

Updated CALPUFF Modeling  
Protocol and Final Report for the  
Subject to BART Analyses  
for PCA Tomahawk Mill

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# Updated CALPUFF Modeling Protocol and Final Report for PCA “Subject to BART” Analysis

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Appendix A – Example CALMET Input File

Appendix B – Example CALPUFF Input File

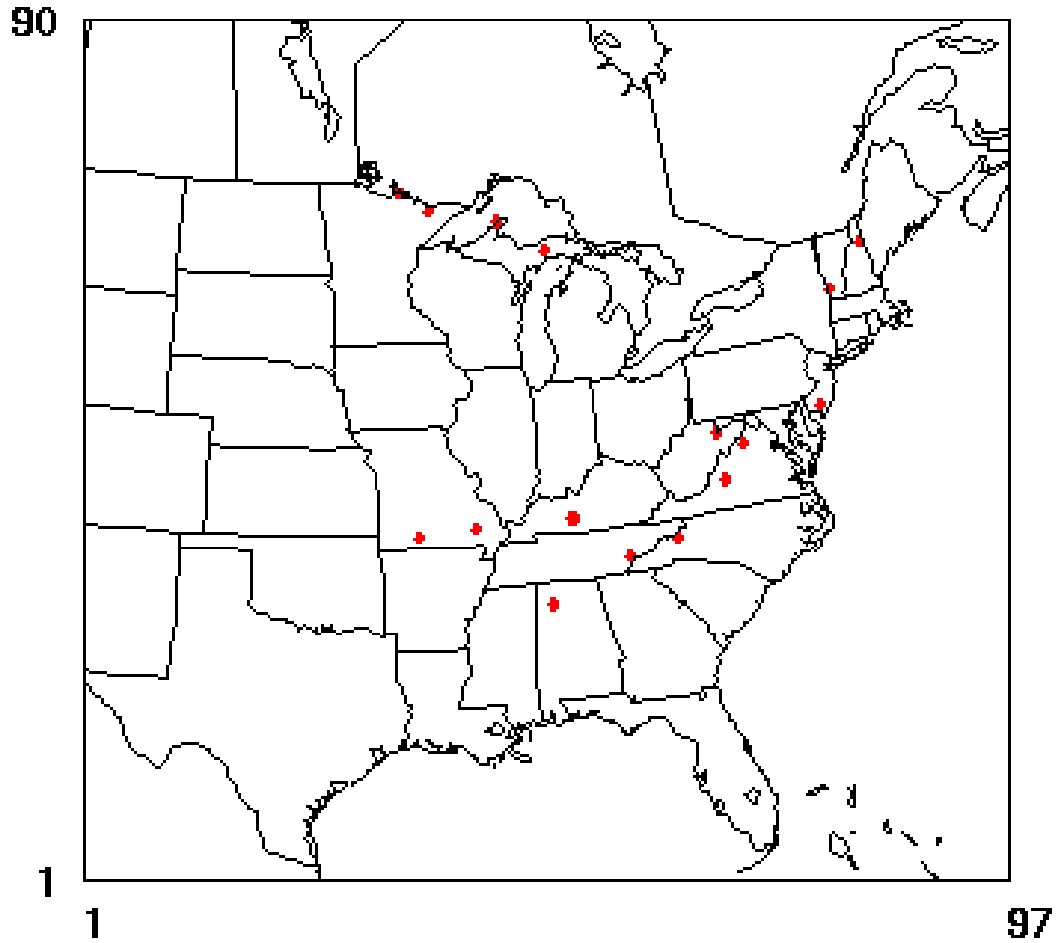
## 1.0 Introduction

The Wisconsin Department of Natural Resources (DNR) performed single source screening modeling in 2006 to evaluate which Best Available Retrofit Technology (BART) eligible sources in the state may “cause or contribute” to visibility impairment and could therefore be subject to BART requirements. This type of analysis is referred to as the “subject to BART” or “BART exemption” analysis. The methods used by DNR are described in the “Single Source Modeling to Support Regional Haze BART Modeling Protocol”, March 21, 2006, prepared by Lake Michigan Air Directors Consortium (LADCO).

DNR applied CALPUFF to each BART eligible source for three annual periods, covering the calendar years 2002 to 2004. The CALPUFF/CALMET modeling domain consisted of ninety-seven 36-km spaced cells in the east-west direction and ninety 36-km spaced cells in the north-south direction (see Figure 1). Meteorological data generated by the MM5 prognostic weather model (provided by LADCO in CALMM5 format) was used to develop the meteorological data for CALPUFF, without additional observational data input (i.e., NOOBS mode). This represents a coarse meteorological grid, suitable for an initial screening analysis.

The results from DNR’s screening analysis indicated that, with the exception of some of the largest electric power stations in the state, the only Class I areas where BART eligible sources may contribute to visibility impairment are the three areas to the north of Wisconsin, the Boundary Waters Canoe Area, Isle Royal National Park, and Seney Wildlife Area. DNR’s screening analysis also indicated that the BART eligible emissions unit at the Packaging Corporation of America Tomahawk Mill (PCA) may “cause or contribute” to visibility impairment at the Seney Class I area (located approximately 280 km to the northeast of the mill), as defined by the 98<sup>th</sup> percentile visibility impacts being greater than 0.5 deciviews (dv) relative to natural background.

Figure 1 - WDNR Screening CALPUFF Domain



NOTE: Class I areas are identified by red marks.

PCA has elected to refine the DNR screening analysis by conducting refined CALPUFF modeling with a 4 km meteorological grid, based on procedures described in the “Visibility Improvement – State and Tribal Association of the Southeast (VISTAS) BART CALPUFF” modeling protocol. The VISTAS protocol was developed by a large group of stakeholders, including EPA, the VISTAS member state agencies and tribes, the Federal Land Managers (FLMs), industry representatives, and consulting experts. Given the resources expended on the VISTAS protocol, it arguably represents the most comprehensive CALPUFF BART protocol ever developed.

PCA had discussions with DNR and EPA Region 5 staff about refining the analysis, and submitted a refined CALPUFF BART modeling protocol in March 2007. EPA reviewed the protocol and provided minor comments to DNR in May 2010. Another facility undergoing Subject to BART review, Thilmany, responded to these comments in a letter “Response to EPA Comments on Thilmany BART Modeling Protocol” dated November 9, 2010. In addition, subsequent to the EPA protocol comments, the US Fish and Wildlife Service and US Forest Service reviewed the protocol and had additional discussions with DNR and EPA. The three EPA comments, and each response, are listed below.

**Comment 1**

Information needs to be presented which clarifies how POSTUTIL was used for ammonia.

**Response 1**

The VISTAS methodology was used, with a MNITRATE switch setting of one (1).

**Comment 2**

The modeling must include a copy of the input and list files associated with runs. The protocol only offers some of the control file information.

**Response 2**

All CALMET, CALPUFF, and CALPOST input and list files, as well as meteorological and geophysical data files, had been provided by PCA to DNR on a portable hard disk drive. PCA will again provide a portable hard disk drive of all modeling files to DNR for the updated BART analysis, for purposes of SIP documentation.

**Comment 3**

The input file ... uses the background ammonia values that were produced for the LADCO protocol. These values are based on modeled estimates. Monitoring data has generally shown much higher ammonia than what was used from the model results. The refined modeling analyses ... must evaluate the visibility impacts using measured background ammonia values.

**Response 3**

The US Fish and Wildlife Service proposed using a constant background ammonia concentration of 3.9 ppb, which was the highest concentration measured at the Mayville Wisconsin station during a mid-1990s EPA study. PCA has agreed to use this single, conservative background value in the absence of daily data.

On November 29, 2010, DNR issued a letter that approved the CALPUFF BART protocol, as well as the emission rates to be used for the BART eligible emission units. DNR requested that the 2007 version of the protocol be updated to include the responses to EPA comments, and text that addresses current CALPUFF modeling guidance. In addition, the updated modeling results could be presented in the protocol (so that all information is located in one document).

This document is the updated modeling protocol and final modeling report for the refined PCA “Subject to BART” analysis. This document does not duplicate the extensive information on visibility, CALPUFF, and EPA guidance on BART modeling procedures that are contained in the VISTAS protocol. Rather, this document describes how data specific to this setting will be used in the refined CALPUFF modeling. Section 1 of the protocol is this introduction. Section 2

briefly summarizes important concepts from the VISTAS protocol. Section 3 describes the specific CALMET data and procedures that were used to develop a refined, 4-km spaced meteorological grid for Wisconsin. Section 4 describes the CALPUFF input data and procedures, and Section 5 discusses the POSTUTIL and CALPOST processing, including the use of annual average natural background visibility. Section 6 presents the emission and stack data that were used for the BART eligible emission units, and Section 7 presents the modeling results. Appendices A and B present example CALMET and CALPUFF input file listings. A portable hard drive is also being submitted that contains all input data, control files, output files, and computer codes used in the analysis.

## **2.0 VISTAS BART Protocol**

The VISTAS member states have prepared a CALPUFF modeling protocol for BART determinations that is fully consistent with the EPA guidelines in 40 CFR Part 51 Appendix W and Appendix Y. The VISTAS protocol describes a two-step CALPUFF analysis methodology for BART exemption analyses. The initial CALPUFF analysis uses a “screening” or coarse resolution meteorological grid to determine if a particular source may be exempted from further BART analyses. The screening results are also used to determine which Class I areas should be included in any refined analysis. Assumptions for the initial screening analysis are conservative so that a source that contributes to visibility impairment is not exempted in error. If a source is shown to contribute to visibility impairment using the initial screening assessment, the source has the option to undertake refined CALPUFF modeling using finer meteorological grids to evaluate further whether it is subject to BART.

VISTAS developed both coarse (12-km grid resolution) and refined (4-km grid resolution) CALMET files. The fine grid CALMET files utilized both MM5 prognostic data, as well as meteorological observational data for the CALMET “Step 2” calculations. Overwater (buoy) data were used in addition to the standard hourly surface meteorological observations, precipitation

observations, and twice-daily upper air sounding data. The VISTAS website provides detailed documentation and supporting information on the refined CALMET processing, including the CALMET and CALPUFF input files that contain model configuration options and settings that have been optimized after consultation with stakeholders and consulting experts. These same CALMET and CALPUFF input files were used as the basis for the refined Wisconsin analyses.

The VISTAS protocol discusses several options for defining the natural background visibility. These include use of the annual average natural background extinction or the background extinction for the 20% best natural conditions days. Based on an email survey of the state agencies from the VISTAS member states, all of the VISTAS states allowed the use of annual average natural background extinction for refined CALPUFF analyses. This is also the guidance provided by EPA Region 5 staff during discussions with DNR and the Wisconsin Paper Council, and the method approved by DNR in their November 29, 2010 letter. Therefore, the refined PCA CALPUFF modeling used the annual average natural background extinction for CALPOST processing.

The VISTAS BART guidance recommends that the threshold value used to define whether a source “contributes” to visibility impairment is a 0.5 dv change from natural background conditions. The 98<sup>th</sup> percentile 24-hr average predicted impact at the Class I area (equal to the 8<sup>th</sup> highest value) is to be compared to this contribution threshold value. According to clarification of the BART guidance received from EPA by the VISTAS workgroup, for a three-year simulation the modeling values to be compared with the threshold are the greatest of the three annual 8<sup>th</sup> highest values, or the 22<sup>nd</sup> highest value over all three years combined, whichever is greater.



### 3.0 CALMET Data and Procedures

EPA has updated the versions of the approved CALPUFF programs since the 2007 modeling protocol was developed, and has provided additional guidance on CALMET “switch” settings. The current approved versions are CALPUFF version 5.8 (level 070623), and CALMET version 5.8 (level 070623), and these versions were used for the updated refined BART CALPUFF analysis. On August 31, 2009, EPA published the memorandum “Clarification of EPA-FLM Recommended Settings for CALMET”. EPA stated that a 4 km fine CALMET grid size was acceptable, and provided recommended CALMET settings. The CALMET settings for the updated refined BART CALPUFF analysis meet all current EPA recommendations.

The first step in development of the refined Wisconsin analysis was to define the CALMET refined modeling domain and grid. As discussed in Section 1, DNR’s screening analysis indicated that there are three Class I areas to the north of Wisconsin, the Boundary Waters Canoe Area, Isle Royal National Park, and Seney Wildlife Area, where the screening analysis predicts some level of visibility impact from Wisconsin BART eligible sources. Therefore, the meteorological grid was developed to ensure coverage for these three Class I areas, as well as for the BART eligible source locations throughout Wisconsin. Standard IWAQM guidance also recommends that the CALMET grid extend a minimum of 50 km beyond any Class I or emission source location, so that puff recirculation can be properly accounted for.

The refined CALMET grid was defined as a 4-km spaced grid with 162 cells in the X-axis and 191 in the Y-axis. The grid projection was Lambert Conformal Conic, using the same projection parameters as in the DNR screening analysis (this allows the use of the same source X and Y coordinates as in the DNR analysis). The projection parameters include RLAT0 = 40N, RLON0 = 97W, XLAT1 = 33N, XLAT2 = 45N, XORIGKM = 247.0, YORIGKM = 240.0, and DATUM = NAS-C.

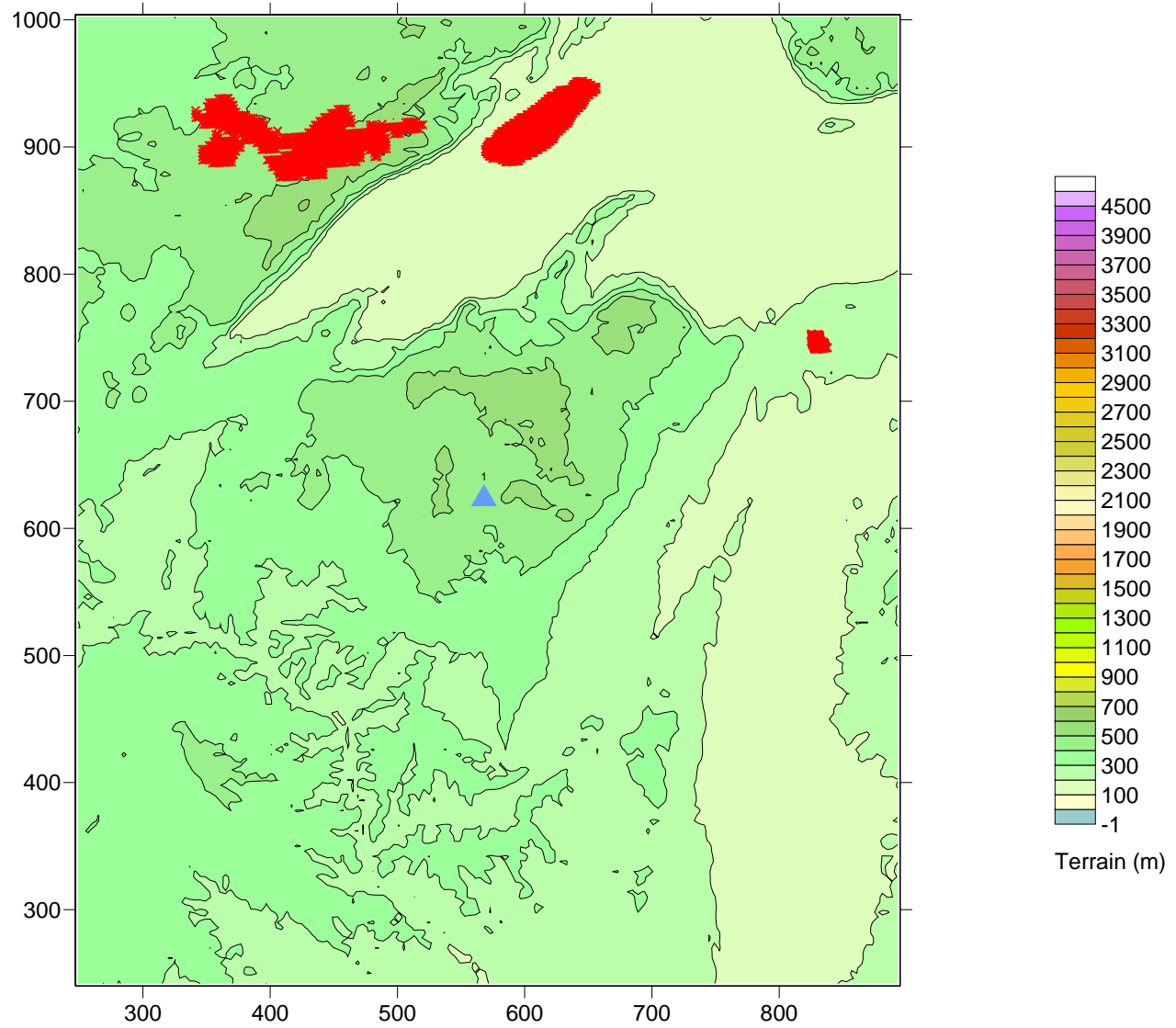
The meteorological data that was input to CALMET included the same three years (2002-2004) of CALMM5 data used by DNR, which was provided by LADCO. However, the refined CALMET runs also considered observation data from 22 surface stations, 16 precipitation stations, 5 overwater buoy stations, and 3 upper air stations. Figure 2 is a plot of the refined CALPUFF modeling domain. Figure 3 presents a plot of the locations of the observation stations in the modeling domain.

The CALMET program MAKEGEO uses USGS GTOPO30 digital terrain data (which includes coverage in Canada) and USGS Composite Theme Grid Land Use and Land Cover (LULC) data to create a gridded data file on geophysical parameters that is used by CALMET. The available USGS LULC coverage does not extend into the Canadian areas of the refined domain. These areas in Canada were assigned to the “forested land” category. Figure 4 presents a plot of the land use data for the modeling domain.

The VISTAS refined CALMET input files discussed in section 2 were used as the basis for the refined Wisconsin CALMET input files. The same CALMET configuration settings and options were utilized, as they are representative of Wisconsin terrain and land use patterns. All settings conform with current EPA and FLM guidance.

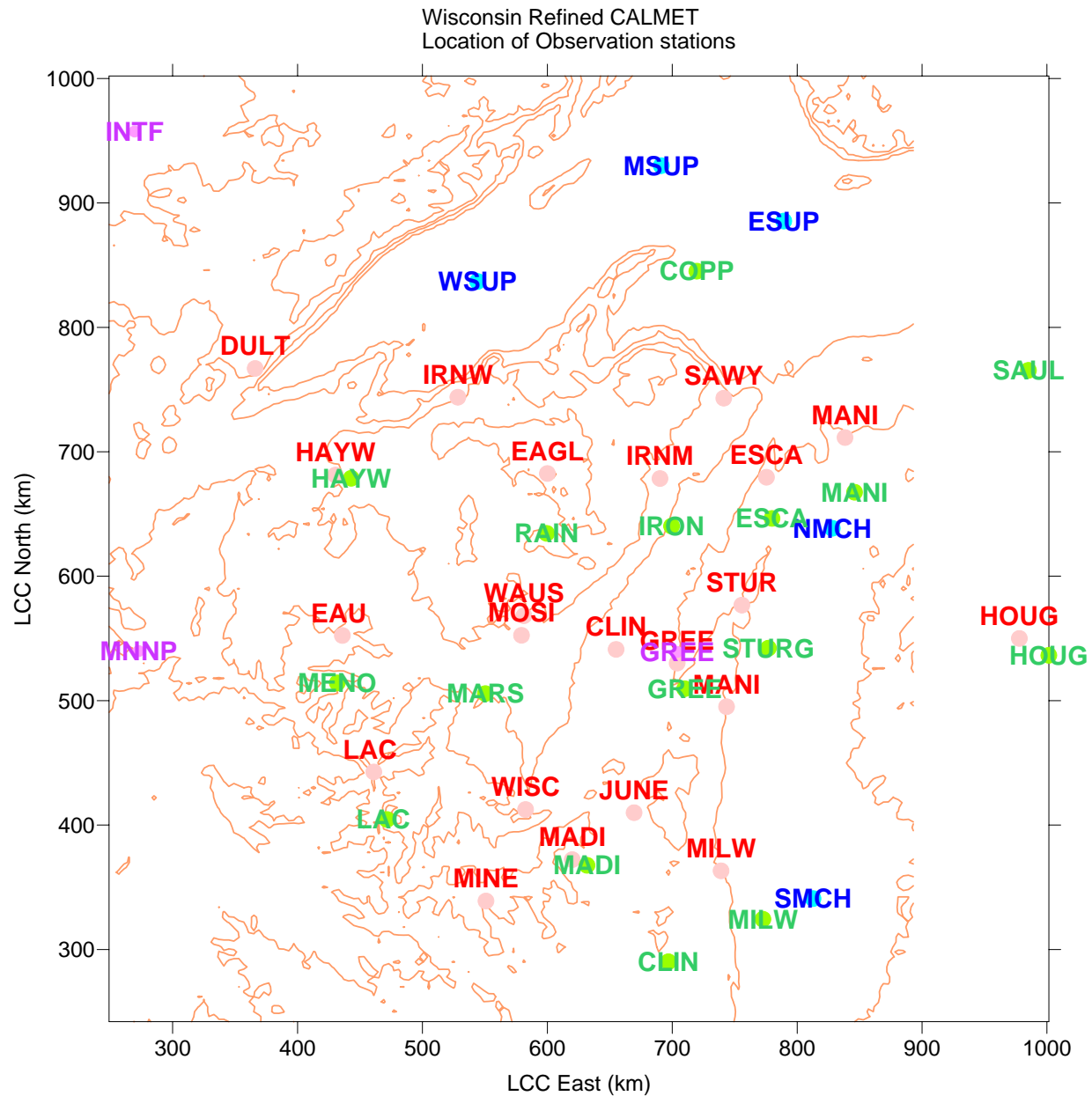
Appendix A presents a listing of a CALMET input file developed for the refined analysis.

Figure 2 – Refined CALPUFF Modeling Domain and Locations of Class I Areas



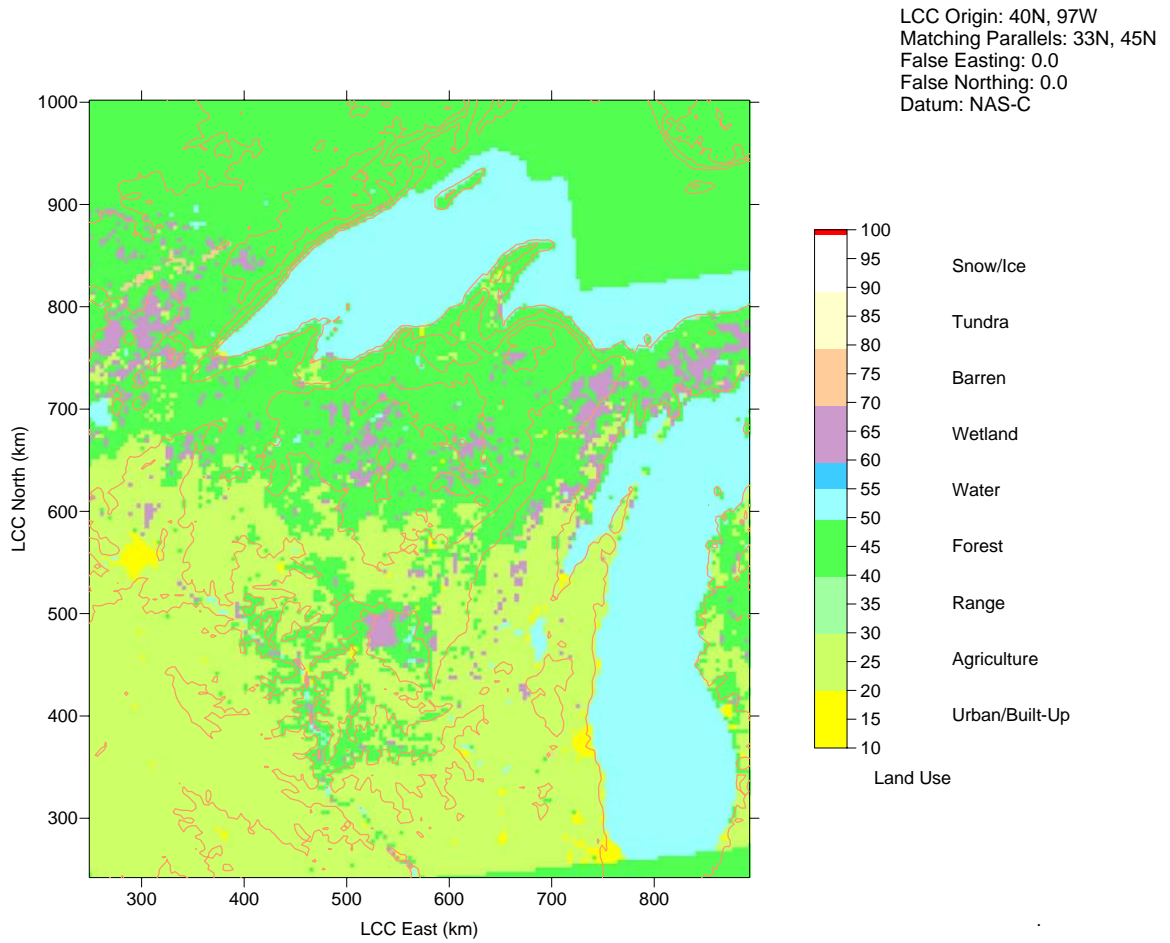
NOTE: The Class I area receptors are in red, the PCA Mill location is in blue, and the X and Y coordinates are the Lambert projection coordinates are in kms.

Figure 3 – Locations of Observational Stations in Refined CALPUFF Modeling Domain



NOTES: Surface stations are in red, overwater (buoy) in blue, precipitation in green, and upper air in purple.

Figure 4 – Land Use Data for Refined CALPUFF Modeling Domain



Some of the notable CALMET configuration settings included:

- CALMET vertical layers set to 10, with cell face heights at 20, 40, 80, 160, 320, 640, 1200, 2000, 3000, and 4000 meters,
- CALMET diagnostic options, including slope flow effects (ISLOPE = 1) with the radius of influence of terrain (TERRAD) set to 15 km,
- EPA recommended CALMET settings for the radius of influence parameters (R1=50, R2=100, RMAX1=100, RMAX2=200, and RMAX3=200).

## 4.0 CALPUFF Procedures

Both the DNR screening and VISTAS refined CALPUFF input files were used as the basis for the refined Wisconsin CALPUFF input files. All CALPUFF settings and options are consistent with EPA and FLM guidance, and the refined VISTAS protocol recommendations. The background ozone (O<sub>3</sub>) concentrations were used directly from the DNR CALPUFF input files. Instead of using monthly background ammonia concentrations from the DNR screening analyses, a single conservative background ammonia concentration of 3.9 ppb was used. Appendix B presents a listing of a CALPUFF input file developed for the analysis.

The DNR screening analysis considered three emitted pollutants, SO<sub>2</sub>, NO<sub>x</sub>, and PM<sub>2.5</sub>. The refined CALPUFF modeling will consider speciated particulate emissions, including coarse PM (PMC), fine PM equivalent to PM<sub>2.5</sub> (PMF), sulfate (SO<sub>4</sub>), nitrate (NO<sub>3</sub>), elemental carbon (EC), and secondary organic aerosols (SOA or OC). Section 6 of this protocol describes how the speciated particulate emissions were calculated.

## 5.0 POSTUTIL and CALPOST Procedures

In CALPUFF, the scavenging of the available ammonia for nitrate formation is computed on a puff-by-puff basis. Because of the possibility of multiple puff overlap, this approach may tend to overestimate the amount of ammonia available to form nitrate. In order to address the cumulative consumption of ammonia, the POSTUTIL program can be used to repartition nitric acid and nitrate. POSTUTIL calculates new NO<sub>3</sub>/HNO<sub>3</sub> equilibrium concentrations for the sample source in the Class I area using the background ammonia. As recommended in the VISTAS procedures for refined CALPUFF modeling, the POSTUTIL “MNITRATE=1” method was used for the refined Wisconsin CALPUFF modeling. Instead of using the monthly values for background ammonia concentrations from the DNR screening analyses, a single conservative background ammonia concentration of 3.9 ppb was used.

The modeling protocol submitted in 2007 proposed to use CALPOST Visibility Method 6 to calculate visibility impacts. However, since that time there have been revisions to CALPOST that allow the use of the updated IMPROVE extinction equation. CALPOST version 6.221 can implement the “Method 8 Mode 5” methodology to calculate visibility impacts, and the FLMs recommend its use in “Federal Land Managers’ Air Quality Related Values Work Group (FLAG) Phase I Report—Revised (2010)” (herein referred to as FLAG 2010). Therefore, this updated methodology will be used for the PCA BART analysis. The new CALPUFF Graphical Users Interface automatically enters the appropriate values, based on the Class I area being selected, for natural background species and relative humidity adjustment factors using the data in Tables 6 through 9 of Section 3.3.5 of FLAG 2010 (which are based on “Guidance for Estimating Natural Visibility Conditions Under the Regional Haze Rule”, EPA-454/B-03-005, September 2003).

## 6.0 Emissions Modeled

EPA’s BART modeling guidance state that the 24-hr maximum actual emissions rates for the period of the modeling simulation (2002-2004) should be used for the analysis. If this data is not available, then the short term “allowable” or “potential” emission rates should be used. PCA provided DNR with source information that has been reviewed to determine the 24-hr maximum actual emissions rates. On November 29, 2010, DNR issued a letter that approved the following emission rates to be used for the PCA BART eligible boiler: 166.3 g/sec (1320 lb/hr) SO<sub>2</sub>, 13.4 g/sec (106 lb/hr) NO<sub>x</sub>, and 4.2 g/sec (33 lb/hr) for PM. The PM emissions have been speciated using the information for coal fired boilers developed by the National Park Service<sup>1</sup>. The speciated PM emissions listed in Table 1 were used in the refined BART modeling. The stack parameters used in the modeling analysis were the same as those used by DNR in their screening analysis.

Table 1 – PCA Speciated PM Emissions

PM Species	Emission Rate (lb/hr)
Coarse PM	3.8
Fine PM	7.8
Elemental Carbon	0.3
Condensable Inorganics (Sulfate)	17.2
Condensable Organics (OC)	4.3

---

<sup>1</sup> <http://www2.nature.nps.gov/air/permits/ect/index.cfm>



## 7.0 Modeling Results

CALPOST computes the daily maximum change in deciviews relative to the annual average natural background visibility, separately for each Class I area. For evaluating the source impact relative to the “Subject to BART” visibility impact threshold of 0.5 dv, the daily maximum changes in visibility are sorted to determine the 98th percentile (8<sup>th</sup> highest) value for each of three years, and the 98th percentile (22<sup>nd</sup> highest) value for the completed three year period. If all 98th percentile values are below the threshold of 0.5 dv based on the refined CALPUFF analysis, the source is exempt from the BART requirements.

Another equivalent metric commonly used for BART exemption modeling is to sum the number of days per year that are above the 0.5 dv threshold for each Class I area separately; if there are 7 or fewer days above the threshold for each year, then the 98<sup>th</sup> percentile value is also below the threshold. This is the metric approved by DNR for use in the BART modeling analysis.

Specifically, for each Class I area separately, if there are less than 8 days per year that the predicted visibility impacts are greater than 0.5 deciviews relative to the annual average natural background visibility, then the BART eligible emission units at the modeled facility will be determined not to be Subject to BART.

Table 2 presents the results from the refined CALPUFF BART exemption modeling. These results demonstrate that there are less than 8 days per year, at each individual Class I area, when the visibility impacts from the PCA BART eligible emission unit are above 0.5 dv. Therefore, the **PCA BART eligible source is not Subject to BART** requirements.

