAN
- JEMIT = -1 SCANFN convolved with absorption (1.0 transmission)
 - = 0 SCANFN convolved with transmission
 - = 1 SCANFN convolved with radiance
 - JV selects choice of scanning function

	JV	Function	Half-width Bound	Default Sample
			1 0	0 5
=	0	rectangular	1.0	0.5
=	1	triangular	2.0	2.0
=	2	gaussian	4.0	4.0
=	3	sinc squared	54.1826	4.0
=	4	sinc	119.332818	4.0
=	5	FOV correction	1.0	0.5
=	6	FOV correction	1.0	0.5

INTRPL

(ISCAN = 2)

RECORD 9.1

DVO, V1, V2, JEMIT, I4PT, IUNIT, IFILST, NIFILS, JUNIT, NPTS 1-10, 11-20, 21-30, 31-35, 36-40, 56-60, 61-65, 66-70, 71-75, 76-80 F10.3, F10.3, F10.3, I5 I5, 15X, I5, I5, I5, I5, I5

DVO wavenumber spacing for interpolated result negative value terminates INTRPL option

V1 beginning wavenumber value for performing INTRPL

V2 ending wavenumber value for performing INTRPL

JEMIT

- = -1 interpolation of absorption (1 transmission)
- = 0 interpolation of transmission
- = 1 interpolation of radiance

I4PT

- = 0 linear interpolation of data points
- = 1 four-point interpolation of data points

IUNIT unit designation of file to be interpolated (default is MFILE)

IFILST initial file from IUNIT to be interpolated

NIFILS number of files to be interpolated starting at IFILST

JUNIT file containing interpolated results (default is JUNIT, file 11)

 \mbox{NPTS} $\,$ number of values to be printed for the beginning and ending of each panel for current interpolated file

REPEAT RECORD 9.1

A '-1.' within columns 1-10 will terminate interpolation.

(ISCAN = 3)

RECORD 10.1

HWHM, V1, V2, JEMIT, JFNin = 38R.nNuVOUT, IUNIT, IFILST, NIFILS, JUNIT, IVX, NOFIX

M, (Half Width Half Maximum)

negative value terminates FFTSCN option

Note: M, $\;$ is the maximum optical path difference of an equivalent interferometer for JFNin < 0.

- V1 beginning wavenumber value for performing FFTSCN
- V2 ending wavenumber value for performing FFTSCN
- JEMIT = 0 convolve with transmittance
 - = 1 convolve with radiance

corporate result of the result

FFTSCN (Continued)

JFNin selects choice of scanning function

JF	'Nin	Scanning Function	Apodization Function	a/HWHM	a/FZ	CR
=	0	boxcar				
=	1	1-v/a, v <a 0, v >a (triangle)</a 	(sin(pi*x*a)/pi*x*a))**2 (sinc squared)	2.0	1.0	40
=	2	exp(-0.5*(v/a)**2) (gauss)	exp(-2*pi*(a*x)**2) (gauss)	0.849322	(NA)	10
=	3	<pre>(sin(pi*x*a)/pi*x*a))**2 (sinc squared)</pre>	1-x*a, x =<1/a 0, x >1/a (triangle)	2.257609	1.0	40
=	4	sin(u)/u (sinc)	1, x =<1/a 0, x >1/a (rectangle)	3.314800	2.0	160
=	5	J(5/2,u)/(u**(5/2)) (Beer)	(1-(x*a)**2)**2	2.100669	0.91728	20
=	6	<pre>sinc(u)+cl*(sinc(u+pi)+ sinc(u-pi)) (Hamming)</pre>	(1+2*c1*cos(pi*x*a))/ (1+2*c1)	2.195676	1.0	20
=	7	<pre>sinc(u)+0.5*(sinc(u+pi)+ sinc(u-pi)) (Hanning)</pre>	(1+cos(pi*x*a))/2	2.0	1.0	20

NORTON-BEER FUNCTIONS:

 $sum\{Ci*(1-(x*a)**2)**i\}$ from i=0 to 4, for
0 =< x =< 1/a
and
0 for x > 1/a

= 8 weak Ci = 0.384093,-0.087577, 2.57027 40 0.703484, 0., 0.