Table 2 – PCA CALPUFF BART Exemption Results Summary

<b>Class I area</b>	<b>Distance (km) from source to Class I area boundary</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2002</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2003</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2004</b>	<b># of days with impact &gt; 0.5 dv in Class I area for 3 year period</b>	<b>Maximum 98<sup>th</sup> Percentile impact over either annual or 3-yr periods</b>
Seney , MI	~ 280	0	0	1	1	0.18
Isle Royal, MN	~ 320	0	0	0	0	0.14
Boundary Waters, MN	~ 350	0	0	0	0	0.13

**Appendix A – Example CALMET Input File**

WI Refined CALMET - 4km Grid - Dec 2010 - with Updated EPA Settings

----- Run title (3 lines) -----  
CALMET MODEL CONTROL FILE  
-----

INPUT GROUP: 0 -- Input and Output File Names

Subgroup (a)

Default Name	Type	File Name
-----	----	-----
GEO.DAT	input	! GEODAT=geo.dat !
SURF.DAT	input	! SRFDAT=SURF02Z.DAT !
CLOUD.DAT	input	* CLDDAT= *
PRECIP.DAT	input	! PRCDAT=PRECIP02.DAT !
MM4.DAT	input	* MM4DAT=E:\MRPO-calmm5\2002\MM5.020102.3D.DAT *
MM4.DAT	input	! MM4DAT=E:\MRPO-calmm5\2002\020102.MM5 !
WT.DAT	input	* WTDAT= *
CALMET.LST	output	! METLST=cal020102.lst !
CALMET.DAT	output	! METDAT=E:\WI-Calmet\2002Output\cal020102.dat !
PACOUT.DAT	output	* PACDAT= *

All file names will be converted to lower case if LCFILES = T  
Otherwise, if LCFILES = F, file names will be converted to UPPER CASE  
T = lower case ! LCFILES = T !  
F = UPPER CASE

NUMBER OF UPPER AIR & OVERWATER STATIONS:

Number of upper air stations (NUSTA) No default ! NUSTA = 3 !  
Number of overwater met stations  
(NOWSTA) No default ! NOWSTA = 5 !

!END!

-----  
Subgroup (b)

---

-----  
Upper air files (one per station)  
-----

Default Name	Type	File Name
UP1.DAT	input	1 ! UPDAT=UPGB0204.DAT! !END!
UP2.DAT	input	2 ! UPDAT=UPMN0204.DAT! !END!
UP3.DAT	input	3 ! UPDAT=UPIN0204.DAT! !END!

-----  
Subgroup (c)  
-----

Overwater station files (one per station)

SEA1.DAT	input	1 ! SEADAT=4500102.DAT! !END!
SEA2.DAT	input	2 ! SEADAT=4500202.DAT! !END!
SEA3.DAT	input	3 ! SEADAT=4500402.DAT! !END!
SEA4.DAT	input	4 ! SEADAT=4500602.DAT! !END!
SEA5.DAT	input	5 ! SEADAT=4500702.DAT! !END!

-----  
Default Name Type File Name  
-----  
-----

-----  
Subgroup (d)  
-----

Other file names  
-----

Default Name	Type	File Name
DIAG.DAT	input	* DIADAT= *
PROG.DAT	input	* PRGDAT= *
TEST.PRT	output	* TSTPRT= *
TEST.OUT	output	* TSTOUT= *
TEST.KIN	output	* TSTKIN= *
TEST.FRD	output	* TSTFRD= *
TEST.SLP	output	* TSTSLP= *

-----  
NOTES: (1) File/path names can be up to 70 characters in length  
(2) Subgroups (a) and (d) must have ONE 'END' (surround by  
delimiters) at the end of the group  
(3) Subgroups (b) and (c) must have an 'END' (surround by

delimiters) at the end of EACH LINE

!END!

-----  
INPUT GROUP: 1 -- General run control parameters  
-----

Starting date: Year (IBYR) -- No default ! IBYR= 2002 !  
Month (IBMO) -- No default ! IBMO= 1 !  
Day (IBDY) -- No default ! IBDY= 2 !  
Hour (IBHR) -- No default ! IBHR= 0 !

Base time zone (IBTZ) -- No default ! IBTZ= 6 !  
PST = 08, MST = 07  
CST = 06, EST = 05

Length of run (hours) (IRLG) -- No default ! IRLG= 24 !

Run type (IRTYPE) -- Default: 1 ! IRTYPE= 1 !

0 = Computes wind fields only  
1 = Computes wind fields and micrometeorological variables  
(u\*, w\*, L, zi, etc.)  
(IRTYPE must be 1 to run CALPUFF or CALGRID)

Compute special data fields required  
by CALGRID (i.e., 3-D fields of W wind  
components and temperature)  
in addition to regular Default: T ! LCALGRD = T !  
fields ? (LCALGRD)  
(LCALGRD must be T to run CALGRID)

Flag to stop run after  
SETUP phase (ITEST) Default: 2 ! ITEST= 2 !  
(Used to allow checking  
of the model inputs, files, etc.)  
ITEST = 1 - STOPS program after SETUP phase  
ITEST = 2 - Continues with execution of  
COMPUTATIONAL phase after SETUP

Test options specified to see if  
they conform to regulatory

values? (MREG)                      No Default                      ! MREG = 1                      !

0 = NO checks are made

1 = Technical options must conform to USEPA guidance

          IMIXH    -1            Maul-Carson convective mixing height  
                                  over land; OCD mixing height overwater  
          ICOARE    0            OCD deltaT method for overwater fluxes  
          THRESHL  0.0           Threshold buoyancy flux over land needed  
                                  to sustain convective mixing height grow

!END!

-----

INPUT GROUP: 2 -- Map Projection and Grid control parameters

-----

Projection for all (X,Y):

-----

Map projection

(PMAP)                              Default: UTM                      ! PMAP = LCC                      !

UTM : Universal Transverse Mercator  
TTM : Tangential Transverse Mercator  
LCC : Lambert Conformal Conic  
PS : Polar Stereographic  
EM : Equatorial Mercator  
LAZA : Lambert Azimuthal Equal Area

False Easting and Northing (km) at the projection origin

(Used only if PMAP= TTM, LCC, or LAZA)

(FEAST)                              Default=0.0                      ! FEAST = 0.000                      !  
(FNORTH)                             Default=0.0                      ! FNORTH = 0.000                      !

UTM zone (1 to 60)

(Used only if PMAP=UTM)

(IUTMZN)                             No Default                      ! IUTMZN = -999                      !

Hemisphere for UTM projection?

(Used only if PMAP=UTM)

(UTMHEM)                             Default: N                      ! UTMHEM = N                      !

N : Northern hemisphere projection  
S : Southern hemisphere projection

Latitude and Longitude (decimal degrees) of projection origin

(Used only if PMAP= TTM, LCC, PS, EM, or LAZA)

(RLAT0)                      No Default              ! RLAT0 = 40N    !

(RLON0)                      No Default              ! RLON0 = 97W    !

TTM : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience

LCC : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience

PS : RLON0 identifies central (grid N/S) meridian of projection  
      RLAT0 selected for convenience

EM : RLON0 identifies central meridian of projection  
      RLAT0 is REPLACED by 0.0N (Equator)

LAZA: RLON0 identifies longitude of tangent-point of mapping plane  
      RLAT0 identifies latitude of tangent-point of mapping plane

Matching parallel(s) of latitude (decimal degrees) for projection

(Used only if PMAP= LCC or PS)

(XLAT1)                      No Default              ! XLAT1 = 33N    !

(XLAT2)                      No Default              ! XLAT2 = 45N    !

LCC : Projection cone slices through Earth's surface at XLAT1 and XLAT2

PS : Projection plane slices through Earth at XLAT1  
      (XLAT2 is not used)

-----

Note: Latitudes and longitudes should be positive, and include a  
      letter N,S,E, or W indicating north or south latitude, and  
      east or west longitude. For example,  
      35.9 N Latitude = 35.9N  
      118.7 E Longitude = 118.7E

Datum-region

-----

The Datum-Region for the coordinates is identified by a character  
string. Many mapping products currently available use the model of the  
Earth known as the World Geodetic System 1984 (WGS-84). Other local  
models may be in use, and their selection in CALMET will make its output  
consistent with local mapping products. The list of Datum-Regions with  
official transformation parameters is provided by the National Imagery and  
Mapping Agency (NIMA).



NIMA Datum - Regions(Examples)

```
-----  
WGS-84   WGS-84 Reference Ellipsoid and Geoid, Global coverage (WGS84)  
NAS-C    NORTH AMERICAN 1927 Clarke 1866 Spheroid, MEAN FOR CONUS (NAD27)  
NAR-C    NORTH AMERICAN 1983 GRS 80 Spheroid, MEAN FOR CONUS (NAD83)  
NWS-84   NWS 6370KM Radius, Sphere  
ESR-S    ESRI REFERENCE 6371KM Radius, Sphere
```

Datum-region for output coordinates

```
(DATUM)           Default: WGS-84      ! DATUM = NAS-C  !
```

Horizontal grid definition:

```
-----  
  
Rectangular grid defined for projection PMAP,  
with X the Easting and Y the Northing coordinate
```

```
      No. X grid cells (NX)      No default      ! NX = 162 !  
      No. Y grid cells (NY)      No default      ! NY = 191 !
```

```
Grid spacing (DGRIDKM)          No default      ! DGRIDKM = 4. !  
                                Units: km
```

Reference grid coordinate of  
SOUTHWEST corner of grid cell (1,1)

```
      X coordinate (XORIGKM)      No default      ! XORIGKM = 247.000 !  
      Y coordinate (YORIGKM)      No default      ! YORIGKM = 240.000 !  
                                Units: km
```

Vertical grid definition:

```
-----  
  
      No. of vertical layers (NZ)  No default      ! NZ = 10  !
```

```
      Cell face heights in arbitrary  
      vertical grid (ZFACE(NZ+1))  No defaults  
                                Units: m
```

```
! ZFACE = 0.,20.,40.,80.,160.,320.,640.,1200.,2000.,3000.,4000. !
```

!END!

-----  
INPUT GROUP: 3 -- Output Options  
-----

DISK OUTPUT OPTION

Save met. fields in an unformatted  
output file ? (LSAVE) Default: T ! LSAVE = T !  
(F = Do not save, T = Save)

Type of unformatted output file:  
(IFORMO) Default: 1 ! IFORMO = 1 !

- 1 = CALPUFF/CALGRID type file (CALMET.DAT)
- 2 = MESOPUFF-II type file (PACOUT.DAT)

LINE PRINTER OUTPUT OPTIONS:

Print met. fields ? (LPRINT) Default: F ! LPRINT = T !  
(F = Do not print, T = Print)  
(NOTE: parameters below control which  
met. variables are printed)

Print interval  
(IPRINF) in hours Default: 1 ! IPRINF = 1 !  
(Meteorological fields are printed  
every 1 hours)

Specify which layers of U, V wind component  
to print (IUVOU(NZ)) -- NOTE: NZ values must be entered  
(0=Do not print, 1=Print)  
(used only if LPRINT=T) Defaults: NZ\*0  
! IUVOU = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !  
-----

Specify which levels of the W wind component to print  
(NOTE: W defined at TOP cell face -- 16 values)  
(IWOUT(NZ)) -- NOTE: NZ values must be entered

---

(0=Do not print, 1=Print)

(used only if LPRINT=T & LCALGRD=T)

-----

Defaults: NZ\*0

! IWOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Specify which levels of the 3-D temperature field to print

(ITOUT(NZ)) -- NOTE: NZ values must be entered

(0=Do not print, 1=Print)

(used only if LPRINT=T & LCALGRD=T)

-----

Defaults: NZ\*0

! ITOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Specify which meteorological fields

to print

(used only if LPRINT=T)

Defaults: 0 (all variables)

-----

Variable

Print ?

(0 = do not print,

1 = print)

-----

! STABILITY	=	0	! - PGT stability class
! USTAR	=	0	! - Friction velocity
! MONIN	=	0	! - Monin-Obukhov length
! MIXHT	=	0	! - Mixing height
! WSTAR	=	0	! - Convective velocity scale
! PRECIP	=	0	! - Precipitation rate
! SENSHEAT	=	0	! - Sensible heat flux
! CONVZI	=	0	! - Convective mixing ht.

Testing and debug print options for micrometeorological module

Print input meteorological data and

internal variables (LDB) Default: F ! LDB = F !

(F = Do not print, T = print)

(NOTE: this option produces large amounts of output)

First time step for which debug data

are printed (NN1)                      Default: 1                ! NN1 = 1 !

Last time step for which debug data  
are printed (NN2)                      Default: 1                ! NN2 = 2 !

Testing and debug print options for wind field module  
(all of the following print options control output to  
wind field module's output files: TEST.PRT, TEST.OUT,  
TEST.KIN, TEST.FRD, and TEST.SLP)

Control variable for writing the test/debug  
wind fields to disk files (IOUTD)  
(0=Do not write, 1=write)              Default: 0                ! IOUTD = 0 !

Number of levels, starting at the surface,  
to print (NZPRN2)                      Default: 1                ! NZPRN2 = 1 !

Print the INTERPOLATED wind components ?  
(IPR0) (0=no, 1=yes)                      Default: 0                ! IPR0 = 0 !

Print the TERRAIN ADJUSTED surface wind  
components ?  
(IPR1) (0=no, 1=yes)                      Default: 0                ! IPR1 = 0 !

Print the SMOOTHED wind components and  
the INITIAL DIVERGENCE fields ?  
(IPR2) (0=no, 1=yes)                      Default: 0                ! IPR2 = 0 !

Print the FINAL wind speed and direction  
fields ?  
(IPR3) (0=no, 1=yes)                      Default: 0                ! IPR3 = 0 !

Print the FINAL DIVERGENCE fields ?  
(IPR4) (0=no, 1=yes)                      Default: 0                ! IPR4 = 0 !

Print the winds after KINEMATIC effects  
are added ?  
(IPR5) (0=no, 1=yes)                      Default: 0                ! IPR5 = 0 !

Print the winds after the FROUDE NUMBER  
adjustment is made ?  
(IPR6) (0=no, 1=yes)                      Default: 0                ! IPR6 = 0 !

Print the winds after SLOPE FLOWS  
are added ?

(IPR7) (0=no, 1=yes)                      Default: 0              ! IPR7 = 0 !

Print the FINAL wind field components ?

(IPR8) (0=no, 1=yes)                      Default: 0              ! IPR8 = 0 !

!END!

-----  
INPUT GROUP: 4 -- Meteorological data options  
-----

NO OBSERVATION MODE                      (NOOBS) Default: 0              ! NOOBS = 0 !

0 = Use surface, overwater, and upper air stations

1 = Use surface and overwater stations (no upper air observations)  
    Use MM5 for upper air data

2 = No surface, overwater, or upper air observations

    Use MM5 for surface, overwater, and upper air data

NUMBER OF SURFACE & PRECIP. METEOROLOGICAL STATIONS

Number of surface stations    (NSSTA) No default              ! NSSTA = 22 !

Number of precipitation stations

(NPSTA=-1: flag for use of MM5 precip data)

(NPSTA) No default              ! NPSTA = 16 !

CLOUD DATA OPTIONS

Gridded cloud fields:

(ICLOUD) Default: 0              ! ICLOUD = 0 !

ICLOUD = 0 - Gridded clouds not used

ICLOUD = 1 - Gridded CLOUD.DAT generated as OUTPUT

ICLOUD = 2 - Gridded CLOUD.DAT read as INPUT

ICLOUD = 3 - Gridded cloud cover from Prognostic Rel. Humidity

FILE FORMATS

Surface meteorological data file format

(IFORMS) Default: 2              ! IFORMS = 2 !

(1 = unformatted (e.g., SMERGE output))

(2 = formatted    (free-formatted user input))

Precipitation data file format

(IFORMP) Default: 2 ! IFORMP = 2 !

(1 = unformatted (e.g., PMERGE output))

(2 = formatted (free-formatted user input))

Cloud data file format

(IFORMC) Default: 2 ! IFORMC = 2 !

(1 = unformatted - CALMET unformatted output)

(2 = formatted - free-formatted CALMET output or user input)

!END!

-----  
INPUT GROUP: 5 -- Wind Field Options and Parameters  
-----

WIND FIELD MODEL OPTIONS

Model selection variable (IWFCOD) Default: 1 ! IWFCOD = 1 !

0 = Objective analysis only

1 = Diagnostic wind module

Compute Froude number adjustment

effects ? (IFRADJ) Default: 1 ! IFRADJ = 1 !

(0 = NO, 1 = YES)

Compute kinematic effects ? (IKINE) Default: 0 ! IKINE = 0 !

(0 = NO, 1 = YES)

Use O'Brien procedure for adjustment

of the vertical velocity ? (IOBR) Default: 0 ! IOBR = 0 !

(0 = NO, 1 = YES)

Compute slope flow effects ? (ISLOPE) Default: 1 ! ISLOPE = 1 !

(0 = NO, 1 = YES)

Extrapolate surface wind observations

to upper layers ? (IEXTRP) Default: -4 ! IEXTRP = -4 !

(1 = no extrapolation is done,

2 = power law extrapolation used,

3 = user input multiplicative factors

for layers 2 - NZ used (see FEXTRP array)  
4 = similarity theory used  
-1, -2, -3, -4 = same as above except layer 1 data  
at upper air stations are ignored

Extrapolate surface winds even

if calm? (ICALM) Default: 0 ! ICALM = 0 !  
(0 = NO, 1 = YES)

Layer-dependent biases modifying the weights of  
surface and upper air stations (BIAS(NZ))

-1<=BIAS<=1

Negative BIAS reduces the weight of upper air stations

(e.g. BIAS=-0.1 reduces the weight of upper air stations  
by 10%; BIAS= -1, reduces their weight by 100 %)

Positive BIAS reduces the weight of surface stations

(e.g. BIAS= 0.2 reduces the weight of surface stations  
by 20%; BIAS=1 reduces their weight by 100%)

Zero BIAS leaves weights unchanged (1/R\*\*2 interpolation)

Default: NZ\*0

! BIAS = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Minimum distance from nearest upper air station  
to surface station for which extrapolation  
of surface winds at surface station will be allowed  
(RMIN2: Set to -1 for IEXTRP = 4 or other situations  
where all surface stations should be extrapolated)

Default: 4. ! RMIN2 = -1.0 !

Use gridded prognostic wind field model  
output fields as input to the diagnostic

wind field model (IPROG) Default: 0 ! IPROG = 14 !  
(0 = No, [IWFCOD = 0 or 1])

1 = Yes, use CSUMM prog. winds as Step 1 field, [IWFCOD = 0]

2 = Yes, use CSUMM prog. winds as initial guess field [IWFCOD = 1]

3 = Yes, use winds from MM4.DAT file as Step 1 field [IWFCOD = 0]

4 = Yes, use winds from MM4.DAT file as initial guess field [IWFCOD = 1]

5 = Yes, use winds from MM4.DAT file as observations [IWFCOD = 1]

13 = Yes, use winds from MM5.DAT file as Step 1 field [IWFCOD = 0]

14 = Yes, use winds from MM5.DAT file as initial guess field [IWFCOD = 1]

15 = Yes, use winds from MM5.DAT file as observations [IWFCOD = 1]

Timestep (hours) of the prognostic

model input data (ISTEPPG) Default: 1 ! ISTEPPG = 1 !

RADIUS OF INFLUENCE PARAMETERS

<p>Use varying radius of influence (if no stations are found within RMAX1,RMAX2, or RMAX3, then the closest station will be used)</p>	<p>Default: F</p>	<p>! LVARY = F!</p>
<p>Maximum radius of influence over land in the surface layer (RMAX1)</p>	<p>No default</p>	<p>! RMAX1 = 100. !</p>
	<p>Units: km</p>	
<p>Maximum radius of influence over land aloft (RMAX2)</p>	<p>No default</p>	<p>! RMAX2 = 200. !</p>
	<p>Units: km</p>	
<p>Maximum radius of influence over water (RMAX3)</p>	<p>No default</p>	<p>! RMAX3 = 200. !</p>
	<p>Units: km</p>	

OTHER WIND FIELD INPUT PARAMETERS

<p>Minimum radius of influence used in the wind field interpolation (RMIN)</p>	<p>Default: 0.1</p>	<p>! RMIN = 0.1 !</p>
	<p>Units: km</p>	
<p>Radius of influence of terrain features (TERRAD)</p>	<p>No default</p>	<p>! TERRAD = 15. !</p>
	<p>Units: km</p>	
<p>Relative weighting of the first guess field and observations in the SURFACE layer (R1) (R1 is the distance from an observational station at which the observation and first guess field are equally weighted)</p>	<p>No default</p>	<p>! R1 = 50. !</p>
	<p>Units: km</p>	
<p>Relative weighting of the first guess field and observations in the layers ALOFT (R2) (R2 is applied in the upper layers in the same manner as R1 is used in the surface layer).</p>	<p>No default</p>	<p>! R2 =100. !</p>
	<p>Units: km</p>	
<p>Relative weighting parameter of the prognostic wind field data (RPROG)</p>	<p>No default</p>	<p>! RPROG = 0. !</p>





X coordinate of BEGINNING  
of each barrier (XBBAR(NBAR))       ! XBBAR = 0. !  
Y coordinate of BEGINNING  
of each barrier (YBBAR(NBAR))       ! YBBAR = 0. !  
  
X coordinate of ENDING  
of each barrier (XEBAR(NBAR))       ! XEBAR = 0. !  
Y coordinate of ENDING  
of each barrier (YEBAR(NBAR))       ! YEBAR = 0. !

DIAGNOSTIC MODULE DATA INPUT OPTIONS

Surface temperature (IDIOPT1)       Default: 0       ! IDIOPT1 = 0 !  
0 = Compute internally from  
    hourly surface observations  
1 = Read preprocessed values from  
    a data file (DIAG.DAT)

Surface met. station to use for  
the surface temperature (ISURFT)   No default       ! ISURFT = 1 !  
(Must be a value from 1 to NSSTA)  
(Used only if IDIOPT1 = 0)

-----

Domain-averaged temperature lapse  
rate (IDIOPT2)                      Default: 0       ! IDIOPT2 = 0 !  
0 = Compute internally from  
    twice-daily upper air observations  
1 = Read hourly preprocessed values  
    from a data file (DIAG.DAT)

Upper air station to use for  
the domain-scale lapse rate (IUPT) No default       ! IUPT = 1 !  
(Must be a value from 1 to NUSTA)  
(Used only if IDIOPT2 = 0)

-----

Depth through which the domain-scale  
lapse rate is computed (ZUPT)       Default: 200.   ! ZUPT = 200. !  
(Used only if IDIOPT2 = 0)       Units: meters

-----

Domain-averaged wind components

(IDIOPT3) Default: 0 ! IDIOPT3 = 0 !

0 = Compute internally from  
twice-daily upper air observations  
1 = Read hourly preprocessed values  
a data file (DIAG.DAT)

Upper air station to use for

the domain-scale winds (IUPWND) Default: -1 ! IUPWND = -1 !

(Must be a value from -1 to NUSTA)

(Used only if IDIOPT3 = 0)

-----

Bottom and top of layer through

which the domain-scale winds

are computed

(ZUPWND(1), ZUPWND(2)) Defaults: 1., 1000. ! ZUPWND= 1., 1000. !

(Used only if IDIOPT3 = 0) Units: meters

-----

Observed surface wind components

for wind field module (IDIOPT4) Default: 0 ! IDIOPT4 = 0 !

0 = Read WS, WD from a surface  
data file (SURF.DAT)

1 = Read hourly preprocessed U, V from  
a data file (DIAG.DAT)

Observed upper air wind components

for wind field module (IDIOPT5) Default: 0 ! IDIOPT5 = 0 !

0 = Read WS, WD from an upper  
air data file (UP1.DAT, UP2.DAT, etc.)

1 = Read hourly preprocessed U, V from  
a data file (DIAG.DAT)

LAKE BREEZE INFORMATION

Use Lake Breeze Module (LLBREZE)

Default: F ! LLBREZE = F !

Number of lake breeze regions (NBOX) ! NBOX = 0 !

X Grid line 1 defining the region of interest

! XG1 = 0. !

X Grid line 2 defining the region of interest

```

! XG2 = 0. !
Y Grid line 1 defining the region of interest
! YG1 = 0. !
Y Grid line 2 defining the region of interest
! YG2 = 0. !

X Point defining the coastline (Straight line)
(XBCST) (KM) Default: none ! XBCST = 0. !

Y Point defining the coastline (Straight line)
(YBCST) (KM) Default: none ! YBCST = 0. !

X Point defining the coastline (Straight line)
(XECST) (KM) Default: none ! XECST = 0. !

Y Point defining the coastline (Straight line)
(YECST) (KM) Default: none ! YECST = 0. !

Number of stations in the region Default: none ! NLB = 0 !
(Surface stations + upper air stations)

Station ID's in the region (METBXID(NLB))
(Surface stations first, then upper air stations)
! METBXID = 0 !

!END!

-----

INPUT GROUP: 6 -- Mixing Height, Temperature and Precipitation Parameters
-----

EMPIRICAL MIXING HEIGHT CONSTANTS

Neutral, mechanical equation
(CONSTB) Default: 1.41 ! CONSTB = 1.41 !
Convective mixing ht. equation
(CONSTE) Default: 0.15 ! CONSTE = 0.15 !
Stable mixing ht. equation
(CONSTN) Default: 2400. ! CONSTN = 2400.!
Overwater mixing ht. equation
(CONSTW) Default: 0.16 ! CONSTW = 0.16 !
```

---

Absolute value of Coriolis  
parameter (FCORIOI)                      Default: 1.E-4   ! FCORIOI = 1.0E-04!  
Units: (1/s)

SPATIAL AVERAGING OF MIXING HEIGHTS

Conduct spatial averaging  
(IAVEZI) (0=no, 1=yes)                    Default: 1       ! IAVEZI = 1 !

Max. search radius in averaging  
process (MNMDAV)                          Default: 1       ! MNMDAV = 1 !  
Units: Grid  
      cells

Half-angle of upwind looking cone  
for averaging (HAFANG)                    Default: 30.     ! HAFANG = 30. !  
Units: deg.

Layer of winds used in upwind  
averaging (ILEVZI)                        Default: 1       ! ILEVZI = 1 !  
(must be between 1 and NZ)

CONVECTIVE MIXING HEIGHT OPTIONS:

Method to compute the convective  
mixing height(IMIXH)                      Default: 1       ! IMIXH = -1 !  
  1: Maul-Carson for land and water cells  
 -1: Maul-Carson for land cells only -  
      OCD mixing height overwater  
  2: Batchvarova and Gryning for land and water cells  
 -2: Batchvarova and Gryning for land cells only  
      OCD mixing height overwater

Threshold buoyancy flux required to  
sustain convective mixing height growth  
overland (THRESHL)                        Default: 0.05    ! THRESHL = 0.00 !  
(expressed as a heat flux                units: W/m3  
per meter of boundary layer)

Threshold buoyancy flux required to  
sustain convective mixing height growth  
overwater (THRESHW)                       Default: 0.05    ! THRESHW = 0.05 !  
(expressed as a heat flux                units: W/m3  
per meter of boundary layer)

Option for overwater lapse rates used  
in convective mixing height growth

```
(ITWPROG)                               Default: 0      ! ITWPROG = 0  !
0 : use SEA.DAT lapse rates and deltaT (or assume neutral
   conditions if missing)
1 : use prognostic lapse rates (only if IPROG>2)
   and SEA.DAT deltaT (or neutral if missing)
2 : use prognostic lapse rates and prognostic delta T
   (only if iprog>12 and 3D.DAT version# 2.0 or higher)
```

Land Use category ocean in 3D.DAT datasets

```
(ILUOC3D)                               Default: 16    ! ILUOC3D = 16  !
Note: if 3D.DAT from MM5 version 3.0, iluoc3d = 16
      if MM4.DAT,           typically iluoc3d = 7
```

OTHER MIXING HEIGHT VARIABLES

```
Minimum potential temperature lapse
rate in the stable layer above the
current convective mixing ht.           Default: 0.001 ! DPTMIN = 0.001 !
(DPTMIN)                               Units: deg. K/m
```

```
Depth of layer above current conv.
mixing height through which lapse
rate is computed (DZZI)                 Default: 200.  ! DZZI = 200.  !
Units: meters
```

```
Minimum overland mixing height           Default: 50.   ! ZIMIN = 50.   !
(ZIMIN)                               Units: meters
```

```
Maximum overland mixing height           Default: 3000. ! ZIMAX = 3000. !
(ZIMAX)                               Units: meters
```

```
Minimum overwater mixing height           Default: 50.   ! ZIMINW = 50.   !
(ZIMINW) -- (Not used if observed
overwater mixing hts. are used)          Units: meters
```

```
Maximum overwater mixing height           Default: 3000. ! ZIMAXW = 3000. !
(ZIMAXW) -- (Not used if observed
overwater mixing hts. are used)          Units: meters
```

OVERWATER SURFACE FLUXES METHOD and PARAMETERS

```
(ICOARE)                               Default: 10    ! ICOARE = 0  !
0 : original deltaT method (OCD)
10: COARE with no wave parameterization (jwave=0, Charnock)
11: COARE with wave option jwave=1 (Oost et al.)
    and default wave properties
-11: COARE with wave option jwave=1 (Oost et al.)
    and observed wave properties (must be in SEA.DAT files)
12: COARE with wave option 2 (Taylor and Yelland)
    and default wave properties
```

-12: COARE with wave option 2 (Taylor and Yelland)  
and observed wave properties (must be in SEA.DAT files)

Coastal/Shallow water length scale (DSHELF)  
(for modified z0 in shallow water)  
( COARE fluxes only)

Default : 0. ! DSHELF = 0. !  
units: km

COARE warm layer computation (IWARM) ! IWARM = 0 !  
1: on - 0: off (must be off if SST measured with  
IR radiometer) Default: 0

COARE cool skin layer computation (ICOOL) ! ICOOL = 0 !  
1: on - 0: off (must be off if SST measured with  
IR radiometer) Default: 0

#### TEMPERATURE PARAMETERS

3D temperature from observations or  
from prognostic data? (ITPROG) Default:0 !ITPROG = 0 !

0 = Use Surface and upper air stations  
(only if NOOBS = 0)  
1 = Use Surface stations (no upper air observations)  
Use MM5 for upper air data  
(only if NOOBS = 0,1)  
2 = No surface or upper air observations  
Use MM5 for surface and upper air data  
(only if NOOBS = 0,1,2)

Interpolation type  
(1 = 1/R ; 2 = 1/R\*\*2) Default:1 ! IRAD = 1 !

Radius of influence for temperature  
interpolation (TRADKM) Default: 500. ! TRADKM = 500. !  
Units: km

Maximum Number of stations to include  
in temperature interpolation (NUMTS) Default: 5 ! NUMTS = 5 !

Conduct spatial averaging of temp-  
eratures (IAVET) (0=no, 1=yes) Default: 1 ! IAVET = 1 !

(will use mixing ht MNMDAV,HAFANG  
so make sure they are correct)

Default temperature gradient            Default: -.0098 ! TGDEFB = -0.0098 !  
below the mixing height over  
water (K/m) (TGDEFB)

Default temperature gradient            Default: -.0045 ! TGDEFA = -0.0045 !  
above the mixing height over  
water (K/m) (TGDEFA)

Beginning (JWAT1) and ending (JWAT2)  
land use categories for temperature                            ! JWAT1 = 55 !  
interpolation over water -- Make                                ! JWAT2 = 55 !  
bigger than largest land use to disable

PRECIP INTERPOLATION PARAMETERS

Method of interpolation (NFLAGP)            Default = 2        ! NFLAGP = 2    !  
(1=1/R,2=1/R\*\*2,3=EXP/R\*\*2)  
Radius of Influence (km) (SIGMAP)        Default = 100.0   ! SIGMAP = 100. !  
(0.0 => use half dist. btwn  
nearest stns w & w/out  
precip when NFLAGP = 3)  
Minimum Precip. Rate Cutoff (mm/hr)    Default = 0.01   ! CUTP = 0.01 !  
(values < CUTP = 0.0 mm/hr)

!END!

-----  
INPUT GROUP: 7 -- Surface meteorological station parameters  
-----

SURFACE STATION VARIABLES

(One record per station -- 0 records in all)

1	2					
Name	ID	X coord.	Y coord.	Time	Anem.	
		(km)	(km)	zone	Ht. (m)	
! SS1	= 'MILW'	26400	739.061	363.200	6	10 !
! SS2	= 'MADI'	26410	619.927	372.210	6	10 !



**Refined BART Modeling Protocol for  
PCA Tomahawk Mill**

---

```

! SS3  ='LA C'   26430   461.012   442.705   6   10  !
! SS4  ='EAU '   26435   435.857   552.352   6   10  !
! SS5  ='GREE'   26450   703.911   530.179   6   10  !
! SS6  ='MANI'   26455   743.477   495.113   6   10  !
! SS7  ='STUR'   26458   755.730   576.760   6   10  !
! SS8  ='WAUS'   26463   581.202   568.247   6   10  !
! SS9  ='MOSI'   26465   579.307   552.489   6   10  !
! SS10 ='CLIN'   26502   655.019   541.168   6   10  !
! SS11 ='WISC'   26503   582.532   412.498   6   10  !
! SS12 ='EAGL'   26504   599.980   682.545   6   10  !
! SS13 ='MINE'   26507   550.817   338.941   6   10  !
! SS14 ='HAYW'   26508   430.347   681.295   6   10  !
! SS15 ='JUNE'   26509   669.429   409.931   6   10  !
! SS16 ='dulu'   27450   366.011   766.977   6   10  !
! SS17 ='MANI'   25408   838.315   711.433   5   10  !
! SS18 ='SAWY'   26284   741.237   743.078   5   10  !
! SS19 ='HOUG'   26380   977.905   549.977   5   10  !
! SS20 ='ESCA'   26480   775.417   679.550   5   10  !
! SS21 ='IRNM'   27437   690.168   678.511   5   10  !
! SS22 ='IRNW'   27445   528.314   743.857   5   10  !

```

```

-----
1
    Four character string for station name
    (MUST START IN COLUMN 9)

```

```

2
    Five digit integer for station ID

```

!END!

```

-----
INPUT GROUP: 8 -- Upper air meteorological station parameters
-----

```

```

UPPER AIR STATION VARIABLES
(One record per station -- 0 records in all)

```

```

      1      2
      Name   ID      X coord.  Y coord.  Time zone
              (km)    (km)
-----
! US1  ='GREE'  14898   703.911   539.179   6   !

```

```
! US2  ='MNNP'   94983   272.681   539.843   6   !
! US3  ='INTF'   14918   269.489   957.416   6   !
```

-----

```
1
  Four character string for station name
  (MUST START IN COLUMN 9)
```

```
2
  Five digit integer for station ID
```

!END!

-----

INPUT GROUP: 9 -- Precipitation station parameters

-----

```
PRECIPITATION STATION VARIABLES
(One record per station -- 0 records in all)
(NOT INCLUDED IF NPSTA = 0)
```

	1	2			
	Name	Station	X coord.	Y coord.	
		Code	(km)	(km)	
! PS1	'COPP'	201780	719.681	845.179	!
! PS2	'ESCA'	202626	779.850	646.454	!
! PS3	'HOUG'	203936	1001.500	536.427	!
! PS4	'IRON'	204090	699.421	640.283	!
! PS5	'MANI'	205073	845.933	667.499	!
! PS6	'SAUL'	207366	985.172	765.448	!
! PS7	'CLIN'	471667	696.889	290.151	!
! PS8	'GREE'	473269	709.933	509.579	!
! PS9	'HAYW'	473511	442.947	678.728	!
! PS10	'LAC '	474370	471.515	404.495	!
! PS11	'MADI'	474961	631.719	367.656	!
! PS12	'MARS'	475120	550.348	505.662	!
! PS13	'MENO'	475335	431.807	514.277	!
! PS14	'MILW'	475479	772.567	324.344	!
! PS15	'RAIN'	476939	599.369	634.460	!
! PS16	'STURG'	478267	776.964	542.174	!

-----  
1

Four character string for station name  
(MUST START IN COLUMN 9)

2

Six digit station code composed of state  
code (first 2 digits) and station ID (last  
4 digits)

!END!

**Appendix B – Example CALPUFF Input File**

Example BART Modeling  
2010 Updated BART Modeling - for PCA, WI.  
Using Refined 4km CALMET Data File and SENE, ISLE, BOWA Class I Receptors

----- Run title (3 lines) -----

CALPUFF MODEL CONTROL FILE  
-----

INPUT GROUP: 0 -- Input and Output File Names

Default Name	Type	File Name	
CALMET.DAT	input	* METDAT =	*
or			
ISCMET.DAT	input	* ISCDAT =	*
or			
PLMMET.DAT	input	* PLMDAT =	*
or			
PROFILE.DAT	input	* PRFDAT =	*
SURFACE.DAT	input	* SFCDAT =	*
RESTARTB.DAT	input	* RSTARTB=	*
-----			
CALPUFF.LST	output	! PUFLST =PCARef02.lst	!
CONC.DAT	output	! CONDAT =PCARef02.dat	!
DFLX.DAT	output	* DFDAT =	*
WFLX.DAT	output	* WFDAT =	*
VISB.DAT	output	* VISDAT =PCA02DNR.VIB	*
RESTARTE.DAT	output	* RSTARTE=	*

Emission Files

PTEMARB.DAT	input	* PTDAT =	*
VOLEMARB.DAT	input	* VOLDAT =	*
BAEMARB.DAT	input	* ARDAT =	*
LNEMARB.DAT	input	* LNDAT =	*

Other Files

OZONE.DAT	input	* OZDAT =	*
VD.DAT	input	* VDDAT =	*
CHEM.DAT	input	* CHEMDAT=	*
H2O2.DAT	input	* H2O2DAT=	*
HILL.DAT	input	* HILDAT=	*
HILLRCT.DAT	input	* RCTDAT=	*
COASTLN.DAT	input	* CSTDAT=	*
FLUXBDY.DAT	input	* BDYDAT=	*
BCON.DAT	input	* BCNDAT=	*
DEBUG.DAT	output	* DEBUG =	*
MASSFLX.DAT	output	* FLXDAT=	*
MASSBAL.DAT	output	* BALDAT=	*
FOG.DAT	output	* FOGDAT=	*

-----  
All file names will be converted to lower case if LCFILES = T  
Otherwise, if LCFILES = F, file names will be converted to UPPER CASE  
T = lower case ! LCFILES = T !  
F = UPPER CASE

NOTE: (1) file/path names can be up to 70 characters in length

Provision for multiple input files  
-----

**Refined BART Modeling Protocol for  
PCA Tomahawk Mill**

---

```

Number of CALMET.DAT files for run (NMETDAT)
          Default: 1          ! NMETDAT = 364 !

Number of PTEMARB.DAT files for run (NPTDAT)
          Default: 0          ! NPTDAT = 0 !

Number of BAEMARB.DAT files for run (NARDAT)
          Default: 0          ! NARDAT = 0 !

Number of VOLEMARB.DAT files for run (NVOLDAT)
          Default: 0          ! NVOLDAT = 0 !

```

!END!

```

-----
Subgroup (0a)
-----

```

The following CALMET.DAT filenames are processed in sequence if NMETDAT>1

Default Name	Type	File Name		
none	input	! METDAT=e:\WI-Calmet\2002output\cal020102.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020103.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020104.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020105.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020106.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020107.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020108.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020109.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020110.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020111.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020112.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020113.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020114.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020115.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020116.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020117.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020118.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020119.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020120.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020121.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020122.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020123.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020124.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020125.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020126.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020127.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020128.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020129.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020130.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020131.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020201.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020202.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020203.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020204.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020205.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020206.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020207.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020208.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020209.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020210.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020211.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020212.DAT	!	!END!













INPUT GROUP: 1 -- General run control parameters

-----

Option to run all periods found  
in the met. file (METRUN) Default: 0 ! METRUN = 0 !

METRUN = 0 - Run period explicitly defined below  
METRUN = 1 - Run all periods in met. file

Starting date: Year (IBYR) -- No default ! IBYR = 2002 !  
(used only if Month (IBMO) -- No default ! IBMO = 1 !  
METRUN = 0) Day (IBDY) -- No default ! IBDY = 02 !  
Hour (IBHR) -- No default ! IBHR = 1 !

Base time zone (XBTZ) -- No default ! XBTZ = 6.0 !  
PST = 8., MST = 7.  
CST = 6., EST = 5.

Length of run (hours) (IRLG) -- No default ! IRLG = 8784 !

Number of chemical species (NSPEC)  
Default: 5 ! NSPEC = 9 !

Number of chemical species  
to be emitted (NSE) Default: 3 ! NSE = 7 !

Flag to stop run after  
SETUP phase (ITEST) Default: 2 ! ITEST = 2 !  
(Used to allow checking  
of the model inputs, files, etc.)  
ITEST = 1 - STOPS program after SETUP phase  
ITEST = 2 - Continues with execution of program  
after SETUP

Restart Configuration:

Control flag (MRESTART) Default: 0 ! MRESTART = 0 !

0 = Do not read or write a restart file  
1 = Read a restart file at the beginning of  
the run  
2 = Write a restart file during run  
3 = Read a restart file at beginning of run  
and write a restart file during run

Number of periods in Restart  
output cycle (NRESPD) Default: 0 ! NRESPD = 0 !

0 = File written only at last period  
>0 = File updated every NRESPD periods

Meteorological Data Format (METFM)  
Default: 1 ! METFM = 1 !

METFM = 1 - CALMET binary file (CALMET.MET)  
METFM = 2 - ISC ASCII file (ISCMET.MET)  
METFM = 3 - AUSPLUME ASCII file (PLMMET.MET)  
METFM = 4 - CTDM plus tower file (PROFILE.DAT) and  
surface parameters file (SURFACE.DAT)

PG sigma-y is adjusted by the factor (AVET/PGTIME)\*\*0.2  
Averaging Time (minutes) (AVET)  
Default: 60.0 ! AVET = 60. !

PG Averaging Time (minutes) (PGTIME)  
Default: 60.0 ! PGTIME = 60. !

!END!

-----  
INPUT GROUP: 2 -- Technical options  
-----

Vertical distribution used in the  
near field (MGAUSS)                   Default: 1       ! MGAUSS = 1   !  
  0 = uniform  
  1 = Gaussian

Terrain adjustment method  
(MCTADJ)                               Default: 3       ! MCTADJ = 3   !  
  0 = no adjustment  
  1 = ISC-type of terrain adjustment  
  2 = simple, CALPUFF-type of terrain  
      adjustment  
  3 = partial plume path adjustment

Subgrid-scale complex terrain  
flag (MCTSG)                           Default: 0       ! MCTSG = 0   !  
  0 = not modeled  
  1 = modeled

Near-field puffs modeled as  
elongated 0 (MSLUG)                   Default: 0       ! MSLUG = 0   !  
  0 = no  
  1 = yes (slug model used)

Transitional plume rise modeled ?  
(MTRANS)                              Default: 1       ! MTRANS = 1   !  
  0 = no (i.e., final rise only)  
  1 = yes (i.e., transitional rise computed)

Stack tip downwash? (MTIP)           Default: 1       ! MTIP = 1   !  
  0 = no (i.e., no stack tip downwash)  
  1 = yes (i.e., use stack tip downwash)

Method used to simulate building  
downwash? (MBDW)                      Default: 1       ! MBDW = 1   !  
  1 = ISC method  
  2 = PRIME method

Vertical wind shear modeled above  
stack top? (MSHEAR)                   Default: 0       ! MSHEAR = 0   !  
  0 = no (i.e., vertical wind shear not modeled)  
  1 = yes (i.e., vertical wind shear modeled)

Puff splitting allowed? (MSPLIT)      Default: 0       ! MSPLIT = 0   !  
  0 = no (i.e., puffs not split)  
  1 = yes (i.e., puffs are split)

Chemical mechanism flag (MCHEM)       Default: 1       ! MCHEM = 1   !  
  0 = chemical transformation not  
      modeled  
  1 = transformation rates computed  
      internally (MESOPUFF II scheme)  
  2 = user-specified transformation  
      rates used  
  3 = transformation rates computed

internally (RIVAD/ARM3 scheme)  
4 = secondary organic aerosol formation  
   computed (MESOPUFF II scheme for OH)

Aqueous phase transformation flag (MAQCHEM)  
(Used only if MCHEM = 1, or 3)      Default: 0      ! MAQCHEM = 0      !  
0 = aqueous phase transformation  
   not modeled  
1 = transformation rates adjusted  
   for aqueous phase reactions

Wet removal modeled ? (MWET)            Default: 1      ! MWET = 1      !  
0 = no  
1 = yes

Dry deposition modeled ? (MDRY)         Default: 1      ! MDRY = 1      !  
0 = no  
1 = yes  
(dry deposition method specified  
  for each species in Input Group 3)

Method used to compute dispersion  
coefficients (MDISP)                    Default: 3      ! MDISP = 3      !  
  
1 = dispersion coefficients computed from measured values  
   of turbulence, sigma v, sigma w  
2 = dispersion coefficients from internally calculated  
   sigma v, sigma w using micrometeorological variables  
   (u\*, w\*, L, etc.)  
3 = PG dispersion coefficients for RURAL areas (computed using  
   the ISCST multi-segment approximation) and MP coefficients in  
   urban areas  
4 = same as 3 except PG coefficients computed using  
   the MESOPUFF II eqns.  
5 = CTDM sigmas used for stable and neutral conditions.  
   For unstable conditions, sigmas are computed as in  
   MDISP = 3, described above. MDISP = 5 assumes that  
   measured values are read

Sigma-v/sigma-theta, sigma-w measurements used? (MTURBVW)  
(Used only if MDISP = 1 or 5)         Default: 3      ! MTURBVW = 3      !  
1 = use sigma-v or sigma-theta measurements  
   from PROFILE.DAT to compute sigma-y  
   (valid for METFM = 1, 2, 3, 4)  
2 = use sigma-w measurements  
   from PROFILE.DAT to compute sigma-z  
   (valid for METFM = 1, 2, 3, 4)  
3 = use both sigma-(v/theta) and sigma-w  
   from PROFILE.DAT to compute sigma-y and sigma-z  
   (valid for METFM = 1, 2, 3, 4)  
4 = use sigma-theta measurements  
   from PLMMET.DAT to compute sigma-y  
   (valid only if METFM = 3)

Back-up method used to compute dispersion  
when measured turbulence data are  
missing (MDISP2)                        Default: 3      ! MDISP2 = 3      !  
(used only if MDISP = 1 or 5)  
2 = dispersion coefficients from internally calculated  
   sigma v, sigma w using micrometeorological variables  
   (u\*, w\*, L, etc.)  
3 = PG dispersion coefficients for RURAL areas (computed using  
   the ISCST multi-segment approximation) and MP coefficients in  
   urban areas  
4 = same as 3 except PG coefficients computed using

---

the MESOPUFF II eqns.

PG sigma-y,z adj. for roughness? Default: 0 ! MROUGH = 0 !  
(MROUGH)  
0 = no  
1 = yes

Partial plume penetration of Default: 1 ! MPARTL = 1 !  
elevated inversion?  
(MPARTL)  
0 = no  
1 = yes

Strength of temperature inversion Default: 0 ! MTINV = 0 !  
provided in PROFILE.DAT extended records?  
(MTINV)  
0 = no (computed from measured/default gradients)  
1 = yes

PDF used for dispersion under convective conditions? Default: 0 ! MPDF = 0 !  
(MPDF)  
0 = no  
1 = yes

Sub-Grid TIBL module used for shore line? Default: 0 ! MSGTIBL = 0 !  
(MSGTIBL)  
0 = no  
1 = yes

Boundary conditions (concentration) modeled? Default: 0 ! MBCON = 0 !  
(MBCON)  
0 = no  
1 = yes, using formatted BCON.DAT file  
2 = yes, using unformatted CONC.DAT file

Analyses of fogging and icing impacts due to emissions from  
arrays of mechanically-forced cooling towers can be performed  
using CALPUFF in conjunction with a cooling tower emissions  
processor (CTEMISS) and its associated postprocessors. Hourly  
emissions of water vapor and temperature from each cooling tower  
cell are computed for the current cell configuration and ambient  
conditions by CTEMISS. CALPUFF models the dispersion of these  
emissions and provides cloud information in a specialized format  
for further analysis. Output to FOG.DAT is provided in either  
'plume mode' or 'receptor mode' format.

Configure for FOG Model output? Default: 0 ! MFOG = 0 !  
(MFOG)  
0 = no  
1 = yes - report results in PLUME Mode format  
2 = yes - report results in RECEPTOR Mode format

TEST options specified to see if  
they conform to regulatory  
values? (MREG) Default: 1 ! MREG = 0 !  
0 = NO checks are made  
1 = Technical options must conform to USEPA  
Long Range Transport (LRT) guidance

---

```

METFM      1 or 2
AVET       60. (min)
PGTIME     60. (min)
MGAUSS     1
MCTADJ     3
MTRANS     1
MTIP       1
MCHEM      1 or 3 (if modeling SOx, NOx)
MWET       1
MDRY       1
MDISP      2 or 3
MPDF       0 if MDISP=3
           1 if MDISP=2
MROUGH     0
MPARTL     1
SYTDEP     550. (m)
MHFTSZ     0

```

!END!

-----  
INPUT GROUP: 3a, 3b -- Species list  
-----

-----  
Subgroup (3a)  
-----

The following species are modeled:

The following species are modeled:

```

! CSPEC =      SO2 !      !END!
! CSPEC =      SO4 !      !END!
! CSPEC =      NOX !      !END!
! CSPEC =      HNO3 !     !END!
! CSPEC =      NO3 !      !END!
! CSPEC =      PMC !      !END!
! CSPEC =      PMF !      !END!
! CSPEC =      EC !       !END!
! CSPEC =      SOA !      !END!

```

SPECIES NAME (Limit: 12 Characters in length)	MODELED (0=NO, 1=YES)	EMITTED (0=NO, 1=YES)	Dry DEPOSITED (0=NO, 1=COMPUTED-GAS 2=COMPUTED-PARTICLE 3=USER-SPECIFIED)	OUTPUT GROUP NUMBER (0=NONE, 1=1st CGRUP, 2=2nd CGRUP, 3= etc.)
! SO2 =	1,	1,	1,	0 !
! SO4 =	1,	1,	2,	0 !
! NOX =	1,	1,	1,	0 !
! HNO3 =	1,	0,	1,	0 !
! NO3 =	1,	0,	2,	0 !
! PMC =	1,	1,	2,	0 !
! PMF =	1,	1,	2,	0 !
! EC =	1,	1,	2,	0 !
! SOA =	1,	1,	2,	0 !

!END!







The computational grid is identical to or a subset of the MET. grid.  
 The lower left (LL) corner of the computational grid is at grid point  
 (IBCOMP, JBCOMP) of the MET. grid. The upper right (UR) corner of the  
 computational grid is at grid point (IECOMP, JECOMP) of the MET. grid.  
 The grid spacing of the computational grid is the same as the MET. grid.

X index of LL corner (IBCOMP) (1 <= IBCOMP <= NX)	No default	! IBCOMP = 1 !
Y index of LL corner (JBCOMP) (1 <= JBCOMP <= NY)	No default	! JBCOMP = 1 !
X index of UR corner (IECOMP) (1 <= IECOMP <= NX)	No default	! IECOMP = 162 !
Y index of UR corner (JECOMP) (1 <= JECOMP <= NY)	No default	! JECOMP = 191 !

SAMPLING Grid (GRIDDED RECEPTORS):

The lower left (LL) corner of the sampling grid is at grid point  
 (IBSAMP, JBSAMP) of the MET. grid. The upper right (UR) corner of the  
 sampling grid is at grid point (IESAMP, JESAMP) of the MET. grid.  
 The sampling grid must be identical to or a subset of the computational  
 grid. It may be a nested grid inside the computational grid.  
 The grid spacing of the sampling grid is DGRIDKM/MESHNDN.

Logical flag indicating if gridded receptors are used (LSAMP) (T=yes, F=no)	Default: T	! LSAMP = F !
X index of LL corner (IBSAMP) (IBCOMP <= IBSAMP <= IECOMP)	No default	! IBSAMP = 1 !
Y index of LL corner (JBSAMP) (JBCOMP <= JBSAMP <= JECOMP)	No default	! JBSAMP = 1 !
X index of UR corner (IESAMP) (IBCOMP <= IESAMP <= IECOMP)	No default	! IESAMP = 162 !
Y index of UR corner (JESAMP) (JBCOMP <= JESAMP <= JECOMP)	No default	! JESAMP = 191 !
Nesting factor of the sampling grid (MESHNDN) (MESHNDN is an integer >= 1)	Default: 1	! MESHNDN = 1 !

!END!

-----  
 INPUT GROUP: 5 -- Output Options  
 -----

FILE	* DEFAULT VALUE	* VALUE THIS RUN
----	-----	-----

```

Concentrations (ICON)           1           ! ICON = 1 !
Dry Fluxes (IDRY)               1           ! IDRY = 0 !
Wet Fluxes (IWET)              1           ! IWET = 0 !
Relative Humidity (IVIS)        1           ! IVIS = 0 !
  (relative humidity file is
  required for visibility
  analysis)
Use data compression option in output file?
(LCOMPRS)                       Default: T           ! LCOMPRS = T !

```

\*

0 = Do not create file, 1 = create file

DIAGNOSTIC MASS FLUX OUTPUT OPTIONS:

```

Mass flux across specified boundaries
for selected species reported hourly?
(IMFLX)                          Default: 0           ! IMFLX = 0 !
  0 = no
  1 = yes (FLUXBDY.DAT and MASSFLX.DAT filenames
  are specified in Input Group 0)

```

```

Mass balance for each species
reported hourly?
(IMBAL)                          Default: 0           ! IMBAL = 0 !
  0 = no
  1 = yes (MASSBAL.DAT filename is
  specified in Input Group 0)

```

LINE PRINTER OUTPUT OPTIONS:

```

Print concentrations (ICPRT)      Default: 0           ! ICPRT = 0 !
Print dry fluxes (IDPRT)        Default: 0           ! IDPRT = 0 !
Print wet fluxes (IWPRT)       Default: 0           ! IWPRT = 0 !
(0 = Do not print, 1 = Print)

```

```

Concentration print interval
(ICFRQ) in hours                 Default: 1           ! ICFRQ = 1 !
Dry flux print interval
(IDFRQ) in hours                 Default: 1           ! IDFRQ = 1 !
Wet flux print interval
(IWFRQ) in hours                 Default: 1           ! IWFRQ = 1 !

```

```

Units for Line Printer Output
(IPRTU)                          Default: 1           ! IPRTU = 3 !
      for
      Concentration   Deposition
  1 =      g/m**3      g/m**2/s
  2 =      mg/m**3     mg/m**2/s
  3 =      ug/m**3     ug/m**2/s
  4 =      ng/m**3     ng/m**2/s
  5 =      Odour Units

```

```

Messages tracking progress of run
written to the screen ?
(IMESG)                          Default: 2           ! IMESG = 2 !
  0 = no
  1 = yes (advection step, puff ID)
  2 = yes (YYYYJJJHH, # old puffs, # emitted puffs)

```

SPECIES (or GROUP for combined species) LIST FOR OUTPUT OPTIONS

**Refined BART Modeling Protocol for  
PCA Tomahawk Mill**

```

----- CONCENTRATIONS ----- ----- DRY FLUXES ----- ----- WET FLUXES -----
-- MASS FLUX --
SPECIES
/GROUP      PRINTED?  SAVED ON DISK?  PRINTED?  SAVED ON DISK?  PRINTED?  SAVED ON DISK?
SAVED ON DISK?
-----
!          SO2 =    0,      1,      0,      0,      0,      0,
0 !
!          SO4 =    0,      1,      0,      0,      0,      0,
0 !
!          NOX =    0,      1,      0,      0,      0,      0,
0 !
!          HNO3 =   0,      1,      0,      0,      0,      0,
0 !
!          NO3 =    0,      1,      0,      0,      0,      0,
0 !
!          PMC =    0,      1,      0,      0,      0,      0,
0 !
!          PMF =    0,      1,      0,      0,      0,      0,
0 !
!          EC =     0,      1,      0,      0,      0,      0,
0 !
!          SOA =    0,      1,      0,      0,      0,      0,
0 !

```

OPTIONS FOR PRINTING "DEBUG" QUANTITIES (much output)

```

Logical for debug output
(LDEBUG)                                Default: F      ! LDEBUG = F !

First puff to track
(IPFDEB)                                Default: 1      ! IPFDEB = 1 !

Number of puffs to track
(NPFDEB)                                Default: 1      ! NPFDEB = 1 !

Met. period to start output
(NN1)                                    Default: 1      ! NN1 = 1 !

Met. period to end output
(NN2)                                    Default: 10     ! NN2 = 10 !

!END!

```

INPUT GROUP: 6a, 6b, & 6c -- Subgrid scale complex terrain inputs

-----  
Subgroup (6a)  
-----

```

Number of terrain features (NHILL)      Default: 0      ! NHILL = 0 !

Number of special complex terrain
receptors (NCTREC)                      Default: 0      ! NCTREC = 0 !

Terrain and CTSG Receptor data for
CTSG hills input in CTDM format ?
(MHILL)                                  No Default     ! MHILL = 2 !
1 = Hill and Receptor data created

```

by CTDM processors & read from  
HILL.DAT and HILLRCT.DAT files  
2 = Hill data created by OPTHILL &  
input below in Subgroup (6b);  
Receptor data in Subgroup (6c)

Factor to convert horizontal dimensions Default: 1.0 ! XHILL2M = 1. !  
to meters (MHILL=1)

Factor to convert vertical dimensions Default: 1.0 ! ZHILL2M = 1. !  
to meters (MHILL=1)

X-origin of CTDM system relative to No Default ! XCTDMKM = 0.0E00 !  
CALPUFF coordinate system, in Kilometers (MHILL=1)

Y-origin of CTDM system relative to No Default ! YCTDMKM = 0.0E00 !  
CALPUFF coordinate system, in Kilometers (MHILL=1)

! END !

-----  
Subgroup (6b)  
-----

1 \*\*  
HILL information

HILL AMAX1 NO. (m)	XC AMAX2 (km) (m)	YC (km)	THETAH (deg.)	ZGRID (m)	RELIEF (m)	EXPO 1 (m)	EXPO 2 (m)	SCALE 1 (m)	SCALE 2 (m)
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----
----	----								

-----  
Subgroup (6c)  
-----

COMPLEX TERRAIN RECEPTOR INFORMATION

XRCT (km)	YRCT (km)	ZRCT (m)	XHH
-----	-----	-----	----

-----  
1

Description of Complex Terrain Variables:

XC, YC = Coordinates of center of hill  
THETAH = Orientation of major axis of hill (clockwise from  
North)  
ZGRID = Height of the 0 of the grid above mean sea  
level  
RELIEF = Height of the crest of the hill above the grid elevation  
EXPO 1 = Hill-shape exponent for the major axis  
EXPO 2 = Hill-shape exponent for the major axis  
SCALE 1 = Horizontal length scale along the major axis  
SCALE 2 = Horizontal length scale along the minor axis  
AMAX = Maximum allowed axis length for the major axis  
BMAX = Maximum allowed axis length for the major axis

XRCT, YRCT = Coordinates of the complex terrain receptors  
ZRCT = Height of the ground (MSL) at the complex terrain  
Receptor

XHH = Hill number associated with each complex terrain receptor  
(NOTE: MUST BE ENTERED AS A REAL NUMBER)

\*\*

NOTE: DATA for each hill and CTSG receptor are treated as a separate  
input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUP: 7 -- Chemical parameters for dry deposition of gases  
-----

SPECIES LAW COEFFICIENT NAME (dimensionless)	DIFFUSIVITY (cm**2/s)	ALPHA STAR	REACTIVITY	MESOPHYLL RESISTANCE (s/cm)	HENRY'S
! SO2 =	0.1509,	1000.,	8.,	0.,	
0.04 !					
! NOX =	0.1656,	1.,	8.,	5.,	
3.5 !					
! HNO3 =	0.1628,	1.,	18.,	0.,	
0.00000008 !					

!END!

-----  
INPUT GROUP: 8 -- Size parameters for dry deposition of particles  
-----

For SINGLE SPECIES, the mean and standard deviation are used to  
compute a deposition velocity for NINT (see group 9) size-ranges,  
and these are then averaged to obtain a mean deposition velocity.

For GROUPED SPECIES, the size distribution should be explicitly  
specified (by the 'species' in the group), and the standard deviation  
for each should be entered as 0. The model will then use the  
deposition velocity for the stated mean diameter.

SPECIES NAME	GEOMETRIC MASS MEAN DIAMETER (microns)	GEOMETRIC STANDARD DEVIATION (microns)
! SO4 =	0.48,	2. !
! NO3 =	0.48,	2. !
! PMC =	2.0,	2.0 !
! PMF =	0.48,	2. !
! EC =	0.48,	2. !
! SOA =	0.48,	2. !

!END!

-----  
INPUT GROUP: 9 -- Miscellaneous dry deposition parameters  
-----

```

Reference cuticle resistance (s/cm)
(RCUTR)                      Default: 30    ! RCUTR = 30.0 !
Reference ground resistance (s/cm)
(RGR)                        Default: 10    ! RGR = 10.0 !
Reference pollutant reactivity
(REACTR)                     Default: 8     ! REACTR = 8.0 !

Number of particle-size intervals used to
evaluate effective particle deposition velocity
(NINT)                        Default: 9     ! NINT = 9 !

Vegetation state in unirrigated areas
(IVEG)                        Default: 1     ! IVEG = 1 !
  IVEG=1 for active and unstressed vegetation
  IVEG=2 for active and stressed vegetation
  IVEG=3 for inactive vegetation

```

!END!

-----

INPUT GROUP: 10 -- Wet Deposition Parameters

-----

Scavenging Coefficient -- Units: (sec)\*\*(-1)

Pollutant	Liquid Precip.	Frozen Precip.
! SO2 =	3.0E-05,	0.0E00 !
! SO4 =	1.0E-04,	3.0E-05 !
! NOX =	0.0E00,	0.0E00 !
! HNO3 =	6.0E-05,	0.0E00 !
! NO3 =	1.0E-04,	3.0E-05 !

!END!

-----

INPUT GROUP: 11 -- Chemistry Parameters

-----

```

Ozone data input option (MOZ)   Default: 1           ! MOZ = 0 !
(Used only if MCHEM = 1, 3, or 4)
  0 = use a monthly background ozone value
  1 = read hourly ozone concentrations from
    the OZONE.DAT data file

```

```

Monthly ozone concentrations
(Used only if MCHEM = 1, 3, or 4 and
MOZ = 0 or MOZ = 1 and all hourly O3 data missing)
(BCKO3) in ppb                    Default: 12*80.