FFTSCN (Continued)

RECORD 10.1 (Continued)

JFNin Scanning Function Apodization Function a/HWHM a/FZ CR

OTHER FUNCTIONS:----a7m24*TR

FFTSCN (Continued)

RECORD 10.1 (Continued)

MRATin ratio of HWHM of the scanning function to the halfwidth of the boxcar; for prescanning with a

RECORD 12.1

CPRGID, CEX

1-60, 79-80

A60, 18X, A2

******	********	******	*******	
******	********	*******	*******	
******	Use Records 12.2A and 12	.3A for PLOT (IOPT = 0,1)	*********	
******	Use Record1 **** = 0,1)	**********	0 0 -33 0 437 Tm /F1.0 1 Tf(******

PLTLBL (Continued)

RECORD 12.3A (IOPT = 0,1)

```
YMIN, YMAX, YSIZE, DELY, NUMSBY, NOENDY, IDEC, JEMIT, JPLOT, LOGPLT, JHDR, JOUT, JPLTFL
1-10, 11-20, 21-30, 31-40,
                           45,
                                50, 55,
                                          60, 65,
                                                       70, 72, 77,
                                                                        80
G10.4, G10.4, F10.3, F10.3, I5, I5, I5, I5, I5, I5, I2,3X, I2,
                                                                       Ι3
```

YMIN is Y value at bottom of Y-axis.

- * IOPT = 1 will cause YMIN to determine the vertical offset (in Y)
- * for the overlayed plot. YMIN is in the units of the previous plot
- axes and any offset is applied with respect to the previously
- * defined axes. (i.e. YMIN = 10. will offset the new plot 10. units up)

YMAX is Y value at top of Y-axis.

(if log plot selected, YMIN and YMAX are the exponent values at the bottom and top of the plot).

YSIZE is the number of inches for the y-axis.

DELY is the number of y units per major division (= 1. when log plot is selected).

NUMSBY is the number of subdivisions per major division of y-axis.

NOENDY controls the plotting of the values at either end of y-axis (= 1 supresses plotting).

IDEC is number of figures after decimal point on linear y-axis.

JEMIT = 0 for transmission,

= 1 for radiance.

JPLOT = 0 plots transmission for JEMIT = 0

- = 1 plots optical depth for JEMIT = 0 = 2 plots attenuation in decibels for JEMIT = 0
- = 0 plots radiance in watts/(cm^2 sr cm^-1) for JEMIT = 1 (if YMIN and YMAX.GT.1, these values are interpreted as brightness temperatures from which minimum and maximum values of radiance are assigned).
- = 1 plots radiance in equivalent brightness temperature for JEMIT = 1

= 0 for linear y-axis, LOGPLT

= 1 for log y-axis.

= 0 for plot of header data ACHT

= 1 for suppression of header data plot

PLTLBL (Continued)

RECORD 12.3A (Continued)

JOUT = ut8 70 43 eJt to system3 eJtfile/screenTj ET Q q $0.24\ 0\ 0\ -0.24\ 18\ 774\ cm$ BT 33 $0\ 0\ -33\ 0\ 324\ cm$ /F

DITTI	(Continued)
PLILIBL	(Continued)

```
(IOPT = 2,3; CEX = 'EX')
RECORD 12.2.1B
        CFILEN(1)
          1-25
           A25
        CFILEN(1) is the external name for the data file associated with JFILE.
RECORD 12.2.2B
        CFILEN(2)
          1-25
          A25
        CFILEN(2) is the external name for the data file associated with LFILE.
RECORD 12.2.3B
        CFILEN(3)
          1-25
          A25
        CFILEN(3) is the external name for the data file associated with the difference/ratio file, MFILE.
______
REPEAT RECORD 12.2A or RECORD 12.2B
           A '-1.' within columns 1-10 will terminate plotting.
```

TABLE III. Structure of file JPLTFL (standard LBLRTM structure).

TABLE III

record 1	/PLTHDR/	<pre>XID(10),SEC,P0,T0,HMOL(64),W(64),WBROAD,D V2V,TBOUND,EMISIV,FSCDID(17),NMOL,NLAYER,</pre>	
record 2	/JPLTFL/	V1P,V2P,DVP,NLIM	
record 3		NLIM values of Y from V1P to V2P V2	V,TBOUND,EMIM

If the NLTE (IHIRAC = 4) option is selected a separate file (TAPE4) must be created. The following instructions describe the contents of a TAPE4 file.

	1 2	2143.272 4260.063	1 1
NO			
	0	0.	1
	1	1878.077	1
	2	3724.067	1

TAPE4(RECORD 1.1)

TIT

1-80