```

```

! BCKO3 = 31.00, 31.00, 31.00, 37.00, 37.00, 37.00, 33.00, 33.00, 33.00, 27.00, 27.00, 27.00
!

```

```

Monthly ammonia concentrations
(Used only if MCHEM = 1, or 3)
(BCKNH3) in ppb                    Default: 12*10.
! BCKNH3 = 12*3.9 !

```



\* BCKNH3 = 00.30, 00.30, 00.30, 00.50, 00.50, 00.50, 00.50, 00.50, 00.50, 00.50, 00.50, 00.50,  
00.50 \*

Nighttime SO2 loss rate (RNITE1)  
in percent/hour                   Default: 0.2                   ! RNITE1 = .2 !

Nighttime NOx loss rate (RNITE2)  
in percent/hour                   Default: 2.0                   ! RNITE2 = 2.0 !

Nighttime HNO3 formation rate (RNITE3)  
in percent/hour                   Default: 2.0                   ! RNITE3 = 2.0 !

H2O2 data input option (MH2O2)   Default: 1                   ! MH2O2 = 1 !  
(Used only if MAQCHEM = 1)  
  0 = use a monthly background H2O2 value  
  1 = read hourly H2O2 concentrations from  
      the H2O2.DAT data file

Monthly H2O2 concentrations  
(Used only if MQACHEM = 1 and  
  MH2O2 = 0 or MH2O2 = 1 and all hourly H2O2 data missing)  
(BCKH2O2) in ppb                   Default: 12\*1.  
\* BCKH2O2 = 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00 \*  
! BCKH2O2 = 0.50, 0.50, 0.50, 2.30, 2.30, 2.30, 3.50, 3.50, 3.50, 0.80, 0.80, 0.80 !

--- Data for SECONDARY ORGANIC AEROSOL (SOA) Option  
(used only if MCHEM = 4)

The SOA module uses monthly values of:  
  Fine particulate concentration in ug/m<sup>3</sup> (BCKPMF)  
  Organic fraction of fine particulate       (OFRAC)  
  VOC / NOX ratio (after reaction)       (VCNX)  
to characterize the air mass when computing  
the formation of SOA from VOC emissions.  
Typical values for several distinct air mass types are:

Month	1	2	3	4	5	6	7	8	9	10	11	12
	Jan	Feb	Mar	Apr	May	Jun	Jul	Aug	Sep	Oct	Nov	Dec
Clean Continental												
BCKPMF	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.
OFRAC	.15	.15	.20	.20	.20	.20	.20	.20	.20	.20	.20	.15
VCNX	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.
Clean Marine (surface)												
BCKPMF	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5
OFRAC	.25	.25	.30	.30	.30	.30	.30	.30	.30	.30	.30	.25
VCNX	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.
Urban - low biogenic (controls present)												
BCKPMF	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.
OFRAC	.20	.20	.25	.25	.25	.25	.25	.25	.20	.20	.20	.20
VCNX	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.
Urban - high biogenic (controls present)												
BCKPMF	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.
OFRAC	.25	.25	.30	.30	.30	.55	.55	.55	.35	.35	.35	.25
VCNX	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.
Regional Plume												
BCKPMF	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.
OFRAC	.20	.20	.25	.35	.25	.40	.40	.40	.30	.30	.30	.20
VCNX	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.

Urban - no controls present

```

BCKPMF 100. 100. 100. 100. 100. 100. 100. 100. 100. 100. 100. 100.
OFRAC  .30  .30  .35  .35  .35  .55  .55  .55  .35  .35  .35  .30
VCNX    2.   2.   2.   2.   2.   2.   2.   2.   2.   2.   2.   2.

```

Default: Clean Continental

```

! BCKPMF = 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00 !
! OFRAC  = 0.15, 0.15, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.15 !
! VCNX   = 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00,
50.00 !

```

!END!

-----

INPUT GROUP: 12 -- Misc. Dispersion and Computational Parameters

-----

Horizontal size of puff (m) beyond which  
time-dependent dispersion equations (Heffter)  
are used to determine sigma-y and  
sigma-z (SYTDEP) Default: 550. !SYTDEP = 5.5E02 !

Switch for using Heffter equation for sigma z  
as above (0 = Not use Heffter; 1 = use Heffter  
(MHFTSZ) Default: 0 ! MHFTSZ = 0 !

Stability class used to determine plume  
growth rates for puffs above the boundary  
layer (JSUP) Default: 5 ! JSUP = 5 !

Vertical dispersion constant for stable  
conditions (k1 in Eqn. 2.7-3) (CONK1) Default: 0.01 ! CONK1 = .01 !

Vertical dispersion constant for neutral/  
unstable conditions (k2 in Eqn. 2.7-4)  
(CONK2) Default: 0.1 ! CONK2 = .1 !

Factor for determining Transition-point from  
Schulman-Scire to Huber-Snyder Building Downwash  
scheme (SS used for Hs < Hb + TBD \* HL)  
(TBD) Default: 0.5 ! TBD = .5 !  
TBD < 0 ==> always use Huber-Snyder  
TBD = 1.5 ==> always use Schulman-Scire  
TBD = 0.5 ==> ISC Transition-point

Range of land use categories for which  
urban dispersion is assumed  
(IURB1, IURB2) Default: 10 ! IURB1 = 10 !  
19 ! IURB2 = 19 !

Site characterization parameters for single-point Met data files -----  
(needed for METFM = 2,3,4)

Land use category for modeling domain  
(ILANDUIN) Default: 20 ! ILANDUIN = 20 !

Roughness length (m) for modeling domain  
(Z0IN) Default: 0.25 ! Z0IN = .25 !

Leaf area index for modeling domain  
(XLAIIN) Default: 3.0 ! XLAIIN = 3.0 !

```

Elevation above sea level (m)
(ELEVIN)                                Default: 0.0    ! ELEVIN = .0 !

Latitude (degrees) for met location
(XLATIN)                                Default: -999.  ! XLATIN = -999.0 !

Longitude (degrees) for met location
(XLONIN)                                Default: -999.  ! XLONIN = -999.0 !

Specialized information for interpreting single-point Met data files -----

Anemometer height (m) (Used only if METFM = 2,3)
(ANEMHT)                                Default: 10.    ! ANEMHT = 10.0 !

Form of lateral turbulence data in PROFILE.DAT file
(Used only if METFM = 4 or MTUREBVW = 1 or 3)
(ISIGMAV)                                Default: 1      ! ISIGMAV = 1 !
    0 = read sigma-theta
    1 = read sigma-v

Choice of mixing heights (Used only if METFM = 4)
(IMIXCTDM)                               Default: 0      ! IMIXCTDM = 0 !
    0 = read PREDICTED mixing heights
    1 = read OBSERVED mixing heights

Maximum length of a slug (met. grid units)
(XMXLEN)                                Default: 1.0    ! XMXLEN = 1.0 !

Maximum travel distance of a puff/slug (in
grid units) during one sampling step
(XSAMPLN)                               Default: 1.0    ! XSAMPLN = 1.0 !

Maximum Number of slugs/puffs release from
one source during one time step
(MXNEW)                                  Default: 99     ! MXNEW = 99 !

Maximum Number of sampling steps for
one puff/slug during one time step
(MXSAM)                                  Default: 99     ! MXSAM = 99 !

Number of iterations used when computing
the transport wind for a sampling step
that includes gradual rise (for CALMET
and PROFILE winds)
(NCOUNT)                               Default: 2      ! NCOUNT = 2 !

Minimum sigma y for a new puff/slug (m)
(SYMIN)                                  Default: 1.0    ! SYMIN = 1.0 !

Minimum sigma z for a new puff/slug (m)
(SZMIN)                                  Default: 1.0    ! SZMIN = 1.0 !

Default minimum turbulence velocities
sigma-v and sigma-w for each
stability class (m/s)
(SVMIN(6) and SWMIN(6))                 Default SVMIN : .50, .50, .50, .50, .50, .50
                                           Default SWMIN : .20, .12, .08, .06, .03, .016

                                           Stability Class :  A    B    C    D    E    F
                                           ---  ---  ---  ---  ---  ---
                                           ! SVMIN = 0.500, 0.500, 0.500, 0.500, 0.500, 0.500!
                                           ! SWMIN = 0.200, 0.120, 0.080, 0.060, 0.030, 0.016!

Divergence criterion for dw/dz across puff
used to initiate adjustment for horizontal

```

```

convergence (1/s)
Partial adjustment starts at CDIV(1), and
full adjustment is reached at CDIV(2)
(CDIV(2))                                Default: 0.0,0.0 ! CDIV = .0, .0 !

Minimum wind speed (m/s) allowed for
non-calm conditions. Also used as minimum
speed returned when using power-law
extrapolation toward surface
(WSCALM)                                Default: 0.5 ! WSCALM = .5 !

Maximum mixing height (m)
(XMAXZI)                                Default: 3000. ! XMAXZI = 3000.0 !

Minimum mixing height (m)
(XMINZI)                                Default: 50. ! XMINZI = 50.0 !

Default wind speed classes --
5 upper bounds (m/s) are entered;
the 6th class has no upper limit
(WSCAT(5))                                Default :
ISC RURAL : 1.54, 3.09, 5.14, 8.23, 10.8 (10.8+)

Wind Speed Class : 1      2      3      4      5
                   ---    ---    ---    ---    ---
                   ! WSCAT = 1.54, 3.09, 5.14, 8.23, 10.80 !

Default wind speed profile power-law
exponents for stabilities 1-6
(PLX0(6))                                Default : ISC RURAL values
ISC RURAL : .07, .07, .10, .15, .35, .55
ISC URBAN : .15, .15, .20, .25, .30, .30

Stability Class :  A      B      C      D      E      F
                   ---    ---    ---    ---    ---    ---
                   ! PLX0 = 0.07, 0.07, 0.10, 0.15, 0.35, 0.55 !

Default potential temperature gradient
for stable classes E, F (degK/m)
(PG0(2))                                Default: 0.020, 0.035
                   ! PG0 = 0.020, 0.035 !

Default plume path coefficients for
each stability class (used when option
for partial plume height terrain adjustment
is selected -- MCTADJ=3)
(PPC(6))                                Stability Class :  A      B      C      D      E      F
Default PPC : .50, .50, .50, .50, .35, .35
                   ---    ---    ---    ---    ---    ---
                   ! PPC = 0.50, 0.50, 0.50, 0.50, 0.35, 0.35 !

Slug-to-puff transition criterion factor
equal to sigma-y/length of slug
(SL2PF)                                Default: 10. ! SL2PF = 10.0 !

Puff-splitting control variables -----

VERTICAL SPLIT
-----

Number of puffs that result every time a puff
is split - nsplit=2 means that 1 puff splits
into 2
(NSPLIT)                                Default: 3 ! NSPLIT = 3 !

```

Time(s) of a day when split puffs are eligible to be split once again; this is typically set once per day, around sunset before nocturnal shear develops.  
 24 values: 0 is midnight (00:00) and 23 is 11 PM (23:00)  
 0=do not re-split 1=eligible for re-split  
 (IRESPLIT(24)) Default: Hour 17 = 1  
 ! IRESPLIT = 0,0 !

Split is allowed only if last hour's mixing height (m) exceeds a minimum value  
 (ZISPLIT) Default: 100. ! ZISPLIT = 100.0 !

Split is allowed only if ratio of last hour's mixing ht to the maximum mixing ht experienced by the puff is less than a maximum value (this postpones a split until a nocturnal layer develops)  
 (ROLDMAX) Default: 0.25 ! ROLDMAX = 0.25 !

HORIZONTAL SPLIT  
 -----

Number of puffs that result every time a puff is split - nsplith=5 means that 1 puff splits into 5  
 (NSPLITH) Default: 5 ! NSPLITH = 5 !

Minimum sigma-y (Grid Cells Units) of puff before it may be split  
 (SYSPLITH) Default: 1.0 ! SYSPLITH = 1.0 !

Minimum puff elongation rate (SYSPLITH/hr) due to wind shear, before it may be split  
 (SHSPLITH) Default: 2. ! SHSPLITH = 2.0 !

Minimum concentration (g/m<sup>3</sup>) of each species in puff before it may be split  
 Enter array of NSPEC values; if a single value is entered, it will be used for ALL species  
 (CNSPLITH) Default: 1.0E-07 ! CNSPLITH = 1.0E-07 !

Integration control variables -----

Fractional convergence criterion for numerical SLUG sampling integration  
 (EPSSLUG) Default: 1.0e-04 ! EPSSLUG = 1.0E-04 !

Fractional convergence criterion for numerical AREA source integration  
 (EPSAREA) Default: 1.0e-06 ! EPSAREA = 1.0E-06 !

Trajectory step-length (m) used for numerical rise integration  
 (DSRISE) Default: 1.0 ! DSRISE = 1.0 !

Boundary Condition (BC) Puff control variables -----

Minimum height (m) to which BC puffs are mixed as they are emitted (MBCON=2 ONLY). Actual height is reset to the current mixing height at the release point if greater than this minimum.  
 (HTMINBC) Default: 500. ! HTMINBC = 500.0 !

Search radius (in BC segment lengths) about a receptor for sampling nearest BC puff. BC puffs are emitted with a spacing of one segment length, so the search radius should be greater than 1.

(RSAMPBC) Default: 4. ! RSAMPBC = 10.0 !  
Near-Surface depletion adjustment to concentration profile used when  
sampling BC puffs?  
(MDEPBC) Default: 1 ! MDEPBC = 1 !  
0 = Concentration is NOT adjusted for depletion  
1 = Adjust Concentration for depletion  
  
!END!

-----  
INPUT GROUPS: 13a, 13b, 13c, 13d -- Point source parameters  
-----

-----  
Subgroup (13a)  
-----

Number of point sources with  
parameters provided below (NPT1) No default ! NPT1 = 1 !  
  
Units used for point source  
emissions below (IPTU) Default: 1 ! IPTU = 1 !  
1 = g/s  
2 = kg/hr  
3 = lb/hr  
4 = tons/yr  
5 = Odour Unit \* m\*\*3/s (vol. flux of odour compound)  
6 = Odour Unit \* m\*\*3/min  
7 = metric tons/yr

Number of source-species  
combinations with variable  
emissions scaling factors  
provided below in (13d) (NSPT1) Default: 0 ! NSPT1 = 0 !

Number of point sources with  
variable emission parameters  
provided in external file (NPT2) No default ! NPT2 = 0 !

(If NPT2 > 0, these point  
source emissions are read from  
the file: PTEMARB.DAT)

!END!

-----  
Subgroup (13b)  
-----

The following species are modeled:

CSPEC = SO2  
CSPEC = SO4  
CSPEC = NOX  
CSPEC = HNO3  
CSPEC = NO3  
CSPEC = PMC  
CSPEC = PMF  
CSPEC = EC  
CSPEC = SOA

a

POINT SOURCE: CONSTANT DATA

```

-----
Source          X          Y          Stack          Base          Stack          Exit          Exit          b          c
No.            Coordinate Coordinate Height Elevation Diameter  Vel.  Temp.  Bldg. Emission
              (km)         (km)         (m)         (m)         (m)         (m/s) (deg. K)  Dwash Rates
-----

```

```

1 ! SRCNAM = P1 !
1 ! X = 568.01, 625.9200, 60.700,360.00, 3.230, 17.0, 468.0, .0,
      166.3E00, 2.17E00, 13.4E00, 0.0E00, 0.0E00, 0.47, 0.98, 0.04, 0.54 !
1 ! FMFAC = 1.0 ! !END!

```

-----

a  
Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.

SRCNAM is a 12-character name for a source (No default)  
X is an array holding the source data listed by the column headings (No default)  
SIGYZI is an array holding the initial sigma-y and sigma-z (m) (Default: 0.,0.)  
FMFAC is a vertical momentum flux factor (0. or 1.0) used to represent the effect of rain-caps or other physical configurations that reduce momentum rise associated with the actual exit velocity. (Default: 1.0 -- full momentum used)

b  
0. = No building downwash modeled, 1. = downwash modeled  
NOTE: must be entered as a REAL number (i.e., with decimal point)

c  
An emission rate must be entered for every pollutant modeled. Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by IPTU (e.g. 1 for g/s).

-----  
Subgroup (13c)  
-----

BUILDING DIMENSION DATA FOR SOURCES SUBJECT TO DOWNWASH

```

Source          a
No.            Effective building height, width, length and X/Y offset (in meters)
              every 10 degrees. LENGTH, XBADJ, and YBADJ are only needed for
              MBDW=2 (PRIME downwash option)
-----

```

-----

a  
Building height, width, length, and X/Y offset from the source are treated as a separate input subgroup for each source and therefore must end with an input group terminator. The X/Y offset is the position, relative to the stack, of the center of the upwind face of the projected building, with the x-axis pointing along the flow direction.

-----

Subgroup (13d)

-----  
a  
POINT SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 13b. Factors entered multiply the rates in 13b. Skip sources here that have constant emissions. For more elaborate variation in source parameters, use PTEMARB.DAT and NPT2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0  
0 = Constant  
1 = Diurnal cycle (24 scaling factors: hours 1-24)  
2 = Monthly cycle (12 scaling factors: months 1-12)  
3 = Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)  
4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)  
5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 14a, 14b, 14c, 14d -- Area source parameters  
-----

-----  
Subgroup (14a)  
-----

Number of polygon area sources with parameters specified below (NAR1) No default ! NAR1 = 0 !

Units used for area source emissions below (IARU) Default: 1 ! IARU = 1 !

- 1 = g/m\*\*2/s
- 2 = kg/m\*\*2/hr
- 3 = lb/m\*\*2/hr
- 4 = tons/m\*\*2/yr
- 5 = Odour Unit \* m/s (vol. flux/m\*\*2 of odour compound)
- 6 = Odour Unit \* m/min
- 7 = metric tons/m\*\*2/yr

Number of source-species combinations with variable emissions scaling factors provided below in (14d) (NSAR1) Default: 0 ! NSAR1 = 0 !

Number of buoyant polygon area sources with variable location and emission



parameters (NAR2) No default ! NAR2 = 0 !  
(If NAR2 > 0, ALL parameter data for  
these sources are read from the file: BAEMARB.DAT)

!END!

-----  
Subgroup (14b)  
-----

a  
AREA SOURCE: CONSTANT DATA  
-----

Source No.	Effect. Height (m)	Base Elevation (m)	Initial Sigma z (m)	Emission Rates
---------------	--------------------------	--------------------------	---------------------------	-------------------

b

-----  
a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.  
b  
An emission rate must be entered for every pollutant modeled.  
Enter emission rate of zero for secondary pollutants that are  
modeled, but not emitted. Units are specified by IARU  
(e.g. 1 for g/m\*\*2/s).

-----  
Subgroup (14c)  
-----

COORDINATES (km) FOR EACH VERTEX(4) OF EACH POLYGON  
-----

Source No.	Ordered list of X followed by list of Y, grouped by source
---------------	--

a

-----  
a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.

-----  
Subgroup (14d)  
-----

a  
AREA SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission  
rates given in 14b. Factors entered multiply the rates in 14b.  
Skip sources here that have constant emissions. For more elaborate  
variation in source parameters, use BAEMARB.DAT and NAR2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0  
0 = Constant  
1 = Diurnal cycle (24 scaling factors: hours 1-24)  
2 = Monthly cycle (12 scaling factors: months 1-12)  
3 = Hour & Season (4 groups of 24 hourly scaling factors,  
where first group is DEC-JAN-FEB)

- 4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)
- 5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a

Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----

INPUT GROUPS: 15a, 15b, 15c -- Line source parameters  
-----

-----  
Subgroup (15a)  
-----

Number of buoyant line sources  
with variable location and emission  
parameters (NLN2)

No default ! NLN2 = 0 !

(If NLN2 > 0, ALL parameter data for  
these sources are read from the file: LNEARB.DAT)

Number of buoyant line sources (NLINES)

No default ! NLINES = 0 !

Units used for line source  
emissions below

(ILNU)

Default: 1 ! ILNU = 1 !

- 1 = g/s  
2 = kg/hr  
3 = lb/hr  
4 = tons/yr  
5 = Odour Unit \* m\*\*3/s (vol. flux of odour compound)  
6 = Odour Unit \* m\*\*3/min  
7 = metric tons/yr

Number of source-species  
combinations with variable  
emissions scaling factors  
provided below in (15c)

(NSLN1) Default: 0 ! NSLN1 = 0 !

Maximum number of segments used to model  
each line (MXNSEG)

Default: 7 ! MXNSEG = 7 !

The following variables are required only if NLINES > 0. They are  
used in the buoyant line source plume rise calculations.

Number of distances at which  
transitional rise is computed

Default: 6 ! NLRISE = 6 !

Average building length (XL)

No default ! XL = .0 !  
(in meters)

Average building height (HBL)

No default ! HBL = .0 !  
(in meters)

```

Average building width (WBL)           No default ! WBL = .0 !
                                         (in meters)

Average line source width (WML)        No default ! WML = .0 !
                                         (in meters)

Average separation between buildings (DXL) No default ! DXL = .0 !
                                         (in meters)

Average buoyancy parameter (FPRIMEL)   No default ! FPRIMEL = .0 !
                                         (in m**4/s**3)

```

!END!

-----  
Subgroup (15b)  
-----

BUOYANT LINE SOURCE: CONSTANT DATA  
-----

Source No.	Beg. X Coordinate (km)	Beg. Y Coordinate (km)	End. X Coordinate (km)	End. Y Coordinate (km)	Release Height (m)	Base Elevation (m)	Emission Rates	a
-----	-----	-----	-----	-----	-----	-----	-----	

a  
Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.

b  
An emission rate must be entered for every pollutant modeled. Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by ILNTU (e.g. 1 for g/s).

-----  
Subgroup (15c)  
-----

a  
BUOYANT LINE SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 15b. Factors entered multiply the rates in 15b. Skip sources here that have constant emissions.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

- 0 = Constant
- 1 = Diurnal cycle (24 scaling factors: hours 1-24)
- 2 = Monthly cycle (12 scaling factors: months 1-12)
- 3 = Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)
- 4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)
- 5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup  
and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 16a, 16b, 16c -- Volume source parameters  
-----

-----  
Subgroup (16a)  
-----

Number of volume sources with  
parameters provided in 16b,c (NVL1)      No default    !    NVL1 = 0    !

Units used for volume source  
emissions below in 16b      (IVLU)      Default: 1    !    IVLU = 1    !

- 1 =            g/s
- 2 =            kg/hr
- 3 =            lb/hr
- 4 =            tons/yr
- 5 =            Odour Unit \* m\*\*3/s (vol. flux of odour compound)
- 6 =            Odour Unit \* m\*\*3/min
- 7 =            metric tons/yr

Number of source-species  
combinations with variable  
emissions scaling factors  
provided below in (16c)      (NSVL1)      Default: 0    !    NSVL1 = 0    !

Number of volume sources with  
variable location and emission  
parameters      (NVL2)      No default    !    NVL2 = 0    !

(If NVL2 > 0, ALL parameter data for  
these sources are read from the VOLEMARB.DAT file(s) )

!END!

-----  
Subgroup (16b)  
-----

a  
VOLUME SOURCE: CONSTANT DATA  
-----

X	Y	Effect.	Base	Initial	Initial	b
Coordinate	Coordinate	Height	Elevation	Sigma y	Sigma z	Emission
(km)	(km)	(m)	(m)	(m)	(m)	Rates
-----	-----	-----	-----	-----	-----	-----

-----  
a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.

b  
An emission rate must be entered for every pollutant modeled.

Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by IVLU (e.g. 1 for g/s).

-----  
Subgroup (16c)  
-----

a  
VOLUME SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 16b. Factors entered multiply the rates in 16b. Skip sources here that have constant emissions. For more elaborate variation in source parameters, use VOLEMARB.DAT and NVL2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

0 = Constant  
1 = Diurnal cycle (24 scaling factors: hours 1-24)  
2 = Monthly cycle (12 scaling factors: months 1-12)  
3 = Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)  
4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12  
5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 17a & 17b -- Non-gridded (discrete) receptor information  
-----

-----  
Subgroup (17a)  
-----

Number of non-gridded receptors (NREC) No default ! NREC = 1995 !  
!END!

Note - three Class I areas  
Seney SENE = 173 receptors  
Isle Royal ISLE = 966 receptors  
Boundary Waters BOWA = 856 receptors  
total number of receptors is 1995

ALL RECEPTOR COORDINATES FOLLOW IN INPUT FILE

Updated CALPUFF Modeling  
Protocol and Final Report for the  
Subject to BART Analyses  
for Thilmany Kaukauna Mill

Prepared by

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December 2010

**Updated CALPUFF Modeling Protocol and Final Report for  
Thilmany “Subject to BART” Analysis**

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Appendix A – Example CALMET Input File

Appendix B – Example CALPUFF Input File

## 1.0 Introduction

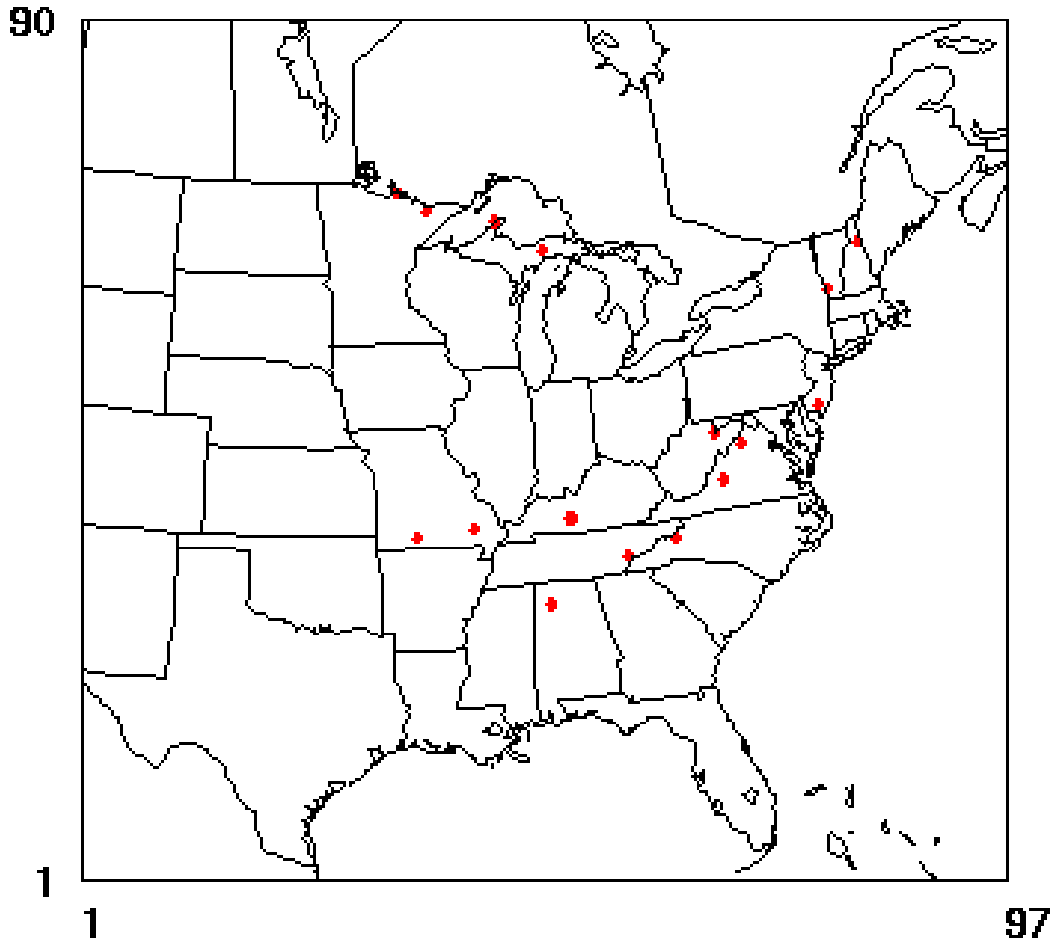
The Wisconsin Department of Natural Resources (DNR) performed single source screening modeling in 2006 to evaluate which Best Available Retrofit Technology (BART) eligible sources in the state may “cause or contribute” to visibility impairment and could therefore be subject to BART requirements. This type of analysis is referred to as the “subject to BART” or “BART exemption” analysis. The methods used by DNR are described in the “Single Source Modeling to Support Regional Haze BART Modeling Protocol”, March 21, 2006, prepared by Lake Michigan Air Directors Consortium (LADCO).

DNR applied CALPUFF to each BART eligible source for three annual periods, covering the calendar years 2002 to 2004. The CALPUFF/CALMET modeling domain consisted of ninety-seven 36-km spaced cells in the east-west direction and ninety 36-km spaced cells in the north-south direction (see Figure 1). Meteorological data generated by the MM5 prognostic weather model (provided by LADCO in CALMM5 format) was used to develop the meteorological data for CALPUFF, without additional observational data input (i.e., NOOBS mode). This represents a coarse meteorological grid, suitable for an initial screening analysis.

The results from DNR’s screening analysis indicated that, with the exception of some of the largest electric power stations in the state, the only Class I areas where BART eligible sources may contribute to visibility impairment are the three areas to the north of Wisconsin, the Boundary Waters Canoe Area, Isle Royal National Park, and Seney Wildlife Area. DNR’s screening analysis also indicated that the BART eligible emissions unit at the Thilmany Kaukauna Mill (Thilmany) may “cause or contribute” to visibility impairment at the Seney Class I area (located approximately 280 km to the northeast of the mill), as defined by the 98<sup>th</sup> percentile visibility impacts being greater than 0.5 deciviews (dv) relative to natural background.



Figure 1 - WDNR Screening CALPUFF Domain



NOTE: Class I areas are identified by red marks.

Thilmany has elected to refine the DNR screening analysis by conducting refined CALPUFF modeling with a 4 km meteorological grid, based on procedures described in the “Visibility Improvement – State and Tribal Association of the Southeast (VISTAS) BART CALPUFF” modeling protocol. The VISTAS protocol was developed by a large group of stakeholders, including EPA, the VISTAS member state agencies and tribes, the Federal Land Managers (FLMs), industry representatives, and consulting experts. Given the resources expended on the VISTAS protocol, it arguably represents the most comprehensive CALPUFF BART protocol ever developed.

Thilmany had discussions with DNR and EPA Region 5 staff about refining the analysis, and submitted a refined CALPUFF BART modeling protocol in March 2007. EPA reviewed the protocol and provided minor comments to DNR in May 2010. Thilmany responded to these comments in a letter “Response to EPA Comments on Thilmany BART Modeling Protocol” dated November 9, 2010. In addition, subsequent to the EPA protocol comments, the US Fish and Wildlife Service and US Forest Service reviewed the protocol and had additional discussions with DNR and EPA. The three EPA comments, and each response, are listed below.

**Comment 1**

Information needs to be presented which clarifies how POSTUTIL was used for ammonia.

**Response 1**

The VISTAS methodology was used, with a MNITRATE switch setting of one (1).

**Comment 2**

The modeling must include a copy of the input and list files associated with runs. The protocol only offers some of the control file information.

**Response 2**

All CALMET, CALPUFF, and CALPOST input and list files, as well as meteorological and geophysical data files, had been provided by Thilmany to DNR on a portable hard disk drive.

Thilmany will again provide a portable hard disk drive of all modeling files to DNR for the updated BART analysis, for purposes of SIP documentation.

**Comment 3**

The Thilmany input file ... uses the background ammonia values that were produced for the LADCO protocol. These values are based on modeled estimates. Monitoring data has generally shown much higher ammonia than what was used from the model results. The refined modeling analyses for Thilmany ... must evaluate the visibility impacts using measured background ammonia values.

**Response 3**

The US Fish and Wildlife Service proposed using a constant background ammonia concentration of 3.9 ppb, which was the highest concentration measured at the Mayville Wisconsin station during a mid-1990s EPA study. Thilmany has agreed to use this single, conservative background value in the absence of daily data.

On November 29, 2010, DNR issued a letter that approved the Thilmany CALPUFF BART protocol, as well as the emission rates to be used for the BART eligible emission units. DNR requested that the 2007 version of the protocol be updated to include the responses to EPA comments, and text that addresses current CALPUFF modeling guidance. In addition, the updated modeling results could be presented in the protocol (so that all information is located in one document).

This document is the updated modeling protocol and final modeling report for the refined Thilmany "Subject to BART" analysis. This document does not duplicate the extensive information on visibility, CALPUFF, and EPA guidance on BART modeling procedures that are contained in the VISTAS protocol. Rather, this document describes how data specific to this setting will be used in the refined CALPUFF modeling. Section 1 of the protocol is this introduction. Section 2 briefly summarizes important concepts from the VISTAS protocol.

Section 3 describes the specific CALMET data and procedures that were used to develop a refined, 4-km spaced meteorological grid for Wisconsin. Section 4 describes the CALPUFF input data and procedures, and Section 5 discusses the POSTUTIL and CALPOST processing, including the use of annual average natural background visibility. Section 6 presents the emission and stack data that were used for the BART eligible emission units, and Section 7 presents the modeling results. Appendices A and B present example CALMET and CALPUFF input file listings. A portable hard drive is also being submitted that contains all input data, control files, output files, and computer codes used in the analysis.

## **2.0 VISTAS BART Protocol**

The VISTAS member states have prepared a CALPUFF modeling protocol for BART determinations that is fully consistent with the EPA guidelines in 40 CFR Part 51 Appendix W and Appendix Y. The VISTAS protocol describes a two-step CALPUFF analysis methodology for BART exemption analyses. The initial CALPUFF analysis uses a “screening” or coarse resolution meteorological grid to determine if a particular source may be exempted from further BART analyses. The screening results are also used to determine which Class I areas should be included in any refined analysis. Assumptions for the initial screening analysis are conservative so that a source that contributes to visibility impairment is not exempted in error. If a source is shown to contribute to visibility impairment using the initial screening assessment, the source has the option to undertake refined CALPUFF modeling using finer meteorological grids to evaluate further whether it is subject to BART.

VISTAS developed both coarse (12-km grid resolution) and refined (4-km grid resolution) CALMET files. The fine grid CALMET files utilized both MM5 prognostic data, as well as meteorological observational data for the CALMET “Step 2” calculations. Overwater (buoy) data were used in addition to the standard hourly surface meteorological observations, precipitation observations, and twice-daily upper air sounding data. The VISTAS website provides detailed

documentation and supporting information on the refined CALMET processing, including the CALMET and CALPUFF input files that contain model configuration options and settings that have been optimized after consultation with stakeholders and consulting experts. These same CALMET and CALPUFF input files were used as the basis for the refined Wisconsin analyses.

The VISTAS protocol discusses several options for defining the natural background visibility. These include use of the annual average natural background extinction or the background extinction for the 20% best natural conditions days. Based on an email survey of the state agencies from the VISTAS member states, all of the VISTAS states allowed the use of annual average natural background extinction for refined CALPUFF analyses. This is also the guidance provided by EPA Region 5 staff during discussions with DNR and the Wisconsin Paper Council, and the method approved by DNR in their November 29, 2010 letter. Therefore, the refined Thilmany CALPUFF modeling used the annual average natural background extinction for CALPOST processing.

The VISTAS BART guidance recommends that the threshold value used to define whether a source “contributes” to visibility impairment is a 0.5 dv change from natural background conditions. The 98<sup>th</sup> percentile 24-hr average predicted impact at the Class I area (equal to the 8<sup>th</sup> highest value) is to be compared to this contribution threshold value. According to clarification of the BART guidance received from EPA by the VISTAS workgroup, for a three-year simulation the modeling values to be compared with the threshold are the greatest of the three annual 8<sup>th</sup> highest values, or the 22<sup>nd</sup> highest value over all three years combined, whichever is greater.

### 3.0 CALMET Data and Procedures

Since the 2007 version of the modeling protocol was developed, EPA has updated the versions of the approved CALPUFF programs, and has provided additional guidance on CALMET “switch” settings. The current approved versions are CALPUFF version 5.8 (level 070623), and CALMET version 5.8 (level 070623), and these versions were used for the updated refined BART CALPUFF analysis. On August 31, 2009, EPA published the memorandum “Clarification of EPA-FLM Recommended Settings for CALMET”. EPA stated that a 4 km fine CALMET grid size was acceptable, and provided recommended CALMET settings. The CALMET settings for the updated refined BART CALPUFF analysis meet all current EPA recommendations.

The first step in development of the refined Wisconsin analysis was to define the CALMET refined modeling domain and grid. As discussed in Section 1, DNR’s screening analysis indicated that there are three Class I areas to the north of Wisconsin, the Boundary Waters Canoe Area, Isle Royal National Park, and Seney Wildlife Area, where the screening analysis predicts some level of visibility impact from Wisconsin BART eligible sources. Therefore, the meteorological grid was developed to ensure coverage for these three Class I areas, as well as for the BART eligible source locations throughout Wisconsin. Standard IWAQM guidance also recommends that the CALMET grid extend a minimum of 50 km beyond any Class I or emission source location, so that puff recirculation can be properly accounted for.

The refined CALMET grid was defined as a 4-km spaced grid with 162 cells in the X-axis and 191 in the Y-axis. The grid projection was Lambert Conformal Conic, using the same projection parameters as in the DNR screening analysis (this allows the use of the same source X and Y coordinates as in the DNR analysis). The projection parameters include RLAT0 = 40N, RLON0 = 97W, XLAT1 = 33N, XLAT2 = 45N, XORIGKM = 247.0, YORIGKM = 240.0, and DATUM = NAS-C.

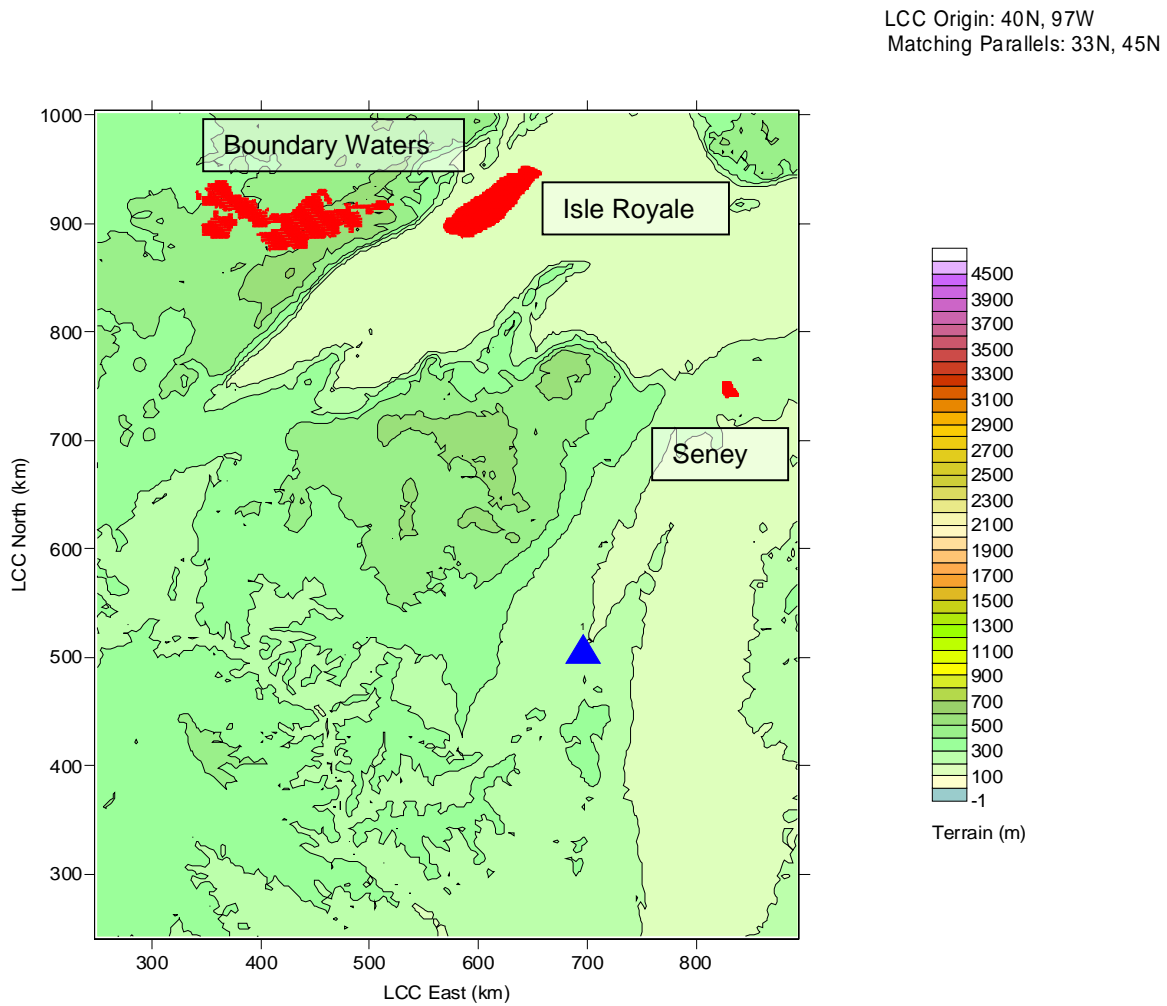
The meteorological data that was input to CALMET included the same three years (2002-2004) of CALMM5 data used by DNR, which was provided by LADCO. However, the refined CALMET runs also considered observation data from 22 surface stations, 16 precipitation stations, 5 overwater buoy stations, and 3 upper air stations. Figure 2 is a plot of the refined CALPUFF modeling domain. Figure 3 presents a plot of the locations of the observation stations in the modeling domain.

The CALMET program MAKEGEO uses USGS GTOPO30 digital terrain data (which includes coverage in Canada) and USGS Composite Theme Grid Land Use and Land Cover (LULC) data to create a gridded data file on geophysical parameters that is used by CALMET. The available USGS LULC coverage does not extend into the Canadian areas of the refined domain. These areas in Canada were assigned to the “forested land” category. Figure 4 presents a plot of the land use data for the modeling domain.

The VISTAS refined CALMET input files discussed in section 2 were used as the basis for the refined Wisconsin CALMET input files. The same CALMET configuration settings and options were utilized, as they are representative of Wisconsin terrain and land use patterns. All settings conform with current EPA and FLM guidance.

Appendix A presents a listing of a CALMET input file developed for the refined analysis.

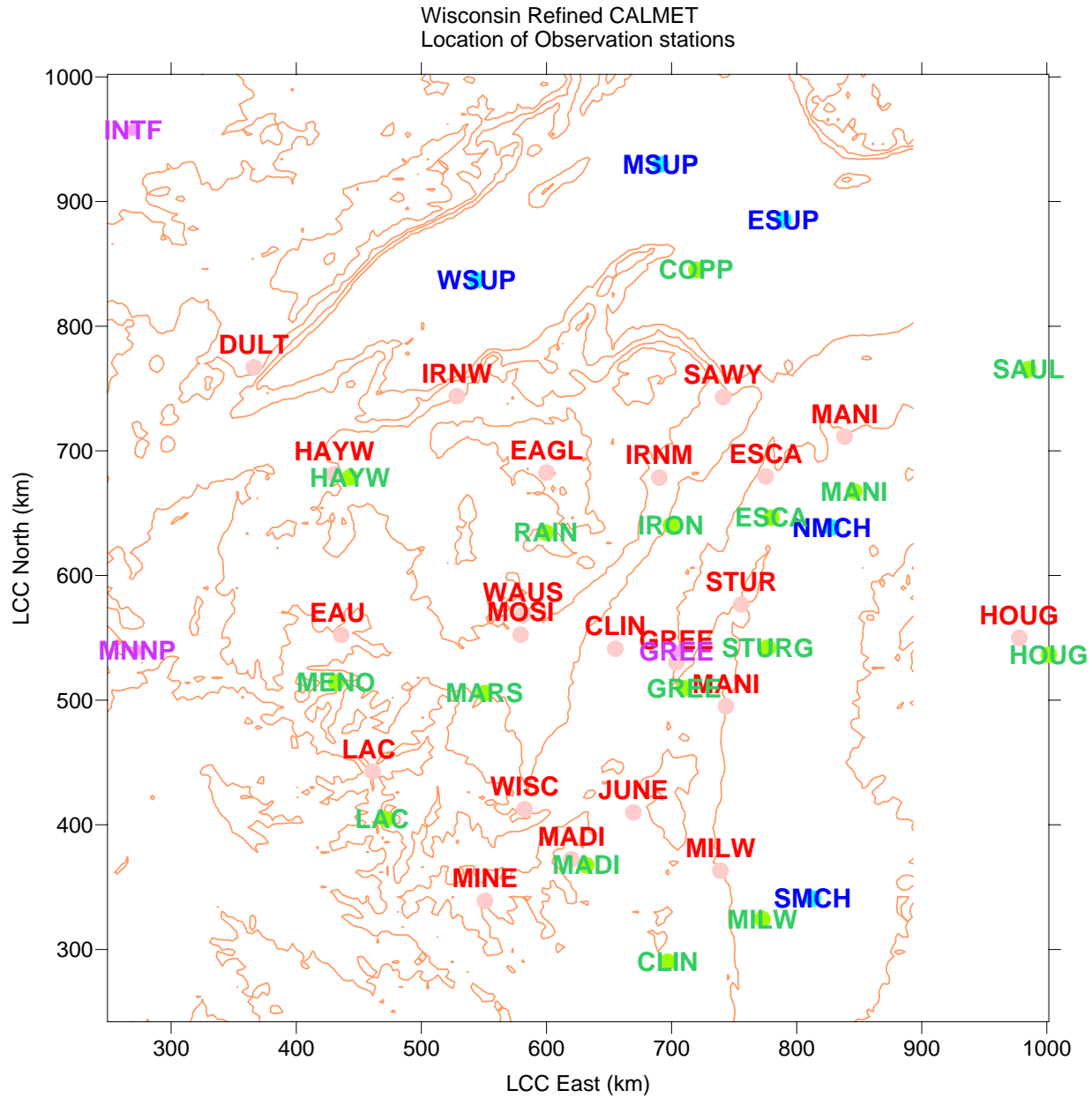
Figure 2 – Refined CALPUFF Modeling Domain and Locations of Class I Areas



NOTE: The Class I area receptors are in red, the Thilmany mill location is in blue, and the X and Y coordinates are the Lambert projection coordinates are in kms.

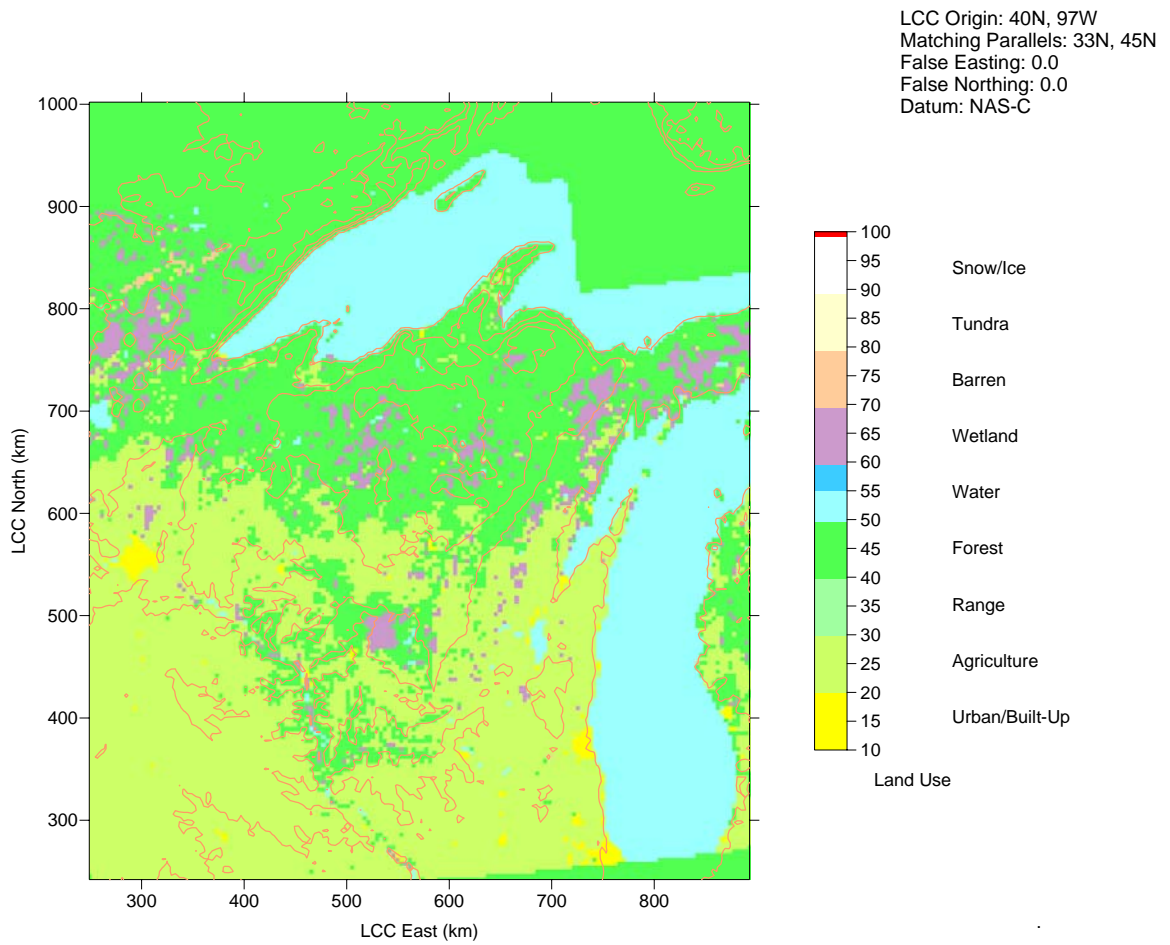


Figure 3 – Locations of Observational Stations in Refined CALPUFF Modeling Domain



NOTES: Surface stations are in red, overwater (bouy) in blue, precipitation in green, and upper air in purple.

Figure 4 – Land Use Data for Refined CALPUFF Modeling Domain



Some of the notable CALMET configuration settings included:

- CALMET vertical layers set to 10, with cell face heights at 20, 40, 80, 160, 320, 640, 1200, 2000, 3000, and 4000 meters,
- CALMET diagnostic options, including slope flow effects (ISLOPE = 1) with the radius of influence of terrain (TERRAD) set to 15 km,
- EPA recommended CALMET settings for the radius of influence parameters (R1=50, R2=100, RMAX1=100, RMAX2=200, and RMAX3=200).

## 4.0 CALPUFF Procedures

Both the DNR screening and VISTAS refined CALPUFF input files were used as the basis for the refined Wisconsin CALPUFF input files. All CALPUFF settings and options are consistent with EPA and FLM guidance, and the refined VISTAS protocol recommendations. The background ozone (O<sub>3</sub>) concentrations were used directly from the DNR CALPUFF input files. Instead of using monthly background ammonia concentrations from the DNR screening analyses, a single conservative background ammonia concentration of 3.9 ppb was used. Appendix B presents a listing of a CALPUFF input file developed for the Thilmany analysis.

The DNR screening analysis considered three emitted pollutants, SO<sub>2</sub>, NO<sub>x</sub>, and PM<sub>2.5</sub>. The refined CALPUFF modeling will consider speciated particulate emissions, including coarse PM (PMC), fine PM equivalent to PM<sub>2.5</sub> (PMF), sulfate (SO<sub>4</sub>), nitrate (NO<sub>3</sub>), elemental carbon (EC), and secondary organic aerosols (SOA or OC). Section 6 of this protocol describes how the speciated particulate emissions were calculated.

## 5.0 POSTUTIL and CALPOST Procedures

In CALPUFF, the scavenging of the available ammonia for nitrate formation is computed on a puff-by-puff basis. Because of the possibility of multiple puff overlap, this approach may tend to overestimate the amount of ammonia available to form nitrate. In order to address the cumulative consumption of ammonia, the POSTUTIL program can be used to repartition nitric acid and nitrate. POSTUTIL calculates new NO<sub>3</sub>/HNO<sub>3</sub> equilibrium concentrations for the sample source in the Class I area using the background ammonia. As recommended in the VISTAS procedures for refined CALPUFF modeling, the POSTUTIL “MNITRATE=1” method was used for the refined Wisconsin CALPUFF modeling. Instead of using the monthly values for background ammonia concentrations from the DNR screening analyses, a single conservative background ammonia concentration of 3.9 ppb was used.

The modeling protocol submitted in 2007 proposed to use CALPOST Visibility Method 6 to calculate visibility impacts. However, since that time there have been revisions to CALPOST that allow the use of the updated IMPROVE extinction equation. CALPOST version 6.221 can implement the “Method 8 Mode 5” methodology to calculate visibility impacts, and the FLMs recommend its use in “Federal Land Managers’ Air Quality Related Values Work Group (FLAG) Phase I Report—Revised (2010)” (herein referred to as FLAG 2010). Therefore, this updated methodology will be used for the Thilmany BART analysis. The new CALPUFF Graphical Users Interface automatically enters the appropriate values, based on the Class I area being selected, for natural background species and relative humidity adjustment factors using the data in Tables 6 through 9 of Section 3.3.5 of FLAG 2010 (which are based on “Guidance for Estimating Natural Visibility Conditions Under the Regional Haze Rule”, EPA-454/B-03-005, September 2003).

## 6.0 Emissions Modeled

EPA’s BART modeling guidance state that the 24-hr maximum actual emissions rates for the period of the modeling simulation (2002-2004) should be used for the analysis. If this data is not available, then the short term “allowable” or “potential” emission rates should be used. Thilmany provided DNR with source information that has been reviewed to determine the 24-hr maximum actual emissions rates. The SO<sub>2</sub> emissions are based on the maximum actual emission rate from the CEMS during the 2002-04 time frame, and the PM and NO<sub>x</sub> emissions represent maximum potential emission rates (which is conservative). On November 29, 2010, DNR issued a letter that approved the following emission rates to be used for the Thilmany BART eligible boiler: 207.9 g/sec (1650 lb/hr) SO<sub>2</sub>, 68.2 g/sec (541 lb/hr) NO<sub>x</sub>, and 11.1 g/sec (88 lb/hr) for PM. The PM emissions have been speciated using the information for coal fired boilers developed by the National Park Service<sup>1</sup>. The speciated PM emissions listed in Table 1 were used in the refined BART modeling. The stack parameters used in the modeling analysis were the same as those used by DNR in their screening analysis.

Table 1 – Thilmany Speciated PM Emissions

PM Species	Emission Rate (lb/hr)
Coarse PM	2.9
Fine PM	3.4
Elemental Carbon	0.1
Condensable Inorganics (Sulfate)	65.3
Condensable Organics (OC)	16.3

---

<sup>1</sup> <http://www2.nature.nps.gov/air/permits/ect/index.cfm>

## 7.0 Modeling Results

CALPOST computes the daily maximum change in deciviews relative to the annual average natural background visibility, separately for each Class I area. For evaluating the source impact relative to the “Subject to BART” visibility impact threshold of 0.5 dv, the daily maximum changes in visibility are sorted to determine the 98th percentile (8<sup>th</sup> highest) value for each of three years, and the 98th percentile (22<sup>nd</sup> highest) value for the completed three year period. If all 98th percentile values are below the threshold of 0.5 dv based on the refined CALPUFF analysis, the source is exempt from the BART requirements.

Another equivalent metric commonly used for BART exemption modeling is to sum the number of days per year that are above the 0.5 dv threshold for each Class I area separately; if there are 7 or fewer days above the threshold for each year, then the 98<sup>th</sup> percentile value is also below the threshold. This is the metric approved by DNR for use in the BART modeling analysis.

Specifically, for each Class I area separately, if there are less than 8 days per year that the predicted visibility impacts are greater than 0.5 deciviews relative to the annual average natural background visibility, then the BART eligible emission units at the modeled facility will be determined not to be Subject to BART.

Table 2 presents the results from the refined CALPUFF BART exemption modeling. These results demonstrate that there are less than 8 days per year, at each individual Class I area, when the visibility impacts from the Thilmany BART eligible source are above 0.5 dv. Therefore, the **Thilmany BART eligible source is not Subject to BART** requirements.

Table 2 – Thilmany CALPUFF BART Exemption Results Summary

<b>Class I area</b>	<b>Distance (km) from source to Class I area boundary</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2002</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2003</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2004</b>	<b># of days with impact &gt; 0.5 dv in Class I area for 3 year period</b>	<b>Maximum 98<sup>th</sup> Percentile impact over either annual or 3-yr periods</b>
Seney , MI	~ 280	2	0	1	3	0.34
Isle Royal, MN	~ 400	0	0	0	0	0.14
Boundary Waters, MN	~ 450	1	0	0	1	0.11

**Appendix A – Example CALMET Input File**



**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

---

WI Refined CALMET - 4km Grid - Dec 2010 - with Updated EPA Settings

----- Run title (3 lines) -----  
CALMET MODEL CONTROL FILE  
-----

INPUT GROUP: 0 -- Input and Output File Names

Subgroup (a)

Default Name	Type	File Name
-----	----	-----
GEO.DAT	input	! GEODAT=geo.dat !
SURF.DAT	input	! SRFDAT=SURF02Z.DAT !
CLOUD.DAT	input	* CLDDAT= *
PRECIP.DAT	input	! PRCDAT=PRECIP02.DAT !
MM4.DAT	input	* MM4DAT=E:\MRPO-calmm5\2002\MM5.020102.3D.DAT *
MM4.DAT	input	! MM4DAT=E:\MRPO-calmm5\2002\020102.MM5 !
WT.DAT	input	* WTDAT= *
CALMET.LST	output	! METLST=cal020102.lst !
CALMET.DAT	output	! METDAT=E:\WI-Calmet\2002Output\cal020102.dat !
PACOUT.DAT	output	* PACDAT= *

All file names will be converted to lower case if LCFILES = T  
Otherwise, if LCFILES = F, file names will be converted to UPPER CASE  
T = lower case ! LCFILES = T !  
F = UPPER CASE

NUMBER OF UPPER AIR & OVERWATER STATIONS:

Number of upper air stations (NUSTA)	No default	! NUSTA = 3 !
Number of overwater met stations	(NOWSTA) No default	! NOWSTA = 5 !

!END!

-----  
Subgroup (b)

---

-----  
Upper air files (one per station)  
-----

Default Name	Type	File Name		
UP1.DAT	input	1 ! UPDAT=UPGB0204.DAT!	!	END!
UP2.DAT	input	2 ! UPDAT=UPMN0204.DAT!	!	END!
UP3.DAT	input	3 ! UPDAT=UPIN0204.DAT!	!	END!

-----  
Subgroup (c)  
-----

Overwater station files (one per station)

SEA1.DAT	input	1 ! SEADAT=4500102.DAT!	!	END!
SEA2.DAT	input	2 ! SEADAT=4500202.DAT!	!	END!
SEA3.DAT	input	3 ! SEADAT=4500402.DAT!	!	END!
SEA4.DAT	input	4 ! SEADAT=4500602.DAT!	!	END!
SEA5.DAT	input	5 ! SEADAT=4500702.DAT!	!	END!

-----  
Default Name Type File Name  
-----  
-----

-----  
Subgroup (d)  
-----

Other file names  
-----

Default Name	Type	File Name		
DIAG.DAT	input	* DIADAT=	*	
PROG.DAT	input	* PRGDAT=	*	
TEST.PRT	output	* TSTPRT=	*	
TEST.OUT	output	* TSTOUT=	*	
TEST.KIN	output	* TSTKIN=	*	
TEST.FRD	output	* TSTFRD=	*	
TEST.SLP	output	* TSTSLP=	*	

-----  
NOTES: (1) File/path names can be up to 70 characters in length  
(2) Subgroups (a) and (d) must have ONE 'END' (surround by  
delimiters) at the end of the group  
(3) Subgroups (b) and (c) must have an 'END' (surround by

delimiters) at the end of EACH LINE

!END!

-----  
INPUT GROUP: 1 -- General run control parameters  
-----

Starting date: Year (IBYR) -- No default ! IBYR= 2002 !  
Month (IBMO) -- No default ! IBMO= 1 !  
Day (IBDY) -- No default ! IBDY= 2 !  
Hour (IBHR) -- No default ! IBHR= 0 !

Base time zone (IBTZ) -- No default ! IBTZ= 6 !  
PST = 08, MST = 07  
CST = 06, EST = 05

Length of run (hours) (IRLG) -- No default ! IRLG= 24 !

Run type (IRTYPE) -- Default: 1 ! IRTYPE= 1 !

0 = Computes wind fields only  
1 = Computes wind fields and micrometeorological variables  
(u\*, w\*, L, zi, etc.)  
(IRTYPE must be 1 to run CALPUFF or CALGRID)

Compute special data fields required  
by CALGRID (i.e., 3-D fields of W wind  
components and temperature)  
in addition to regular Default: T ! LCALGRD = T !  
fields ? (LCALGRD)  
(LCALGRD must be T to run CALGRID)

Flag to stop run after  
SETUP phase (ITEST) Default: 2 ! ITEST= 2 !  
(Used to allow checking  
of the model inputs, files, etc.)  
ITEST = 1 - STOPS program after SETUP phase  
ITEST = 2 - Continues with execution of  
COMPUTATIONAL phase after SETUP

Test options specified to see if  
they conform to regulatory

values? (MREG)                      No Default                      ! MREG = 1                      !

0 = NO checks are made

1 = Technical options must conform to USEPA guidance

          IMIXH    -1            Maul-Carson convective mixing height  
                                  over land; OCD mixing height overwater  
          ICOARE    0            OCD deltaT method for overwater fluxes  
          THRESHL  0.0           Threshold buoyancy flux over land needed  
                                  to sustain convective mixing height grow

!END!

-----

INPUT GROUP: 2 -- Map Projection and Grid control parameters

-----

Projection for all (X,Y):

-----

Map projection

(PMAP)                                      Default: UTM                      ! PMAP = LCC                      !

UTM : Universal Transverse Mercator  
TTM : Tangential Transverse Mercator  
LCC : Lambert Conformal Conic  
PS : Polar Stereographic  
EM : Equatorial Mercator  
LAZA : Lambert Azimuthal Equal Area

False Easting and Northing (km) at the projection origin

(Used only if PMAP= TTM, LCC, or LAZA)

(FEAST)                                      Default=0.0                      ! FEAST = 0.000                      !  
(FNORTH)                                     Default=0.0                      ! FNORTH = 0.000                      !

UTM zone (1 to 60)

(Used only if PMAP=UTM)

(IUTMZN)                                     No Default                      ! IUTMZN = -999                      !

Hemisphere for UTM projection?

(Used only if PMAP=UTM)

(UTMHEM)                                     Default: N                      ! UTMHEM = N                      !

N : Northern hemisphere projection  
S : Southern hemisphere projection

Latitude and Longitude (decimal degrees) of projection origin

(Used only if PMAP= TTM, LCC, PS, EM, or LAZA)

(RLAT0)                      No Default              ! RLAT0 = 40N    !

(RLON0)                      No Default              ! RLON0 = 97W    !

TTM : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience

LCC : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience

PS : RLON0 identifies central (grid N/S) meridian of projection  
      RLAT0 selected for convenience

EM : RLON0 identifies central meridian of projection  
      RLAT0 is REPLACED by 0.0N (Equator)

LAZA: RLON0 identifies longitude of tangent-point of mapping plane  
      RLAT0 identifies latitude of tangent-point of mapping plane

Matching parallel(s) of latitude (decimal degrees) for projection

(Used only if PMAP= LCC or PS)

(XLAT1)                      No Default              ! XLAT1 = 33N    !

(XLAT2)                      No Default              ! XLAT2 = 45N    !

LCC : Projection cone slices through Earth's surface at XLAT1 and XLAT2

PS : Projection plane slices through Earth at XLAT1  
      (XLAT2 is not used)

-----

Note: Latitudes and longitudes should be positive, and include a  
      letter N,S,E, or W indicating north or south latitude, and  
      east or west longitude. For example,  
      35.9 N Latitude = 35.9N  
      118.7 E Longitude = 118.7E

Datum-region

-----

The Datum-Region for the coordinates is identified by a character  
string. Many mapping products currently available use the model of the  
Earth known as the World Geodetic System 1984 (WGS-84). Other local  
models may be in use, and their selection in CALMET will make its output  
consistent with local mapping products. The list of Datum-Regions with  
official transformation parameters is provided by the National Imagery and  
Mapping Agency (NIMA).

NIMA Datum - Regions(Examples)

```

-----
WGS-84   WGS-84 Reference Ellipsoid and Geoid, Global coverage (WGS84)
NAS-C    NORTH AMERICAN 1927 Clarke 1866 Spheroid, MEAN FOR CONUS (NAD27)
NAR-C    NORTH AMERICAN 1983 GRS 80 Spheroid, MEAN FOR CONUS (NAD83)
NWS-84   NWS 6370KM Radius, Sphere
ESR-S    ESRI REFERENCE 6371KM Radius, Sphere
  
```

Datum-region for output coordinates

```

(DATUM)                Default: WGS-84   ! DATUM = NAS-C   !
  
```

Horizontal grid definition:

```

-----

Rectangular grid defined for projection PMAP,
with X the Easting and Y the Northing coordinate
  
```

```

         No. X grid cells (NX)      No default   ! NX = 162  !
         No. Y grid cells (NY)      No default   ! NY = 191  !

Grid spacing (DGRIDKM)              No default   ! DGRIDKM = 4.  !
                                     Units: km
  
```

Reference grid coordinate of  
SOUTHWEST corner of grid cell (1,1)

```

         X coordinate (XORIGKM)      No default   ! XORIGKM = 247.000  !
         Y coordinate (YORIGKM)      No default   ! YORIGKM = 240.000  !
                                     Units: km
  
```

Vertical grid definition:

```

-----

No. of vertical layers (NZ)         No default   ! NZ = 10  !

Cell face heights in arbitrary
vertical grid (ZFACE(NZ+1))         No defaults
                                     Units: m
! ZFACE = 0.,20.,40.,80.,160.,320.,640.,1200.,2000.,3000.,4000.  !
  
```

!END!

-----  
INPUT GROUP: 3 -- Output Options  
-----

DISK OUTPUT OPTION

Save met. fields in an unformatted  
output file ? (LSAVE) Default: T ! LSAVE = T !  
(F = Do not save, T = Save)

Type of unformatted output file:  
(IFORMO) Default: 1 ! IFORMO = 1 !

- 1 = CALPUFF/CALGRID type file (CALMET.DAT)
- 2 = MESOPUFF-II type file (PACOUT.DAT)

LINE PRINTER OUTPUT OPTIONS:

Print met. fields ? (LPRINT) Default: F ! LPRINT = T !  
(F = Do not print, T = Print)  
(NOTE: parameters below control which  
met. variables are printed)

Print interval  
(IPRINF) in hours Default: 1 ! IPRINF = 1 !  
(Meteorological fields are printed  
every 1 hours)

Specify which layers of U, V wind component  
to print (IUVOUT(NZ)) -- NOTE: NZ values must be entered  
(0=Do not print, 1=Print)  
(used only if LPRINT=T) Defaults: NZ\*0  
! IUVOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !  
-----

Specify which levels of the W wind component to print  
(NOTE: W defined at TOP cell face -- 16 values)  
(IWOUT(NZ)) -- NOTE: NZ values must be entered

---

(0=Do not print, 1=Print)  
(used only if LPRINT=T & LCALGRD=T)  
-----  
Defaults: NZ\*0  
! IWOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Specify which levels of the 3-D temperature field to print  
(ITOUT(NZ)) -- NOTE: NZ values must be entered  
(0=Do not print, 1=Print)  
(used only if LPRINT=T & LCALGRD=T)  
-----  
Defaults: NZ\*0  
! ITOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Specify which meteorological fields  
to print  
(used only if LPRINT=T) Defaults: 0 (all variables)  
-----

Variable	Print ?	
	(0 = do not print,	
	1 = print)	
-----	-----	
! STABILITY =	0	! - PGT stability class
! USTAR =	0	! - Friction velocity
! MONIN =	0	! - Monin-Obukhov length
! MIXHT =	0	! - Mixing height
! WSTAR =	0	! - Convective velocity scale
! PRECIP =	0	! - Precipitation rate
! SENSHEAT =	0	! - Sensible heat flux
! CONVZI =	0	! - Convective mixing ht.

Testing and debug print options for micrometeorological module

Print input meteorological data and  
internal variables (LDB) Default: F ! LDB = F !  
(F = Do not print, T = print)  
(NOTE: this option produces large amounts of output)

First time step for which debug data



are printed (NN1)                                  Default: 1                                  ! NN1 = 1 !

Last time step for which debug data  
are printed (NN2)                                  Default: 1                                  ! NN2 = 2 !

Testing and debug print options for wind field module  
(all of the following print options control output to  
wind field module's output files: TEST.PRT, TEST.OUT,  
TEST.KIN, TEST.FRD, and TEST.SLP)

Control variable for writing the test/debug  
wind fields to disk files (IOUTD)  
(0=Do not write, 1=write)                                  Default: 0                                  ! IOUTD = 0 !

Number of levels, starting at the surface,  
to print (NZPRN2)                                  Default: 1                                  ! NZPRN2 = 1 !

Print the INTERPOLATED wind components ?  
(IPR0) (0=no, 1=yes)                                  Default: 0                                  ! IPR0 = 0 !

Print the TERRAIN ADJUSTED surface wind  
components ?  
(IPR1) (0=no, 1=yes)                                  Default: 0                                  ! IPR1 = 0 !

Print the SMOOTHED wind components and  
the INITIAL DIVERGENCE fields ?  
(IPR2) (0=no, 1=yes)                                  Default: 0                                  ! IPR2 = 0 !

Print the FINAL wind speed and direction  
fields ?  
(IPR3) (0=no, 1=yes)                                  Default: 0                                  ! IPR3 = 0 !

Print the FINAL DIVERGENCE fields ?  
(IPR4) (0=no, 1=yes)                                  Default: 0                                  ! IPR4 = 0 !

Print the winds after KINEMATIC effects  
are added ?  
(IPR5) (0=no, 1=yes)                                  Default: 0                                  ! IPR5 = 0 !

Print the winds after the FROUDE NUMBER  
adjustment is made ?  
(IPR6) (0=no, 1=yes)                                  Default: 0                                  ! IPR6 = 0 !

Print the winds after SLOPE FLOWS  
are added ?

(IPR7) (0=no, 1=yes)                      Default: 0                      ! IPR7 = 0 !

Print the FINAL wind field components ?

(IPR8) (0=no, 1=yes)                      Default: 0                      ! IPR8 = 0 !

!END!

-----  
INPUT GROUP: 4 -- Meteorological data options  
-----

NO OBSERVATION MODE                      (NOOBS) Default: 0                      ! NOOBS = 0 !

0 = Use surface, overwater, and upper air stations

1 = Use surface and overwater stations (no upper air observations)  
    Use MM5 for upper air data

2 = No surface, overwater, or upper air observations  
    Use MM5 for surface, overwater, and upper air data

NUMBER OF SURFACE & PRECIP. METEOROLOGICAL STATIONS

Number of surface stations    (NSSTA) No default                      ! NSSTA = 22 !

Number of precipitation stations

(NPSTA=-1: flag for use of MM5 precip data)

(NPSTA) No default                      ! NPSTA = 16 !

CLOUD DATA OPTIONS

Gridded cloud fields:

(ICLOUD) Default: 0                      ! ICLOUD = 0 !

ICLOUD = 0 - Gridded clouds not used

ICLOUD = 1 - Gridded CLOUD.DAT generated as OUTPUT

ICLOUD = 2 - Gridded CLOUD.DAT read as INPUT

ICLOUD = 3 - Gridded cloud cover from Prognostic Rel. Humidity

FILE FORMATS

Surface meteorological data file format

(IFORMS) Default: 2                      ! IFORMS = 2 !

(1 = unformatted (e.g., SMERGE output))

(2 = formatted    (free-formatted user input))

Precipitation data file format

(IFORMP) Default: 2 ! IFORMP = 2 !

(1 = unformatted (e.g., PMERGE output))

(2 = formatted (free-formatted user input))

Cloud data file format

(IFORMC) Default: 2 ! IFORMC = 2 !

(1 = unformatted - CALMET unformatted output)

(2 = formatted - free-formatted CALMET output or user input)

!END!

-----  
INPUT GROUP: 5 -- Wind Field Options and Parameters  
-----

WIND FIELD MODEL OPTIONS

Model selection variable (IWFCOD) Default: 1 ! IWFCOD = 1 !

0 = Objective analysis only

1 = Diagnostic wind module

Compute Froude number adjustment

effects ? (IFRADJ) Default: 1 ! IFRADJ = 1 !

(0 = NO, 1 = YES)

Compute kinematic effects ? (IKINE) Default: 0 ! IKINE = 0 !

(0 = NO, 1 = YES)

Use O'Brien procedure for adjustment

of the vertical velocity ? (IOBR) Default: 0 ! IOBR = 0 !

(0 = NO, 1 = YES)

Compute slope flow effects ? (ISLOPE) Default: 1 ! ISLOPE = 1 !

(0 = NO, 1 = YES)

Extrapolate surface wind observations

to upper layers ? (IEXTRP) Default: -4 ! IEXTRP = -4 !

(1 = no extrapolation is done,

2 = power law extrapolation used,

3 = user input multiplicative factors

for layers 2 - NZ used (see FEXTRP array)  
4 = similarity theory used  
-1, -2, -3, -4 = same as above except layer 1 data  
at upper air stations are ignored

Extrapolate surface winds even

if calm? (ICALM) Default: 0 ! ICALM = 0 !  
(0 = NO, 1 = YES)

Layer-dependent biases modifying the weights of  
surface and upper air stations (BIAS(NZ))

-1<=BIAS<=1

Negative BIAS reduces the weight of upper air stations

(e.g. BIAS=-0.1 reduces the weight of upper air stations  
by 10%; BIAS= -1, reduces their weight by 100 %)

Positive BIAS reduces the weight of surface stations

(e.g. BIAS= 0.2 reduces the weight of surface stations  
by 20%; BIAS=1 reduces their weight by 100%)

Zero BIAS leaves weights unchanged (1/R\*\*2 interpolation)

Default: NZ\*0

! BIAS = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Minimum distance from nearest upper air station  
to surface station for which extrapolation  
of surface winds at surface station will be allowed  
(RMIN2: Set to -1 for IEXTRP = 4 or other situations  
where all surface stations should be extrapolated)

Default: 4. ! RMIN2 = -1.0 !

Use gridded prognostic wind field model  
output fields as input to the diagnostic

wind field model (IPROG) Default: 0 ! IPROG = 14 !  
(0 = No, [IWFCOD = 0 or 1])

1 = Yes, use CSUMM prog. winds as Step 1 field, [IWFCOD = 0]

2 = Yes, use CSUMM prog. winds as initial guess field [IWFCOD = 1]

3 = Yes, use winds from MM4.DAT file as Step 1 field [IWFCOD = 0]

4 = Yes, use winds from MM4.DAT file as initial guess field [IWFCOD = 1]

5 = Yes, use winds from MM4.DAT file as observations [IWFCOD = 1]

13 = Yes, use winds from MM5.DAT file as Step 1 field [IWFCOD = 0]

14 = Yes, use winds from MM5.DAT file as initial guess field [IWFCOD = 1]

15 = Yes, use winds from MM5.DAT file as observations [IWFCOD = 1]

Timestep (hours) of the prognostic

model input data (ISTEPPG) Default: 1 ! ISTEPPG = 1 !

---

RADIUS OF INFLUENCE PARAMETERS

Use varying radius of influence           Default: F           ! LVARY = F!  
(if no stations are found within RMAX1,RMAX2,  
  or RMAX3, then the closest station will be used)

Maximum radius of influence over land  
in the surface layer (RMAX1)           No default           ! RMAX1 = 100. !  
  Units: km

Maximum radius of influence over land  
aloft (RMAX2)                          No default           ! RMAX2 = 200. !  
  Units: km

Maximum radius of influence over water  
(RMAX3)                                No default           ! RMAX3 = 200. !  
  Units: km

OTHER WIND FIELD INPUT PARAMETERS

Minimum radius of influence used in  
the wind field interpolation (RMIN)    Default: 0.1       ! RMIN = 0.1 !  
  Units: km

Radius of influence of terrain  
features (TERRAD)                      No default       ! TERRAD = 15. !  
  Units: km

Relative weighting of the first  
guess field and observations in the  
SURFACE layer (R1)                    No default       ! R1 = 50. !  
(R1 is the distance from an           Units: km  
observational station at which the  
observation and first guess field are  
equally weighted)

Relative weighting of the first  
guess field and observations in the  
layers ALOFT (R2)                      No default       ! R2 =100. !  
(R2 is applied in the upper layers    Units: km  
in the same manner as R1 is used in  
the surface layer).

Relative weighting parameter of the  
prognostic wind field data (RPROG)   No default       ! RPROG = 0. !

(Used only if IPROG = 1) Units: km

-----

Maximum acceptable divergence in the  
divergence minimization procedure

(DIVLIM) Default: 5.E-6 ! DIVLIM= 5.0E-06 !

Maximum number of iterations in the  
divergence min. procedure (NITER)

Default: 50 ! NITER = 50 !

Number of passes in the smoothing  
procedure (NSMTH(NZ))

NOTE: NZ values must be entered

Default: 2,(mxnz-1)\*4 ! NSMTH =

2 , 4 , 4 , 4 , 4 , 4 , 4 , 4 , 4 , 4 , 4 !

Maximum number of stations used in  
each layer for the interpolation of  
data to a grid point (NINTR2(NZ))

NOTE: NZ values must be entered Default: 99. ! NINTR2 =

99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 !

Critical Froude number (CRITFN)

Default: 1.0 ! CRITFN = 1. !

Empirical factor controlling the  
influence of kinematic effects

(ALPHA) Default: 0.1 ! ALPHA = 0.1 !

Multiplicative scaling factor for  
extrapolation of surface observations

to upper layers (FEXTR2(NZ)) Default: NZ\*0.0

! FEXTR2 = 0., 0., 0., 0., 0., 0., 0., 0., 0., 0. !

(Used only if IEXTRP = 3 or -3)

#### BARRIER INFORMATION

Number of barriers to interpolation  
of the wind fields (NBAR)

Default: 0 ! NBAR = 0 !

THE FOLLOWING 4 VARIABLES ARE INCLUDED

ONLY IF NBAR > 0

NOTE: NBAR values must be entered No defaults

for each variable Units: km

X coordinate of BEGINNING  
of each barrier (XBBAR(NBAR))       ! XBBAR = 0. !  
Y coordinate of BEGINNING  
of each barrier (YBBAR(NBAR))       ! YBBAR = 0. !  
  
X coordinate of ENDING  
of each barrier (XEBAR(NBAR))       ! XEBAR = 0. !  
Y coordinate of ENDING  
of each barrier (YEBAR(NBAR))       ! YEBAR = 0. !

DIAGNOSTIC MODULE DATA INPUT OPTIONS

Surface temperature (IDIOPT1)       Default: 0       ! IDIOPT1 = 0 !  
0 = Compute internally from  
    hourly surface observations  
1 = Read preprocessed values from  
    a data file (DIAG.DAT)

Surface met. station to use for  
the surface temperature (ISURFT)   No default       ! ISURFT = 1 !  
(Must be a value from 1 to NSSTA)  
(Used only if IDIOPT1 = 0)

-----

Domain-averaged temperature lapse  
rate (IDIOPT2)                      Default: 0       ! IDIOPT2 = 0 !  
0 = Compute internally from  
    twice-daily upper air observations  
1 = Read hourly preprocessed values  
    from a data file (DIAG.DAT)

Upper air station to use for  
the domain-scale lapse rate (IUPT) No default       ! IUPT = 1 !  
(Must be a value from 1 to NUSTA)  
(Used only if IDIOPT2 = 0)

-----

Depth through which the domain-scale  
lapse rate is computed (ZUPT)       Default: 200.   ! ZUPT = 200. !  
(Used only if IDIOPT2 = 0)       Units: meters

-----

Domain-averaged wind components

(IDIOPT3) Default: 0 ! IDIOPT3 = 0 !

0 = Compute internally from  
twice-daily upper air observations  
1 = Read hourly preprocessed values  
a data file (DIAG.DAT)

Upper air station to use for

the domain-scale winds (IUPWND) Default: -1 ! IUPWND = -1 !

(Must be a value from -1 to NUSTA)

(Used only if IDIOPT3 = 0)

-----

Bottom and top of layer through

which the domain-scale winds

are computed

(ZUPWND(1), ZUPWND(2)) Defaults: 1., 1000. ! ZUPWND= 1., 1000. !

(Used only if IDIOPT3 = 0) Units: meters

-----

Observed surface wind components

for wind field module (IDIOPT4) Default: 0 ! IDIOPT4 = 0 !

0 = Read WS, WD from a surface  
data file (SURF.DAT)

1 = Read hourly preprocessed U, V from  
a data file (DIAG.DAT)

Observed upper air wind components

for wind field module (IDIOPT5) Default: 0 ! IDIOPT5 = 0 !

0 = Read WS, WD from an upper  
air data file (UP1.DAT, UP2.DAT, etc.)

1 = Read hourly preprocessed U, V from  
a data file (DIAG.DAT)

LAKE BREEZE INFORMATION

Use Lake Breeze Module (LLBREZE)

Default: F ! LLBREZE = F !

Number of lake breeze regions (NBOX) ! NBOX = 0 !

X Grid line 1 defining the region of interest

! XG1 = 0. !

X Grid line 2 defining the region of interest



**Refined BART Modeling Protocol for  
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```

! XG2 = 0. !
Y Grid line 1 defining the region of interest
! YG1 = 0. !
Y Grid line 2 defining the region of interest
! YG2 = 0. !

X Point defining the coastline (Straight line)
(XBCST) (KM) Default: none ! XBCST = 0. !

Y Point defining the coastline (Straight line)
(YBCST) (KM) Default: none ! YBCST = 0. !

X Point defining the coastline (Straight line)
(XECST) (KM) Default: none ! XECST = 0. !

Y Point defining the coastline (Straight line)
(YECST) (KM) Default: none ! YECST = 0. !

Number of stations in the region Default: none ! NLB = 0 !
(Surface stations + upper air stations)

Station ID's in the region (METBXID(NLB))
(Surface stations first, then upper air stations)
! METBXID = 0 !

!END!

-----

INPUT GROUP: 6 -- Mixing Height, Temperature and Precipitation Parameters
-----

EMPIRICAL MIXING HEIGHT CONSTANTS

Neutral, mechanical equation
(CONSTB) Default: 1.41 ! CONSTB = 1.41 !
Convective mixing ht. equation
(CONSTE) Default: 0.15 ! CONSTE = 0.15 !
Stable mixing ht. equation
(CONSTN) Default: 2400. ! CONSTN = 2400.!
Overwater mixing ht. equation
(CONSTW) Default: 0.16 ! CONSTW = 0.16 !
```

---

Absolute value of Coriolis  
parameter (FCORIOI)                      Default: 1.E-4   ! FCORIOI = 1.0E-04!  
Units: (1/s)

SPATIAL AVERAGING OF MIXING HEIGHTS

Conduct spatial averaging  
(IAVEZI) (0=no, 1=yes)                    Default: 1       ! IAVEZI = 1 !

Max. search radius in averaging  
process (MNMDAV)                            Default: 1       ! MNMDAV = 1 !  
Units: Grid  
      cells

Half-angle of upwind looking cone  
for averaging (HAFANG)                     Default: 30.     ! HAFANG = 30. !  
Units: deg.

Layer of winds used in upwind  
averaging (ILEVZI)                         Default: 1       ! ILEVZI = 1 !  
(must be between 1 and NZ)

CONVECTIVE MIXING HEIGHT OPTIONS:

Method to compute the convective  
mixing height(IMIXH)                        Default: 1       ! IMIXH = -1 !

- 1: Maul-Carson for land and water cells
- 1: Maul-Carson for land cells only -  
    OCD mixing height overwater
- 2: Batchvarova and Gryning for land and water cells
- 2: Batchvarova and Gryning for land cells only  
    OCD mixing height overwater

Threshold buoyancy flux required to  
sustain convective mixing height growth  
overland (THRESHL)                         Default: 0.05   ! THRESHL = 0.00 !  
(expressed as a heat flux                 units: W/m3  
per meter of boundary layer)

Threshold buoyancy flux required to  
sustain convective mixing height growth  
overwater (THRESHW)                        Default: 0.05   ! THRESHW = 0.05 !  
(expressed as a heat flux                 units: W/m3  
per meter of boundary layer)

Option for overwater lapse rates used  
in convective mixing height growth

(ITWPROG) Default: 0 ! ITWPROG = 0 !  
0 : use SEA.DAT lapse rates and deltaT (or assume neutral  
conditions if missing)  
1 : use prognostic lapse rates (only if IPROG>2)  
and SEA.DAT deltaT (or neutral if missing)  
2 : use prognostic lapse rates and prognostic delta T  
(only if iprog>12 and 3D.DAT version# 2.0 or higher)

Land Use category ocean in 3D.DAT datasets

(ILUOC3D) Default: 16 ! ILUOC3D = 16 !  
Note: if 3D.DAT from MM5 version 3.0, iluoc3d = 16  
if MM4.DAT, typically iluoc3d = 7

#### OTHER MIXING HEIGHT VARIABLES

Minimum potential temperature lapse  
rate in the stable layer above the  
current convective mixing ht. Default: 0.001 ! DPTMIN = 0.001 !  
(DPTMIN) Units: deg. K/m

Depth of layer above current conv.  
mixing height through which lapse  
rate is computed (DZZI) Default: 200. ! DZZI = 200. !  
Units: meters

Minimum overland mixing height Default: 50. ! ZIMIN = 50. !  
(ZIMIN) Units: meters

Maximum overland mixing height Default: 3000. ! ZIMAX = 3000. !  
(ZIMAX) Units: meters

Minimum overwater mixing height Default: 50. ! ZIMINW = 50. !  
(ZIMINW) -- (Not used if observed  
overwater mixing hts. are used) Units: meters

Maximum overwater mixing height Default: 3000. ! ZIMAXW = 3000. !  
(ZIMAXW) -- (Not used if observed  
overwater mixing hts. are used) Units: meters

#### OVERWATER SURFACE FLUXES METHOD and PARAMETERS

(ICOARE) Default: 10 ! ICOARE = 0 !  
0: original deltaT method (OCD)  
10: COARE with no wave parameterization (jwave=0, Charnock)  
11: COARE with wave option jwave=1 (Oost et al.)  
and default wave properties  
-11: COARE with wave option jwave=1 (Oost et al.)  
and observed wave properties (must be in SEA.DAT files)  
12: COARE with wave option 2 (Taylor and Yelland)  
and default wave properties

-12: COARE with wave option 2 (Taylor and Yelland)  
and observed wave properties (must be in SEA.DAT files)

Coastal/Shallow water length scale (DSHELF)  
(for modified z0 in shallow water)  
( COARE fluxes only)

Default : 0.            ! DSHELF = 0. !  
units: km

COARE warm layer computation (IWARM)            ! IWARM = 0 !  
1: on - 0: off (must be off if SST measured with  
IR radiometer)            Default: 0

COARE cool skin layer computation (ICOOL)            ! ICOOL = 0 !  
1: on - 0: off (must be off if SST measured with  
IR radiometer)            Default: 0

TEMPERATURE PARAMETERS

3D temperature from observations or  
from prognostic data? (ITPROG)            Default:0            !ITPROG = 0 !

- 0 = Use Surface and upper air stations  
(only if NOOBS = 0)
- 1 = Use Surface stations (no upper air observations)  
Use MM5 for upper air data  
(only if NOOBS = 0,1)
- 2 = No surface or upper air observations  
Use MM5 for surface and upper air data  
(only if NOOBS = 0,1,2)

Interpolation type  
(1 = 1/R ; 2 = 1/R\*\*2)            Default:1            ! IRAD = 1 !

Radius of influence for temperature  
interpolation (TRADKM)            Default: 500.            ! TRADKM = 500. !  
Units: km

Maximum Number of stations to include  
in temperature interpolation (NUMTS)            Default: 5            ! NUMTS = 5 !

Conduct spatial averaging of temp-  
eratures (IAVET) (0=no, 1=yes)            Default: 1            ! IAVET = 1 !

(will use mixing ht MNMDAV,HAFANG  
so make sure they are correct)

Default temperature gradient            Default: -.0098 ! TGDEFB = -0.0098 !  
below the mixing height over  
water (K/m) (TGDEFB)

Default temperature gradient            Default: -.0045 ! TGDEFA = -0.0045 !  
above the mixing height over  
water (K/m) (TGDEFA)

Beginning (JWAT1) and ending (JWAT2)  
land use categories for temperature            ! JWAT1 = 55 !  
interpolation over water -- Make            ! JWAT2 = 55 !  
bigger than largest land use to disable

PRECIP INTERPOLATION PARAMETERS

Method of interpolation (NFLAGP)            Default = 2            ! NFLAGP = 2 !  
(1=1/R,2=1/R\*\*2,3=EXP/R\*\*2)  
Radius of Influence (km) (SIGMAP)          Default = 100.0        ! SIGMAP = 100. !  
(0.0 => use half dist. btwn  
nearest stns w & w/out  
precip when NFLAGP = 3)  
Minimum Precip. Rate Cutoff (mm/hr)        Default = 0.01        ! CUTP = 0.01 !  
(values < CUTP = 0.0 mm/hr)

!END!

-----  
INPUT GROUP: 7 -- Surface meteorological station parameters  
-----

SURFACE STATION VARIABLES

(One record per station -- 0 records in all)

1	2					
Name	ID	X coord.	Y coord.	Time	Anem.	
		(km)	(km)	zone	Ht. (m)	
! SS1	= 'MILW'	26400	739.061	363.200	6	10 !
! SS2	= 'MADI'	26410	619.927	372.210	6	10 !

**Refined BART Modeling Protocol for  
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```

! SS3  ='LA C'   26430   461.012   442.705   6   10  !
! SS4  ='EAU '   26435   435.857   552.352   6   10  !
! SS5  ='GREE'   26450   703.911   530.179   6   10  !
! SS6  ='MANI'   26455   743.477   495.113   6   10  !
! SS7  ='STUR'   26458   755.730   576.760   6   10  !
! SS8  ='WAUS'   26463   581.202   568.247   6   10  !
! SS9  ='MOSI'   26465   579.307   552.489   6   10  !
! SS10 ='CLIN'   26502   655.019   541.168   6   10  !
! SS11 ='WISC'   26503   582.532   412.498   6   10  !
! SS12 ='EAGL'   26504   599.980   682.545   6   10  !
! SS13 ='MINE'   26507   550.817   338.941   6   10  !
! SS14 ='HAYW'   26508   430.347   681.295   6   10  !
! SS15 ='JUNE'   26509   669.429   409.931   6   10  !
! SS16 ='dulu'   27450   366.011   766.977   6   10  !
! SS17 ='MANI'   25408   838.315   711.433   5   10  !
! SS18 ='SAWY'   26284   741.237   743.078   5   10  !
! SS19 ='HOUG'   26380   977.905   549.977   5   10  !
! SS20 ='ESCA'   26480   775.417   679.550   5   10  !
! SS21 ='IRNM'   27437   690.168   678.511   5   10  !
! SS22 ='IRNW'   27445   528.314   743.857   5   10  !

```

```

-----
1
    Four character string for station name
    (MUST START IN COLUMN 9)

```

```

2
    Five digit integer for station ID

```

!END!

```

-----
INPUT GROUP: 8 -- Upper air meteorological station parameters
-----

```

```

UPPER AIR STATION VARIABLES
(One record per station -- 0 records in all)

```

```

      1      2
      Name   ID      X coord.  Y coord.  Time zone
              (km)    (km)
-----
! US1  ='GREE'  14898   703.911   539.179   6   !

```

Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill

---

```
! US2 = 'MNNP'   94983   272.681   539.843   6   !
! US3 = 'INTF'   14918   269.489   957.416   6   !
```

-----

1  
Four character string for station name  
(MUST START IN COLUMN 9)

2  
Five digit integer for station ID

!END!

-----

INPUT GROUP: 9 -- Precipitation station parameters

-----

PRECIPITATION STATION VARIABLES  
(One record per station -- 0 records in all)  
(NOT INCLUDED IF NPSTA = 0)

	1	2		
	Name	Station	X coord.	Y coord.
		Code	(km)	(km)
! PS1	= 'COPP'	201780	719.681	845.179 !
! PS2	= 'ESCA'	202626	779.850	646.454 !
! PS3	= 'HOUG'	203936	1001.500	536.427 !
! PS4	= 'IRON'	204090	699.421	640.283 !
! PS5	= 'MANI'	205073	845.933	667.499 !
! PS6	= 'SAUL'	207366	985.172	765.448 !
! PS7	= 'CLIN'	471667	696.889	290.151 !
! PS8	= 'GREE'	473269	709.933	509.579 !
! PS9	= 'HAYW'	473511	442.947	678.728 !
! PS10	= 'LAC '	474370	471.515	404.495 !
! PS11	= 'MADI'	474961	631.719	367.656 !
! PS12	= 'MARS'	475120	550.348	505.662 !
! PS13	= 'MENO'	475335	431.807	514.277 !
! PS14	= 'MILW'	475479	772.567	324.344 !
! PS15	= 'RAIN'	476939	599.369	634.460 !
! PS16	= 'STURG'	478267	776.964	542.174 !

-----  
1

Four character string for station name  
(MUST START IN COLUMN 9)

2

Six digit station code composed of state  
code (first 2 digits) and station ID (last  
4 digits)

!END!



**Appendix B – Example CALPUFF Input File**

**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

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Thilmany BART Modeling - SO2, PM speciation  
Using Refined 4km CALMET Data File and SENE, ISLE, BOWA Class I Receptors

----- Run title (3 lines) -----

CALPUFF MODEL CONTROL FILE  
-----

INPUT GROUP: 0 -- Input and Output File Names

-----  
Default Name    Type            File Name  
-----  
CALMET.DAT    input        \* METDAT =            \*  
          or  
ISCMET.DAT    input        \* ISCDAT =            \*  
          or  
PLMMET.DAT    input        \* PLMDAT =            \*  
          or  
PROFILE.DAT    input        \* PRFDAT =            \*  
SURFACE.DAT    input        \* SFCDAT =            \*  
RESTARTB.DAT  input        \* RSTARTB=            \*

-----  
CALPUFF.LST    output      ! PUFLST =ThilRef02.lst !  
CONC.DAT        output      ! CONDAT =ThilRef02.dat !  
DFLX.DAT        output      \* DFDAT =              \*  
WFLX.DAT        output      \* WFDAT =              \*  
VISB.DAT        output      \* VISDAT =Thil02DNR.VIB \*  
RESTARTE.DAT  output      \* RSTARTE=            \*

-----  
Emission Files  
-----

PTEMARB.DAT    input        \* PTDAT =              \*  
VOLEMARB.DAT  input        \* VOLDAT =              \*  
BAEMARB.DAT    input        \* ARDAT =              \*  
LNEMARB.DAT    input        \* LNDAT =              \*

-----  
Other Files  
-----

OZONE.DAT      input        \* OZDAT =              \*  
VD.DAT          input        \* VDDAT =              \*  
CHEM.DAT        input        \* CHEMDAT=            \*  
H2O2.DAT        input        \* H2O2DAT=            \*  
HILL.DAT        input        \* HILDAT=              \*  
HILLRCT.DAT    input        \* RCTDAT=              \*  
COASTLN.DAT    input        \* CSTDAT=              \*  
FLUXBDY.DAT    input        \* BDYDAT=              \*  
BCON.DAT        input        \* BCNDAT=              \*  
DEBUG.DAT        output      \* DEBUG =              \*  
MASSFLX.DAT    output      \* FLXDAT=              \*  
MASSBAL.DAT    output      \* BALDAT=              \*  
FOG.DAT        output      \* FOGDAT=              \*

-----  
All file names will be converted to lower case if LCFILES = T  
Otherwise, if LCFILES = F, file names will be converted to UPPER CASE  
      T = lower case            ! LCFILES = T !  
      F = UPPER CASE

NOTE: (1) file/path names can be up to 70 characters in length

Provision for multiple input files  
-----

**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

---

Number of CALMET.DAT files for run (NMETDAT)  
Default: 1 ! NMETDAT = 364 !

Number of PTEMARB.DAT files for run (NPTDAT)  
Default: 0 ! NPTDAT = 0 !

Number of BAEMARB.DAT files for run (NARDAT)  
Default: 0 ! NARDAT = 0 !

Number of VOLEMARB.DAT files for run (NVOLDAT)  
Default: 0 ! NVOLDAT = 0 !

!END!

-----  
Subgroup (0a)  
-----

The following CALMET.DAT filenames are processed in sequence if NMETDAT>1

Default Name	Type	File Name	
none	input	! METDAT=e:\WI-Calmet\2002output\cal020102.DAT	! !END!
NOTE: the remaining CALMET data files names are listed here ....			
none	input	! METDAT=e:\WI-Calmet\2002output\cal021231.DAT	! !END!

-----  
INPUT GROUP: 1 -- General run control parameters  
-----

Option to run all periods found  
in the met. file (METRUN) Default: 0 ! METRUN = 0 !

METRUN = 0 - Run period explicitly defined below  
METRUN = 1 - Run all periods in met. file

Starting date: Year (IBYR) -- No default ! IBYR = 2002 !  
(used only if Month (IBMO) -- No default ! IBMO = 1 !  
METRUN = 0) Day (IBDY) -- No default ! IDBY = 02 !  
Hour (IBHR) -- No default ! IBHR = 1 !

Base time zone (XBTZ) -- No default ! XBTZ = 6.0 !  
PST = 8., MST = 7.  
CST = 6., EST = 5.

Length of run (hours) (IRLG) -- No default ! IRLG = 8736 !

Number of chemical species (NSPEC)  
Default: 5 ! NSPEC = 9 !

Number of chemical species  
to be emitted (NSE) Default: 3 ! NSE = 7 !

Flag to stop run after  
SETUP phase (ITEST) Default: 2 ! ITEST = 2 !  
(Used to allow checking  
of the model inputs, files, etc.)  
ITEST = 1 - STOPS program after SETUP phase  
ITEST = 2 - Continues with execution of program  
after SETUP

Restart Configuration:

Control flag (MRESTART) Default: 0 ! MRESTART = 0 !

0 = Do not read or write a restart file  
1 = Read a restart file at the beginning of  
the run  
2 = Write a restart file during run  
3 = Read a restart file at beginning of run  
and write a restart file during run

Number of periods in Restart  
output cycle (NRESPD) Default: 0 ! NRESPD = 0 !

0 = File written only at last period  
>0 = File updated every NRESPD periods

Meteorological Data Format (METFM)  
Default: 1 ! METFM = 1 !

METFM = 1 - CALMET binary file (CALMET.MET)  
METFM = 2 - ISC ASCII file (ISCMET.MET)  
METFM = 3 - AUSPLUME ASCII file (PLMMET.MET)  
METFM = 4 - CTDM plus tower file (PROFILE.DAT) and  
surface parameters file (SURFACE.DAT)

PG sigma-y is adjusted by the factor (AVET/PGTIME)\*\*0.2  
Averaging Time (minutes) (AVET)  
Default: 60.0 ! AVET = 60. !

PG Averaging Time (minutes) (PGTIME)  
Default: 60.0 ! PGTIME = 60. !

!END!

-----  
INPUT GROUP: 2 -- Technical options  
-----

Vertical distribution used in the  
near field (MGAUSS) Default: 1 ! MGAUSS = 1 !  
0 = uniform  
1 = Gaussian

Terrain adjustment method  
(MCTADJ) Default: 3 ! MCTADJ = 3 !  
0 = no adjustment  
1 = ISC-type of terrain adjustment  
2 = simple, CALPUFF-type of terrain  
adjustment  
3 = partial plume path adjustment

Subgrid-scale complex terrain  
flag (MCTSG) Default: 0 ! MCTSG = 0 !  
0 = not modeled  
1 = modeled

Near-field puffs modeled as  
elongated 0 (MSLUG) Default: 0 ! MSLUG = 0 !  
0 = no  
1 = yes (slug model used)

Transitional plume rise modeled ?  
(MTRANS) Default: 1 ! MTRANS = 1 !  
0 = no (i.e., final rise only)

1 = yes (i.e., transitional rise computed)

Stack tip downwash? (MTIP)                    Default: 1        ! MTIP = 1 !  
0 = no (i.e., no stack tip downwash)  
1 = yes (i.e., use stack tip downwash)

Method used to simulate building  
downwash? (MBDW)                            Default: 1        ! MBDW = 1 !  
1 = ISC method  
2 = PRIME method

Vertical wind shear modeled above  
stack top? (MSHEAR)                        Default: 0        ! MSHEAR = 0 !  
0 = no (i.e., vertical wind shear not modeled)  
1 = yes (i.e., vertical wind shear modeled)

Puff splitting allowed? (MSPLIT)            Default: 0        ! MSPLIT = 0 !  
0 = no (i.e., puffs not split)  
1 = yes (i.e., puffs are split)

Chemical mechanism flag (MCHEM)            Default: 1        ! MCHEM = 1 !  
0 = chemical transformation not  
    modeled  
1 = transformation rates computed  
    internally (MESOPUFF II scheme)  
2 = user-specified transformation  
    rates used  
3 = transformation rates computed  
    internally (RIVAD/ARM3 scheme)  
4 = secondary organic aerosol formation  
    computed (MESOPUFF II scheme for OH)

Aqueous phase transformation flag (MAQCHEM)  
(Used only if MCHEM = 1, or 3)            Default: 0        ! MAQCHEM = 0 !  
0 = aqueous phase transformation  
    not modeled  
1 = transformation rates adjusted  
    for aqueous phase reactions

Wet removal modeled ? (MWET)              Default: 1        ! MWET = 1 !  
0 = no  
1 = yes

Dry deposition modeled ? (MDRY)            Default: 1        ! MDRY = 1 !  
0 = no  
1 = yes  
(dry deposition method specified  
  for each species in Input Group 3)

Method used to compute dispersion  
coefficients (MDISP)                        Default: 3        ! MDISP = 3 !  
  
1 = dispersion coefficients computed from measured values  
  of turbulence, sigma v, sigma w  
2 = dispersion coefficients from internally calculated  
  sigma v, sigma w using micrometeorological variables  
  (u\*, w\*, L, etc.)  
3 = PG dispersion coefficients for RURAL areas (computed using  
  the ISCST multi-segment approximation) and MP coefficients in  
  urban areas  
4 = same as 3 except PG coefficients computed using  
  the MESOPUFF II eqns.  
5 = CTDM sigmas used for stable and neutral conditions.  
  For unstable conditions, sigmas are computed as in  
  MDISP = 3, described above. MDISP = 5 assumes that

measured values are read

Sigma-v/sigma-theta, sigma-w measurements used? (MTURBVW)  
(Used only if MDISP = 1 or 5)           Default: 3           ! MTURBVW = 3 !  
1 = use sigma-v or sigma-theta measurements  
    from PROFILE.DAT to compute sigma-y  
    (valid for METFM = 1, 2, 3, 4)  
2 = use sigma-w measurements  
    from PROFILE.DAT to compute sigma-z  
    (valid for METFM = 1, 2, 3, 4)  
3 = use both sigma-(v/theta) and sigma-w  
    from PROFILE.DAT to compute sigma-y and sigma-z  
    (valid for METFM = 1, 2, 3, 4)  
4 = use sigma-theta measurements  
    from PLMMET.DAT to compute sigma-y  
    (valid only if METFM = 3)

Back-up method used to compute dispersion  
when measured turbulence data are  
missing (MDISP2)                        Default: 3           ! MDISP2 = 3 !  
(used only if MDISP = 1 or 5)  
2 = dispersion coefficients from internally calculated  
    sigma v, sigma w using micrometeorological variables  
    (u\*, w\*, L, etc.)  
3 = PG dispersion coefficients for RURAL areas (computed using  
    the ISCST multi-segment approximation) and MP coefficients in  
    urban areas  
4 = same as 3 except PG coefficients computed using  
    the MESOPUFF II eqns.

PG sigma-y,z adj. for roughness?       Default: 0           ! MROUGH = 0 !  
(MROUGH)  
0 = no  
1 = yes

Partial plume penetration of  
elevated inversion?                    Default: 1           ! MPARTL = 1 !  
(MPARTL)  
0 = no  
1 = yes

Strength of temperature inversion      Default: 0           ! MTINV = 0 !  
provided in PROFILE.DAT extended records?  
(MTINV)  
0 = no (computed from measured/default gradients)  
1 = yes

PDF used for dispersion under convective conditions?  
  Default: 0           ! MPDF = 0 !  
(MPDF)  
0 = no  
1 = yes

Sub-Grid TIBL module used for shore line?  
  Default: 0           ! MSGTIBL = 0 !  
(MSGTIBL)  
0 = no  
1 = yes

Boundary conditions (concentration) modeled?  
  Default: 0           ! MBCON = 0 !  
(MBCON)  
0 = no  
1 = yes, using formatted BCON.DAT file  
2 = yes, using unformatted CONC.DAT file

---

Analyses of fogging and icing impacts due to emissions from arrays of mechanically-forced cooling towers can be performed using CALPUFF in conjunction with a cooling tower emissions processor (CTEMISS) and its associated postprocessors. Hourly emissions of water vapor and temperature from each cooling tower cell are computed for the current cell configuration and ambient conditions by CTEMISS. CALPUFF models the dispersion of these emissions and provides cloud information in a specialized format for further analysis. Output to FOG.DAT is provided in either 'plume mode' or 'receptor mode' format.

Configure for FOG Model output? Default: 0 ! MFOG = 0 !  
(MFOG)  
0 = no  
1 = yes - report results in PLUME Mode format  
2 = yes - report results in RECEPTOR Mode format

TEST options specified to see if they conform to regulatory values? (MREG) Default: 1 ! MREG = 0 !

0 = NO checks are made  
1 = Technical options must conform to USEPA  
Long Range Transport (LRT) guidance  
METFM 1 or 2  
AVET 60. (min)  
PGTIME 60. (min)  
MGAUSS 1  
MCTADJ 3  
MTRANS 1  
MTIP 1  
MCHEM 1 or 3 (if modeling SOx, NOx)  
MWET 1  
MDRY 1  
MDISP 2 or 3  
MPDF 0 if MDISP=3  
1 if MDISP=2  
MROUGH 0  
MPARTL 1  
SYTDEP 550. (m)  
MHFTSZ 0

!END!

-----  
INPUT GROUP: 3a, 3b -- Species list  
-----

-----  
Subgroup (3a)  
-----

The following species are modeled:

The following species are modeled:

! CSPEC = SO2 ! !END!  
! CSPEC = SO4 ! !END!

```
! CSPEC =      NOX !      !END!
! CSPEC =      HNO3 !     !END!
! CSPEC =      NO3 !     !END!
! CSPEC =      PMC !     !END!
! CSPEC =      PMF !     !END!
! CSPEC =      EC !     !END!
! CSPEC =      SOA !     !END!
```

SPECIES NAME (Limit: 12 Characters in length)	MODELED (0=NO, 1=YES)	EMITTED (0=NO, 1=YES)	Dry DEPOSITED (0=NO, 1=COMPUTED-GAS 2=COMPUTED-PARTICLE 3=USER-SPECIFIED)	OUTPUT GROUP NUMBER (0=NONE, 1=1st CGRUP, 2=2nd CGRUP, 3= etc.)
! SO2 =	1,	1,	1,	0 !
! SO4 =	1,	1,	2,	0 !
! NOX =	1,	1,	1,	0 !
! HNO3 =	1,	0,	1,	0 !
! NO3 =	1,	0,	2,	0 !
! PMC =	1,	1,	2,	0 !
! PMF =	1,	1,	2,	0 !
! EC =	1,	1,	2,	0 !
! SOA =	1,	1,	2,	0 !

!END!

-----  
Subgroup (3b)  
-----

The following names are used for Species-Groups in which results for certain species are combined (added) prior to output. The CGRUP name will be used as the species name in output files. Use this feature to model specific particle-size distributions by treating each size-range as a separate species. Order must be consistent with 3(a) above.

-----  
INPUT GROUP: 4 -- Map Projection and Grid control parameters  
-----

Projection for all (X,Y):  
-----

Map projection

(PMAP)                                   Default: UTM     ! PMAP = LCC !

UTM : Universal Transverse Mercator  
TTM : Tangential Transverse Mercator  
LCC : Lambert Conformal Conic  
PS : Polar Stereographic  
EM : Equatorial Mercator  
LAZA : Lambert Azimuthal Equal Area

False Easting and Northing (km) at the projection origin

(Used only if PMAP= TTM, LCC, or LAZA)

(FEAST)                                Default=0.0     ! FEAST = 0.000 !  
(FNORTH)                               Default=0.0     ! FNORTH = 0.000 !

UTM zone (1 to 60)



(Used only if PMAP=UTM)  
(IUTMZN)                      No Default                      ! IUTMZN = 0                      !

Hemisphere for UTM projection?  
(Used only if PMAP=UTM)  
(UTMHEM)                      Default: N                      ! UTMHEM = N                      !  
    N   : Northern hemisphere projection  
    S   : Southern hemisphere projection

Latitude and Longitude (decimal degrees) of projection origin  
(Used only if PMAP= TTM, LCC, PS, EM, or LAZA)  
(RLAT0)                      No Default                      ! RLAT0 = 40N                      !  
(RLON0)                      No Default                      ! RLON0 = 97W                      !

- TTM : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience
- LCC : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience
- PS : RLON0 identifies central (grid N/S) meridian of projection  
      RLAT0 selected for convenience
- EM : RLON0 identifies central meridian of projection  
      RLAT0 is REPLACED by 0.0N (Equator)
- LAZA: RLON0 identifies longitude of tangent-point of mapping plane  
      RLAT0 identifies latitude of tangent-point of mapping plane

Matching parallel(s) of latitude (decimal degrees) for projection  
(Used only if PMAP= LCC or PS)  
(XLAT1)                      No Default                      ! XLAT1 = 33N                      !  
(XLAT2)                      No Default                      ! XLAT2 = 45N                      !

- LCC : Projection cone slices through Earth's surface at XLAT1 and XLAT2
- PS : Projection plane slices through Earth at XLAT1  
      (XLAT2 is not used)

-----  
Note: Latitudes and longitudes should be positive, and include a  
      letter N,S,E, or W indicating north or south latitude, and  
      east or west longitude. For example,  
      35.9 N Latitude = 35.9N  
      118.7 E Longitude = 118.7E

Datum-region  
-----

The Datum-Region for the coordinates is identified by a character  
string. Many mapping products currently available use the model of the  
Earth known as the World Geodetic System 1984 (WGS-84). Other local  
models may be in use, and their selection in CALMET will make its output  
consistent with local mapping products. The list of Datum-Regions with  
official transformation parameters is provided by the National Imagery and  
Mapping Agency (NIMA).

NIMA Datum - Regions(Examples)  
-----

- WGS-84   WGS-84 Reference Ellipsoid and Geoid, Global coverage (WGS84)
- NAS-C    NORTH AMERICAN 1927 Clarke 1866 Spheroid, MEAN FOR CONUS (NAD27)
- NAR-C    NORTH AMERICAN 1983 GRS 80 Spheroid, MEAN FOR CONUS (NAD83)
- NWS-84   NWS 6370KM Radius, Sphere
- ESR-S    ESRI REFERENCE 6371KM Radius, Sphere

Datum-region for output coordinates  
(DATUM)                      Default: WGS-84                      ! DATUM = NAS-C                      !

METEOROLOGICAL Grid:

Rectangular grid defined for projection PMAP,  
with X the Easting and Y the Northing coordinate

No. X grid cells (NX)	No default	! NX = 162 !
No. Y grid cells (NY)	No default	! NY = 191 !
No. vertical layers (NZ)	No default	! NZ = 10 !
Grid spacing (DGRIDKM)	No default	! DGRIDKM = 4. !
	Units: km	

Cell face heights in arbitrary  
vertical grid (ZFACE(NZ+1)) No defaults  
Units: m  
! ZFACE = 0.,20.,40.,80.,160.,320.,640.,1200.,2000.,3000.,4000. !

Reference Coordinates  
of SOUTHWEST corner of  
grid cell(1, 1):

X coordinate (XORIGKM)	No default	! XORIGKM = 247. !
Y coordinate (YORIGKM)	No default	! YORIGKM = 240. !
	Units: km	

COMPUTATIONAL Grid:

The computational grid is identical to or a subset of the MET. grid.  
The lower left (LL) corner of the computational grid is at grid point  
(IBCOMP, JBCOMP) of the MET. grid. The upper right (UR) corner of the  
computational grid is at grid point (IECOMP, JECOMP) of the MET. grid.  
The grid spacing of the computational grid is the same as the MET. grid.

X index of LL corner (IBCOMP) (1 <= IBCOMP <= NX)	No default	! IBCOMP = 1 !
Y index of LL corner (JBCOMP) (1 <= JBCOMP <= NY)	No default	! JBCOMP = 1 !
X index of UR corner (IECOMP) (1 <= IECOMP <= NX)	No default	! IECOMP = 162 !
Y index of UR corner (JECOMP) (1 <= JECOMP <= NY)	No default	! JECOMP = 191 !

SAMPLING Grid (GRIDDED RECEPTORS):

The lower left (LL) corner of the sampling grid is at grid point  
(IBSAMP, JBSAMP) of the MET. grid. The upper right (UR) corner of the  
sampling grid is at grid point (IESAMP, JESAMP) of the MET. grid.  
The sampling grid must be identical to or a subset of the computational  
grid. It may be a nested grid inside the computational grid.  
The grid spacing of the sampling grid is DGRIDKM/MESHNDN.

Logical flag indicating if gridded receptors are used (LSAMP) (T=yes, F=no)	Default: T	! LSAMP = F !
X index of LL corner (IBSAMP) (IBCOMP <= IBSAMP <= IECOMP)	No default	! IBSAMP = 1 !

**Refined BART Modeling Protocol for  
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```

Y index of LL corner (JBSAMP)      No default      ! JBSAMP = 1  !
  (JBCOMP <= JBSAMP <= JECOMP)

X index of UR corner (IESAMP)      No default      ! IESAMP = 162 !
  (IBCOMP <= IESAMP <= IECOMP)

Y index of UR corner (JESAMP)      No default      ! JESAMP = 191 !
  (JBCOMP <= JESAMP <= JECOMP)

Nesting factor of the sampling
grid (MESH DN)                      Default: 1      ! MESH DN = 1  !
(MESH DN is an integer >= 1)

!END!

```

-----

INPUT GROUP: 5 -- Output Options

-----

FILE	* DEFAULT VALUE	* VALUE THIS RUN
----	-----	-----
Concentrations (ICON)	1	! ICON = 1  !
Dry Fluxes (IDRY)	1	! IDRY = 0  !
Wet Fluxes (IWET)	1	! IWET = 0  !
Relative Humidity (IVIS) (relative humidity file is required for visibility analysis)	1	! IVIS = 0  !
Use data compression option in output file? (LCOMPRS)	Default: T	! LCOMPRS = T !

\*  
0 = Do not create file, 1 = create file

DIAGNOSTIC MASS FLUX OUTPUT OPTIONS:

```

Mass flux across specified boundaries
for selected species reported hourly?
(IMFLX)                      Default: 0      ! IMFLX = 0  !
  0 = no
  1 = yes (FLUXBDY.DAT and MASSFLX.DAT filenames
           are specified in Input Group 0)

```

```

Mass balance for each species
reported hourly?
(IMBAL)                      Default: 0      ! IMBAL = 0  !
  0 = no
  1 = yes (MASSBAL.DAT filename is
           specified in Input Group 0)

```

LINE PRINTER OUTPUT OPTIONS:

```

Print concentrations (ICPRT)    Default: 0      ! ICPRT = 0  !
Print dry fluxes (IDPRT)       Default: 0      ! IDPRT = 0  !
Print wet fluxes (IWPRT)       Default: 0      ! IWPRT = 0  !
(0 = Do not print, 1 = Print)

```

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---

Concentration print interval  
(ICFRQ) in hours                   Default: 1                   ! ICFRQ = 1 !  
Dry flux print interval  
(IDFRQ) in hours                   Default: 1                   ! IDFRQ = 1 !  
Wet flux print interval  
(IWFRQ) in hours                   Default: 1                   ! IWFRQ = 1 !

Units for Line Printer Output  
(IPRTU)                               Default: 1                   ! IPRTU = 3 !  
                  for                    for  
                  Concentration        Deposition  
1 =            g/m\*\*3                g/m\*\*2/s  
2 =            mg/m\*\*3               mg/m\*\*2/s  
3 =            ug/m\*\*3               ug/m\*\*2/s  
4 =            ng/m\*\*3               ng/m\*\*2/s  
5 =            Odour Units

Messages tracking progress of run  
written to the screen ?  
(IMESG)                               Default: 2                   ! IMESG = 2 !  
0 = no  
1 = yes (advection step, puff ID)  
2 = yes (YYYYJJJHH, # old puffs, # emitted puffs)

SPECIES (or GROUP for combined species) LIST FOR OUTPUT OPTIONS

-- MASS FLUX --	---- CONCENTRATIONS ----		----- DRY FLUXES -----		----- WET FLUXES -----	
SPECIES /GROUP SAVED ON DISK?	PRINTED?	SAVED ON DISK?	PRINTED?	SAVED ON DISK?	PRINTED?	SAVED ON DISK?
-----	-----	-----	-----	-----	-----	-----
!            SO2 =	0,	1,	0,	0,	0,	0,
0 !						
!            SO4 =	0,	1,	0,	0,	0,	0,
0 !						
!            NOX =	0,	1,	0,	0,	0,	0,
0 !						
!            HNO3 =	0,	1,	0,	0,	0,	0,
0 !						
!            NO3 =	0,	1,	0,	0,	0,	0,
0 !						
!            PMC =	0,	1,	0,	0,	0,	0,
0 !						
!            PMF =	0,	1,	0,	0,	0,	0,
0 !						
!            EC =	0,	1,	0,	0,	0,	0,
0 !						
!            SOA =	0,	1,	0,	0,	0,	0,
0 !						

OPTIONS FOR PRINTING "DEBUG" QUANTITIES (much output)

Logical for debug output  
(LDEBUG)                               Default: F                   ! LDEBUG = F !  
  
First puff to track  
(IPFDEB)                               Default: 1                   ! IPFDEB = 1 !  
  
Number of puffs to track  
(NPFDEB)                               Default: 1                   ! NPFDEB = 1 !

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```

Met. period to start output
(NN1)                               Default: 1      ! NN1 = 1  !

Met. period to end output
(NN2)                               Default: 10    ! NN2 = 10 !

!END!

```

-----

INPUT GROUP: 6a, 6b, & 6c -- Subgrid scale complex terrain inputs

-----

-----

Subgroup (6a)

-----

```

Number of terrain features (NHILL)   Default: 0      ! NHILL = 0  !

Number of special complex terrain
receptors (NCTREC)                   Default: 0      ! NCTREC = 0  !

Terrain and CTSG Receptor data for
CTSG hills input in CTDM format ?
(MHILL)                               No Default     ! MHILL = 2  !
1 = Hill and Receptor data created
  by CTDM processors & read from
  HILL.DAT and HILLRCT.DAT files
2 = Hill data created by OPTHILL &
  input below in Subgroup (6b);
  Receptor data in Subgroup (6c)

Factor to convert horizontal dimensions
to meters (MHILL=1)                   Default: 1.0    ! XHILL2M = 1. !

Factor to convert vertical dimensions
to meters (MHILL=1)                   Default: 1.0    ! ZHILL2M = 1. !

X-origin of CTDM system relative to
CALPUFF coordinate system, in Kilometers (MHILL=1) No Default     ! XCTDMKM = 0.0E00 !

Y-origin of CTDM system relative to
CALPUFF coordinate system, in Kilometers (MHILL=1) No Default     ! YCTDMKM = 0.0E00 !

```

! END !

-----

Subgroup (6b)

-----

```

                1 **
HILL information

HILL      XC      YC      THETAH  ZGRID  RELIEF   EXPO 1   EXPO 2   SCALE 1   SCALE 2
AMAX1     AMAX2
NO.       (km)    (km)    (deg.)  (m)    (m)     (m)     (m)     (m)     (m)
(m)       (m)
-----
-----

```

-----

Subgroup (6c)

---

-----  
COMPLEX TERRAIN RECEPTOR INFORMATION

XRCT (km)	YRCT (km)	ZRCT (m)	XHH
-----	-----	-----	----

-----  
1

Description of Complex Terrain Variables:

XC, YC = Coordinates of center of hill  
 THETAH = Orientation of major axis of hill (clockwise from North)  
 ZGRID = Height of the 0 of the grid above mean sea level  
 RELIEF = Height of the crest of the hill above the grid elevation  
 EXPO 1 = Hill-shape exponent for the major axis  
 EXPO 2 = Hill-shape exponent for the minor axis  
 SCALE 1 = Horizontal length scale along the major axis  
 SCALE 2 = Horizontal length scale along the minor axis  
 AMAX = Maximum allowed axis length for the major axis  
 BMAX = Maximum allowed axis length for the minor axis  
  
 XRCT, YRCT = Coordinates of the complex terrain receptors  
 ZRCT = Height of the ground (MSL) at the complex terrain Receptor  
 XHH = Hill number associated with each complex terrain receptor  
 (NOTE: MUST BE ENTERED AS A REAL NUMBER)

\*\*

NOTE: DATA for each hill and CTSG receptor are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUP: 7 -- Chemical parameters for dry deposition of gases

SPECIES LAW COEFFICIENT NAME (dimensionless)	DIFFUSIVITY (cm**2/s)	ALPHA STAR	REACTIVITY	MESOPHYLL RESISTANCE (s/cm)	HENRY'S
-----	-----	-----	-----	-----	-----
! SO2 =	0.1509,	1000.,	8.,	0.,	
0.04 !					
! NOX =	0.1656,	1.,	8.,	5.,	
3.5 !					
! HNO3 =	0.1628,	1.,	18.,	0.,	
0.00000008 !					
!END!					

-----  
INPUT GROUP: 8 -- Size parameters for dry deposition of particles

For SINGLE SPECIES, the mean and standard deviation are used to compute a deposition velocity for NINT (see group 9) size-ranges,

and these are then averaged to obtain a mean deposition velocity.

For GROUPED SPECIES, the size distribution should be explicitly specified (by the 'species' in the group), and the standard deviation for each should be entered as 0. The model will then use the deposition velocity for the stated mean diameter.

SPECIES NAME	GEOMETRIC MASS MEAN DIAMETER (microns)	GEOMETRIC STANDARD DEVIATION (microns)
-----	-----	-----
! SO4 =	0.48,	2. !
! NO3 =	0.48,	2. !
! PMC =	2.0,	2.0 !
! PMF =	0.48,	2. !
! EC =	0.48,	2. !
! SOA =	0.48,	2. !

!END!

-----

INPUT GROUP: 9 -- Miscellaneous dry deposition parameters

-----

Reference cuticle resistance (s/cm)  
(RCUTR) Default: 30 ! RCUTR = 30.0 !

Reference ground resistance (s/cm)  
(RGR) Default: 10 ! RGR = 10.0 !

Reference pollutant reactivity  
(REACTR) Default: 8 ! REACTR = 8.0 !

Number of particle-size intervals used to  
evaluate effective particle deposition velocity  
(NINT) Default: 9 ! NINT = 9 !

Vegetation state in unirrigated areas  
(IVEG) Default: 1 ! IVEG = 1 !  
IVEG=1 for active and unstressed vegetation  
IVEG=2 for active and stressed vegetation  
IVEG=3 for inactive vegetation

!END!

-----

INPUT GROUP: 10 -- Wet Deposition Parameters

-----

Scavenging Coefficient -- Units: (sec)\*\*(-1)

Pollutant	Liquid Precip.	Frozen Precip.
-----	-----	-----
! SO2 =	3.0E-05,	0.0E00 !
! SO4 =	1.0E-04,	3.0E-05 !
! NOX =	0.0E00,	0.0E00 !
! HNO3 =	6.0E-05,	0.0E00 !
! NO3 =	1.0E-04,	3.0E-05 !

!END!

-----  
INPUT GROUP: 11 -- Chemistry Parameters  
-----

Ozone data input option (MOZ)       Default: 1                   ! MOZ = 0 !  
(Used only if MCHEM = 1, 3, or 4)  
0 = use a monthly background ozone value  
1 = read hourly ozone concentrations from  
the OZONE.DAT data file

Monthly ozone concentrations  
(Used only if MCHEM = 1, 3, or 4 and  
MOZ = 0 or MOZ = 1 and all hourly O3 data missing)  
(BCKO3) in ppb                       Default: 12\*80.

! BCKO3 = 31.00, 31.00, 31.00, 37.00, 37.00, 37.00, 33.00, 33.00, 33.00, 27.00, 27.00, 27.00

!

Monthly ammonia concentrations  
(Used only if MCHEM = 1, or 3)  
(BCKNH3) in ppb                       Default: 12\*10.

! BCKNH3 = 12\*3.9 !

Nighttime SO2 loss rate (RNITE1)  
in percent/hour                       Default: 0.2                   ! RNITE1 = .2 !

Nighttime NOx loss rate (RNITE2)  
in percent/hour                       Default: 2.0                   ! RNITE2 = 2.0 !

Nighttime HNO3 formation rate (RNITE3)  
in percent/hour                       Default: 2.0                   ! RNITE3 = 2.0 !

H2O2 data input option (MH2O2)       Default: 1                   ! MH2O2 = 1 !  
(Used only if MQACHEM = 1)  
0 = use a monthly background H2O2 value  
1 = read hourly H2O2 concentrations from  
the H2O2.DAT data file

Monthly H2O2 concentrations  
(Used only if MQACHEM = 1 and  
MH2O2 = 0 or MH2O2 = 1 and all hourly H2O2 data missing)  
(BCKH2O2) in ppb                       Default: 12\*1.

\* BCKH2O2 = 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00 \*

! BCKH2O2 = 0.50, 0.50, 0.50, 2.30, 2.30, 2.30, 3.50, 3.50, 3.50, 0.80, 0.80, 0.80 !

--- Data for SECONDARY ORGANIC AEROSOL (SOA) Option  
(used only if MCHEM = 4)

The SOA module uses monthly values of:  
Fine particulate concentration in ug/m<sup>3</sup> (BCKPMF)  
Organic fraction of fine particulate (OFRAC)  
VOC / NOX ratio (after reaction) (VCNX)

to characterize the air mass when computing  
the formation of SOA from VOC emissions.  
Typical values for several distinct air mass types are:

Month	1	2	3	4	5	6	7	8	9	10	11	12
	Jan	Feb	Mar	Apr	May	Jun	Jul	Aug	Sep	Oct	Nov	Dec



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---

Clean Continental

BCKPMF	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.
OFRAC	.15	.15	.20	.20	.20	.20	.20	.20	.20	.20	.20	.15
VCNX	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.

Clean Marine (surface)

BCKPMF	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5
OFRAC	.25	.25	.30	.30	.30	.30	.30	.30	.30	.30	.30	.25
VCNX	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.

Urban - low biogenic (controls present)

BCKPMF	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.
OFRAC	.20	.20	.25	.25	.25	.25	.25	.25	.20	.20	.20	.20
VCNX	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.

Urban - high biogenic (controls present)

BCKPMF	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.
OFRAC	.25	.25	.30	.30	.30	.55	.55	.55	.35	.35	.35	.25
VCNX	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.

Regional Plume

BCKPMF	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.
OFRAC	.20	.20	.25	.35	.25	.40	.40	.40	.30	.30	.30	.20
VCNX	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.

Urban - no controls present

BCKPMF	100.	100.	100.	100.	100.	100.	100.	100.	100.	100.	100.	100.
OFRAC	.30	.30	.35	.35	.35	.55	.55	.55	.35	.35	.35	.30
VCNX	2.	2.	2.	2.	2.	2.	2.	2.	2.	2.	2.	2.

Default: Clean Continental

! BCKPMF = 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00 !  
! OFRAC = 0.15, 0.15, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.15 !  
! VCNX = 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00 !

!END!

-----  
INPUT GROUP: 12 -- Misc. Dispersion and Computational Parameters  
-----

Horizontal size of puff (m) beyond which  
time-dependent dispersion equations (Heffter)  
are used to determine sigma-y and  
sigma-z (SYTDEP) Default: 550. ! SYTDEP = 5.5E02 !

Switch for using Heffter equation for sigma z  
as above (0 = Not use Heffter; 1 = use Heffter  
(MHFTSZ) Default: 0 ! MHFTSZ = 0 !

Stability class used to determine plume  
growth rates for puffs above the boundary  
layer (JSUP) Default: 5 ! JSUP = 5 !

Vertical dispersion constant for stable  
conditions (k1 in Eqn. 2.7-3) (CONK1) Default: 0.01 ! CONK1 = .01 !

Vertical dispersion constant for neutral/  
unstable conditions (k2 in Eqn. 2.7-4)

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(CONK2) Default: 0.1 ! CONK2 = .1 !

Factor for determining Transition-point from  
Schulman-Scire to Huber-Snyder Building Downwash  
scheme (SS used for Hs < Hb + TBD \* HL)  
(TBD) Default: 0.5 ! TBD = .5 !  
TBD < 0 ==> always use Huber-Snyder  
TBD = 1.5 ==> always use Schulman-Scire  
TBD = 0.5 ==> ISC Transition-point

Range of land use categories for which  
urban dispersion is assumed  
(IURB1, IURB2) Default: 10 ! IURB1 = 10 !  
19 ! IURB2 = 19 !

Site characterization parameters for single-point Met data files -----  
(needed for METFM = 2,3,4)

Land use category for modeling domain  
(ILANDUIN) Default: 20 ! ILANDUIN = 20 !

Roughness length (m) for modeling domain  
(Z0IN) Default: 0.25 ! Z0IN = .25 !

Leaf area index for modeling domain  
(XLAIIN) Default: 3.0 ! XLAIIN = 3.0 !

Elevation above sea level (m)  
(ELEVIN) Default: 0.0 ! ELEVIN = .0 !

Latitude (degrees) for met location  
(XLATIN) Default: -999. ! XLATIN = -999.0 !

Longitude (degrees) for met location  
(XLONIN) Default: -999. ! XLONIN = -999.0 !

Specialized information for interpreting single-point Met data files -----

Anemometer height (m) (Used only if METFM = 2,3)  
(ANEMHT) Default: 10. ! ANEMHT = 10.0 !

Form of lateral turbulence data in PROFILE.DAT file  
(Used only if METFM = 4 or MTURBVW = 1 or 3)  
(ISIGMAV) Default: 1 ! ISIGMAV = 1 !  
0 = read sigma-theta  
1 = read sigma-v

Choice of mixing heights (Used only if METFM = 4)  
(IMIXCTDM) Default: 0 ! IMIXCTDM = 0 !  
0 = read PREDICTED mixing heights  
1 = read OBSERVED mixing heights

Maximum length of a slug (met. grid units)  
(MXLEN) Default: 1.0 ! MXLEN = 1.0 !

Maximum travel distance of a puff/slug (in  
grid units) during one sampling step  
(XSAMLEN) Default: 1.0 ! XSAMLEN = 1.0 !

Maximum Number of slugs/puffs release from  
one source during one time step  
(MXNEW) Default: 99 ! MXNEW = 99 !

Maximum Number of sampling steps for  
one puff/slug during one time step

---

**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

---

```

(MXSAM)                                Default: 99      ! MXSAM = 99  !

Number of iterations used when computing
the transport wind for a sampling step
that includes gradual rise (for CALMET
and PROFILE winds)
(NCOUNT)                              Default: 2      ! NCOUNT = 2  !

Minimum sigma y for a new puff/slug (m)
(SYMIN)                                Default: 1.0    ! SYMIN = 1.0  !

Minimum sigma z for a new puff/slug (m)
(SZMIN)                                Default: 1.0    ! SZMIN = 1.0  !

Default minimum turbulence velocities
sigma-v and sigma-w for each
stability class (m/s)
(SVMIN(6) and SWMIN(6))
      Default SVMIN : .50, .50, .50, .50, .50, .50
      Default SWMIN : .20, .12, .08, .06, .03, .016

      Stability Class :  A    B    C    D    E    F
                       ---  ---  ---  ---  ---  ---
      ! SVMIN = 0.500, 0.500, 0.500, 0.500, 0.500, 0.500!
      ! SWMIN = 0.200, 0.120, 0.080, 0.060, 0.030, 0.016!

Divergence criterion for dw/dz across puff
used to initiate adjustment for horizontal
convergence (1/s)
Partial adjustment starts at CDIV(1), and
full adjustment is reached at CDIV(2)
(CDIV(2))                              Default: 0.0,0.0 ! CDIV = .0, .0 !

Minimum wind speed (m/s) allowed for
non-calm conditions. Also used as minimum
speed returned when using power-law
extrapolation toward surface
(WSCALM)                               Default: 0.5    ! WSCALM = .5  !

Maximum mixing height (m)
(XMAXZI)                               Default: 3000.  ! XMAXZI = 3000.0 !

Minimum mixing height (m)
(XMINZI)                               Default: 50.    ! XMINZI = 50.0  !

Default wind speed classes --
5 upper bounds (m/s) are entered;
the 6th class has no upper limit
(WSCAT(5))
      Default      :
      ISC RURAL   : 1.54, 3.09, 5.14, 8.23, 10.8 (10.8+)

      Wind Speed Class :  1    2    3    4    5
                       ---  ---  ---  ---  ---
      ! WSCAT = 1.54, 3.09, 5.14, 8.23, 10.80 !

Default wind speed profile power-law
exponents for stabilities 1-6
(PLX0(6))
      Default      : ISC RURAL values
      ISC RURAL   : .07, .07, .10, .15, .35, .55
      ISC URBAN   : .15, .15, .20, .25, .30, .30

      Stability Class :  A    B    C    D    E    F
                       ---  ---  ---  ---  ---  ---
      ! PLX0 = 0.07, 0.07, 0.10, 0.15, 0.35, 0.55 !

Default potential temperature gradient

```

**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

---

```
for stable classes E, F (degK/m)
(PYG0(2))                               Default: 0.020, 0.035
                                         ! PYG0 = 0.020, 0.035 !

Default plume path coefficients for
each stability class (used when option
for partial plume height terrain adjustment
is selected -- MCTADJ=3)
(PPC(6))                                Stability Class :  A    B    C    D    E    F
                                         Default  PPC :  .50, .50, .50, .50, .35, .35
                                         ---    ---    ---    ---    ---    ---
                                         !  PPC = 0.50, 0.50, 0.50, 0.50, 0.35, 0.35 !

Slug-to-puff transition criterion factor
equal to sigma-y/length of slug
(SL2PF)                                  Default: 10.          ! SL2PF = 10.0 !

Puff-splitting control variables -----

VERTICAL SPLIT
-----

Number of puffs that result every time a puff
is split - nsplit=2 means that 1 puff splits
into 2
(NSPLIT)                                 Default: 3           ! NSPLIT = 3 !

Time(s) of a day when split puffs are eligible to
be split once again; this is typically set once
per day, around sunset before nocturnal shear develops.
24 values: 0 is midnight (00:00) and 23 is 11 PM (23:00)
0=do not re-split    1=eligible for re-split
(IRESPLIT(24))       Default: Hour 17 = 1
! IRESPLIT = 0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0,0 !

Split is allowed only if last hour's mixing
height (m) exceeds a minimum value
(ZISPLIT)                                 Default: 100.        ! ZISPLIT = 100.0 !

Split is allowed only if ratio of last hour's
mixing ht to the maximum mixing ht experienced
by the puff is less than a maximum value (this
postpones a split until a nocturnal layer develops)
(ROLDMAX)                                 Default: 0.25        ! ROLDMAX = 0.25 !

HORIZONTAL SPLIT
-----

Number of puffs that result every time a puff
is split - nsplith=5 means that 1 puff splits
into 5
(NSPLITH)                                 Default: 5           ! NSPLITH = 5 !

Minimum sigma-y (Grid Cells Units) of puff
before it may be split
(SYSPLITH)                                 Default: 1.0         ! SYSPLITH = 1.0 !

Minimum puff elongation rate (SYSPLITH/hr) due to
wind shear, before it may be split
(SHSPLITH)                                 Default: 2.          ! SHSPLITH = 2.0 !

Minimum concentration (g/m^3) of each
species in puff before it may be split
Enter array of NSPEC values; if a single value is
```

**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

---

```
entered, it will be used for ALL species
(CNSPLITH)                      Default:  1.0E-07   ! CNSPLITH = 1.0E-07 !

Integration control variables -----

Fractional convergence criterion for numerical SLUG
sampling integration
(EPSSLUG)                      Default:  1.0e-04   ! EPSSLUG = 1.0E-04 !

Fractional convergence criterion for numerical AREA
source integration
(EPSAREA)                      Default:  1.0e-06   ! EPSAREA = 1.0E-06 !

Trajectory step-length (m) used for numerical rise
integration
(DSRISE)                       Default:  1.0       ! DSRISE = 1.0 !

Boundary Condition (BC) Puff control variables -----

Minimum height (m) to which BC puffs are mixed as they are emitted
(MBCON=2 ONLY). Actual height is reset to the current mixing height
at the release point if greater than this minimum.
(HTMINBC)                      Default:  500.     ! HTMINBC = 500.0 !

Search radius (in BC segment lengths) about a receptor for sampling
nearest BC puff. BC puffs are emitted with a spacing of one segment
length, so the search radius should be greater than 1.
(RSAMPEC)                      Default:  4.       ! RSAMPEC = 10.0 !

Near-Surface depletion adjustment to concentration profile used when
sampling BC puffs?
(MDEPBC)                      Default:  1       ! MDEPBC = 1 !
  0 = Concentration is NOT adjusted for depletion
  1 = Adjust Concentration for depletion

!END!

-----

INPUT GROUPS: 13a, 13b, 13c, 13d -- Point source parameters
-----

-----
Subgroup (13a)
-----

Number of point sources with
parameters provided below      (NPT1) No default ! NPT1 = 1 !

Units used for point source
emissions below                (IPTU) Default: 1 ! IPTU = 1 !
  1 =      g/s
  2 =      kg/hr
  3 =      lb/hr
  4 =      tons/yr
  5 =      Odour Unit * m**3/s (vol. flux of odour compound)
  6 =      Odour Unit * m**3/min
  7 =      metric tons/yr

Number of source-species
combinations with variable
emissions scaling factors
provided below in (13d)        (NSPT1) Default: 0 ! NSPT1 = 0 !
```

Number of point sources with  
variable emission parameters  
provided in external file (NPT2) No default ! NPT2 = 0 !

(If NPT2 > 0, these point  
source emissions are read from  
the file: PTEMARB.DAT)

!END!

-----  
Subgroup (13b)  
-----

The following species are modeled:

CSPEC = SO2  
CSPEC = SO4  
CSPEC = NOX  
CSPEC = HNO3  
CSPEC = NO3  
CSPEC = PMC  
CSPEC = PMF  
CSPEC = EC  
CSPEC = SOA

a  
POINT SOURCE: CONSTANT DATA  
-----

Source No.	X Coordinate (km)	Y Coordinate (km)	Stack Height (m)	Base Elevation (m)	Stack Diameter (m)	Exit Vel. (m/s)	Exit Temp. (deg. K)	b Bldg. Dwash	c Emission Rates
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

```

1 ! SRCNAM = P1 !
1 ! X = 695.75, 507.2100, 88.390,216.46, 2.600, 20.0, 445.0, .0,
208.0E00, 8.23, 68.1E00, 0.0E00, 0.0E00, 0.37, 0.43, 0.02, 2.06 !
1 ! FMFAC = 1.0 ! !END!

```

a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.

SRCNAM is a 12-character name for a source  
(No default)  
X is an array holding the source data listed by the column headings  
(No default)  
SIGYZI is an array holding the initial sigma-y and sigma-z (m)  
(Default: 0.,0.)  
FMFAC is a vertical momentum flux factor (0. or 1.0) used to represent  
the effect of rain-caps or other physical configurations that  
reduce momentum rise associated with the actual exit velocity.  
(Default: 1.0 -- full momentum used)

b  
0. = No building downwash modeled, 1. = downwash modeled  
NOTE: must be entered as a REAL number (i.e., with decimal point)

c  
An emission rate must be entered for every pollutant modeled.

Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by IPTU (e.g. 1 for g/s).

-----  
Subgroup (13c)  
-----

BUILDING DIMENSION DATA FOR SOURCES SUBJECT TO DOWNWASH  
-----

Source No. Effective building height, width, length and X/Y offset (in meters) every 10 degrees. LENGTH, XBADJ, and YBADJ are only needed for MBDW=2 (PRIME downwash option) <sup>a</sup>

-----

<sup>a</sup> Building height, width, length, and X/Y offset from the source are treated as a separate input subgroup for each source and therefore must end with an input group terminator. The X/Y offset is the position, relative to the stack, of the center of the upwind face of the projected building, with the x-axis pointing along the flow direction.

-----  
Subgroup (13d)  
-----

<sup>a</sup>  
POINT SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 13b. Factors entered multiply the rates in 13b. Skip sources here that have constant emissions. For more elaborate variation in source parameters, use PTEMARB.DAT and NPT2 > 0.

IVARY determines the type of variation, and is source-specific:

(IVARY) Default: 0  
0 = Constant  
1 = Diurnal cycle (24 scaling factors: hours 1-24)  
2 = Monthly cycle (12 scaling factors: months 1-12)  
3 = Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)  
4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)  
5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----

<sup>a</sup> Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----

**Refined BART Modeling Protocol for  
Thilmany Kaukauna Mill**

---

INPUT GROUPS: 14a, 14b, 14c, 14d -- Area source parameters

-----  
Subgroup (14a)  
-----

Number of polygon area sources with  
parameters specified below (NAR1)           No default ! NAR1 = 0 !

Units used for area source  
emissions below                           (IARU)           Default: 1 ! IARU = 1 !

- 1 =           g/m\*\*2/s
- 2 =           kg/m\*\*2/hr
- 3 =           lb/m\*\*2/hr
- 4 =           tons/m\*\*2/yr
- 5 =           Odour Unit \* m/s (vol. flux/m\*\*2 of odour compound)
- 6 =           Odour Unit \* m/min
- 7 =           metric tons/m\*\*2/yr

Number of source-species  
combinations with variable  
emissions scaling factors  
provided below in (14d)                   (NSAR1) Default: 0 ! NSAR1 = 0 !

Number of buoyant polygon area sources  
with variable location and emission  
parameters (NAR2)                       No default ! NAR2 = 0 !  
(If NAR2 > 0, ALL parameter data for  
these sources are read from the file: BAEMARB.DAT)

!END!

-----  
Subgroup (14b)  
-----

a  
AREA SOURCE: CONSTANT DATA  
-----

Source No.	Effect. Height (m)	Base Elevation (m)	Initial Sigma z (m)	Emission Rates
-----	-----	-----	-----	-----

- a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.
- b  
An emission rate must be entered for every pollutant modeled.  
Enter emission rate of zero for secondary pollutants that are  
modeled, but not emitted. Units are specified by IARU  
(e.g. 1 for g/m\*\*2/s).

-----  
Subgroup (14c)  
-----

COORDINATES (km) FOR EACH VERTEX(4) OF EACH POLYGON  
-----

Source No.	Ordered list of X followed by list of Y, grouped by source	a
-----	-----	-----



-----  
a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.

-----  
Subgroup (14d)  
-----

a  
AREA SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission  
rates given in 14b. Factors entered multiply the rates in 14b.  
Skip sources here that have constant emissions. For more elaborate  
variation in source parameters, use BAEMARB.DAT and NAR2 > 0.

IVARY determines the type of variation, and is source-specific:

(IVARY) Default: 0  
0 = Constant  
1 = Diurnal cycle (24 scaling factors: hours 1-24)  
2 = Monthly cycle (12 scaling factors: months 1-12)  
3 = Hour & Season (4 groups of 24 hourly scaling factors,  
where first group is DEC-JAN-FEB)  
4 = Speed & Stab. (6 groups of 6 scaling factors, where  
first group is Stability Class A,  
and the speed classes have upper  
bounds (m/s) defined in Group 12  
5 = Temperature (12 scaling factors, where temperature  
classes have upper bounds (C) of:  
0, 5, 10, 15, 20, 25, 30, 35, 40,  
45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup  
and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 15a, 15b, 15c -- Line source parameters  
-----

-----  
Subgroup (15a)  
-----

Number of buoyant line sources  
with variable location and emission  
parameters (NLN2) No default ! NLN2 = 0 !

(If NLN2 > 0, ALL parameter data for  
these sources are read from the file: LNEMARB.DAT)

Number of buoyant line sources (NLINES) No default ! NLINES = 0 !

Units used for line source  
emissions below (ILNU) Default: 1 ! ILNU = 1 !  
1 = g/s

2 = kg/hr  
 3 = lb/hr  
 4 = tons/yr  
 5 = Odour Unit \* m\*\*3/s (vol. flux of odour compound)  
 6 = Odour Unit \* m\*\*3/min  
 7 = metric tons/yr

Number of source-species combinations with variable emissions scaling factors provided below in (15c) (NSLN1) Default: 0 ! NSLN1 = 0 !

Maximum number of segments used to model each line (MXNSEG) Default: 7 ! MXNSEG = 7 !

The following variables are required only if NLINES > 0. They are used in the buoyant line source plume rise calculations.

Number of distances at which transitional rise is computed Default: 6 ! NLRISE = 6 !

Average building length (XL) No default ! XL = .0 ! (in meters)

Average building height (HBL) No default ! HBL = .0 ! (in meters)

Average building width (WBL) No default ! WBL = .0 ! (in meters)

Average line source width (WML) No default ! WML = .0 ! (in meters)

Average separation between buildings (DXL) No default ! DXL = .0 ! (in meters)

Average buoyancy parameter (FPRIMEL) No default ! FPRIMEL = .0 ! (in m\*\*4/s\*\*3)

!END!

-----  
 Subgroup (15b)  
 -----

BUOYANT LINE SOURCE: CONSTANT DATA  
 -----

Source No.	Beg. X Coordinate (km)	Beg. Y Coordinate (km)	End. X Coordinate (km)	End. Y Coordinate (km)	Release Height (m)	Base Elevation (m)	Emission Rates
-----	-----	-----	-----	-----	-----	-----	-----

a  
 Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.

b  
 An emission rate must be entered for every pollutant modeled. Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by ILNTU (e.g. 1 for g/s).

-----  
Subgroup (15c)  
-----

a  
BUOYANT LINE SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 15b. Factors entered multiply the rates in 15b. Skip sources here that have constant emissions.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

0 =	Constant
1 =	Diurnal cycle (24 scaling factors: hours 1-24)
2 =	Monthly cycle (12 scaling factors: months 1-12)
3 =	Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)
4 =	Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)
5 =	Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 16a, 16b, 16c -- Volume source parameters  
-----

-----  
Subgroup (16a)  
-----

Number of volume sources with parameters provided in 16b,c (NVL1) No default ! NVL1 = 0 !

Units used for volume source emissions below in 16b (IVLU) Default: 1 ! IVLU = 1 !

1 =	g/s
2 =	kg/hr
3 =	lb/hr
4 =	tons/yr
5 =	Odour Unit * m**3/s (vol. flux of odour compound)
6 =	Odour Unit * m**3/min
7 =	metric tons/yr

Number of source-species combinations with variable emissions scaling factors provided below in (16c) (NSVL1) Default: 0 ! NSVL1 = 0 !

Number of volume sources with variable location and emission

parameters (NVL2) No default ! NVL2 = 0 !

(If NVL2 > 0, ALL parameter data for  
these sources are read from the VOLEMARB.DAT file(s) )

!END!

-----  
Subgroup (16b)  
-----

a  
VOLUME SOURCE: CONSTANT DATA  
-----

X	Y	Effect.	Base	Initial	Initial	b
Coordinate	Coordinate	Height	Elevation	Sigma y	Sigma z	Emission
(km)	(km)	(m)	(m)	(m)	(m)	Rates
-----	-----	-----	-----	-----	-----	-----

-----  
a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.

b  
An emission rate must be entered for every pollutant modeled.  
Enter emission rate of zero for secondary pollutants that are  
modeled, but not emitted. Units are specified by IVLU  
(e.g. 1 for g/s).

-----  
Subgroup (16c)  
-----

a  
VOLUME SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission  
rates given in 16b. Factors entered multiply the rates in 16b.  
Skip sources here that have constant emissions. For more elaborate  
variation in source parameters, use VOLEMARB.DAT and NVL2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

0 =	Constant
1 =	Diurnal cycle (24 scaling factors: hours 1-24)
2 =	Monthly cycle (12 scaling factors: months 1-12)
3 =	Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)
4 =	Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)
5 =	Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup  
and therefore must end with an input group terminator.

---

-----  
INPUT GROUPS: 17a & 17b -- Non-gridded (discrete) receptor information  
-----

-----  
Subgroup (17a)  
-----

Number of non-gridded receptors (NREC) No default ! NREC = 1995 !

!END!

-----  
Subgroup (17b)  
-----

-----  
NON-GRIDDED (DISCRETE) RECEPTOR DATA<sup>a</sup>  
-----

Receptor No.	X Coordinate (km)	Y Coordinate (km)	Ground Elevation (m)	Height Above Ground (m)
-----------------	-------------------------	-------------------------	----------------------------	-------------------------------

Note - three Class I areas  
Seney SENE = 173 receptors  
Isle Royal ISLE = 966 receptors  
Boundary Waters BOWA = 856 receptors  
total number of receptors is 1995

ALL RECEPTOR COORDINATES FOLLOW IN INPUT FILE

Updated CALPUFF Modeling  
Protocol and Final Report for the  
Subject to BART Analyses  
for Wausau Paper Mills, LLC  
Mosinee Mill

Prepared by

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December 2010

# Updated CALPUFF Modeling Protocol and Final Report for Wausau “Subject to BART” Analysis

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Appendix A – Example CALMET Input File

Appendix B – Example CALPUFF Input File

## 1.0 Introduction

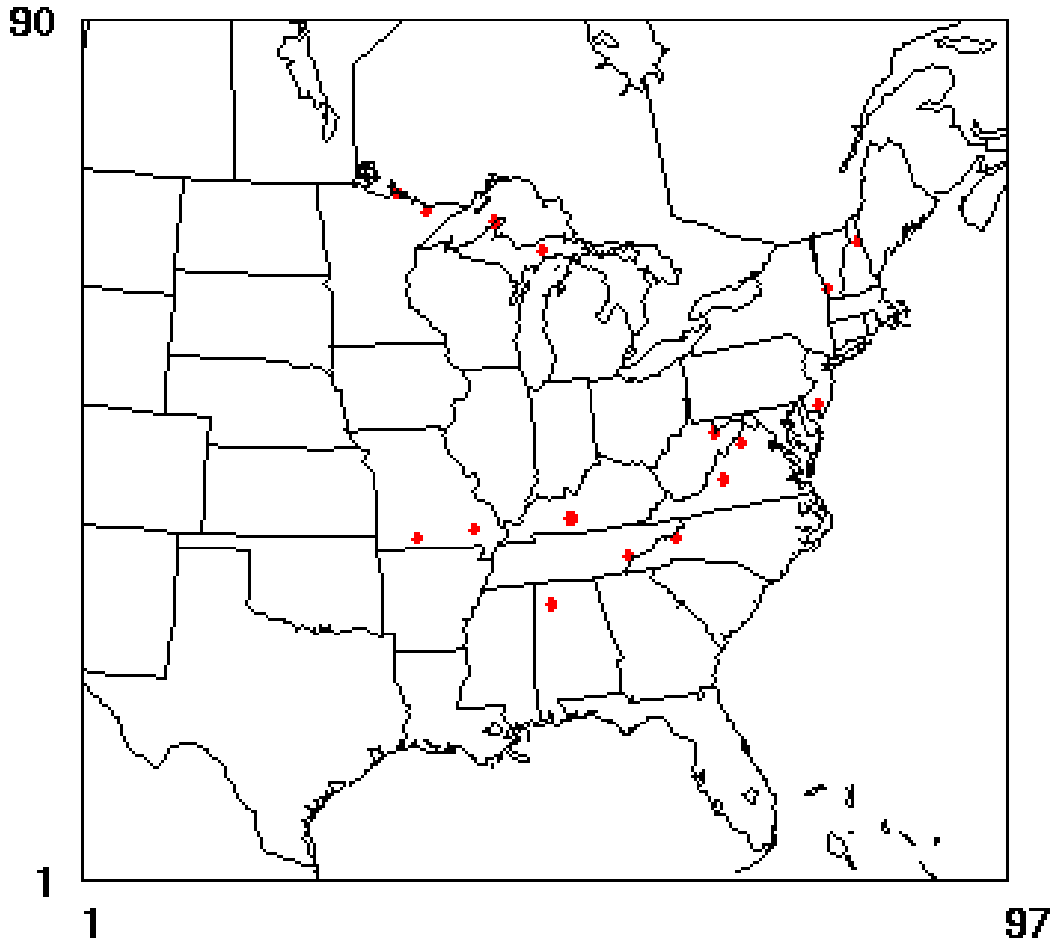
The Wisconsin Department of Natural Resources (DNR) performed single source screening modeling in 2006 to evaluate which Best Available Retrofit Technology (BART) eligible sources in the state may “cause or contribute” to visibility impairment and could therefore be subject to BART requirements. This type of analysis is referred to as the “subject to BART” or “BART exemption” analysis. The methods used by DNR are described in the “Single Source Modeling to Support Regional Haze BART Modeling Protocol”, March 21, 2006, prepared by Lake Michigan Air Directors Consortium (LADCO).

DNR applied CALPUFF to each BART eligible source for three annual periods, covering the calendar years 2002 to 2004. The CALPUFF/CALMET modeling domain consisted of ninety-seven 36-km spaced cells in the east-west direction and ninety 36-km spaced cells in the north-south direction (see Figure 1). Meteorological data generated by the MM5 prognostic weather model (provided by LADCO in CALMM5 format) was used to develop the meteorological data for CALPUFF, without additional observational data input (i.e., NOOBS mode). This represents a coarse meteorological grid, suitable for an initial screening analysis.

The results from DNR’s screening analysis indicated that, with the exception of some of the largest electric power stations in the state, the only Class I areas where BART eligible sources may contribute to visibility impairment are the three areas to the north of Wisconsin, the Boundary Waters Canoe Area, Isle Royal National Park, and Seney Wildlife Area. DNR’s screening analysis also indicated that the BART eligible emissions unit at the Wausau Paper Mills LLC Mosinee Mill (Wausau) may “cause or contribute” to visibility impairment at the Seney Class I area (located approximately 320 km to the northeast of the mill), as defined by the 98<sup>th</sup> percentile visibility impacts being greater than 0.5 deciviews (dv) relative to natural background.



Figure 1 - WDNR Screening CALPUFF Domain



NOTE: Class I areas are identified by red marks.

Wausau has elected to refine the DNR screening analysis by conducting refined CALPUFF modeling with a 4 km meteorological grid, based on procedures described in the “Visibility Improvement – State and Tribal Association of the Southeast (VISTAS) BART CALPUFF” modeling protocol. The VISTAS protocol was developed by a large group of stakeholders, including EPA, the VISTAS member state agencies and tribes, the Federal Land Managers (FLMs), industry representatives, and consulting experts. Given the resources expended on the VISTAS protocol, it arguably represents the most comprehensive CALPUFF BART protocol ever developed.

Wausau had discussions with DNR and EPA Region 5 staff about refining the analysis, and submitted a refined CALPUFF BART modeling protocol in March 2007. EPA reviewed the protocol and provided minor comments to DNR in May 2010. Another facility undergoing Subject to BART review, Thilmany, responded to these comments in a letter “Response to EPA Comments on Thilmany BART Modeling Protocol” dated November 9, 2010. In addition, subsequent to the EPA protocol comments, the US Fish and Wildlife Service and US Forest Service reviewed the protocol and had additional discussions with DNR and EPA. The three EPA comments, and each response, are listed below.

**Comment 1**

Information needs to be presented which clarifies how POSTUTIL was used for ammonia.

**Response 1**

The VISTAS methodology was used, with a MNITRATE switch setting of one (1).

**Comment 2**

The modeling must include a copy of the input and list files associated with runs. The protocol only offers some of the control file information.

**Response 2**

All CALMET, CALPUFF, and CALPOST input and list files, as well as meteorological and geophysical data files, had been provided by Wausau to DNR on a portable hard disk drive. Wausau will again provide a portable hard disk drive of all modeling files to DNR for the updated BART analysis, for purposes of SIP documentation.

**Comment 3**

The input file ... uses the background ammonia values that were produced for the LADCO protocol. These values are based on modeled estimates. Monitoring data has generally shown much higher ammonia than what was used from the model results. The refined modeling analyses ... must evaluate the visibility impacts using measured background ammonia values.

**Response 3**

The US Fish and Wildlife Service proposed using a constant background ammonia concentration of 3.9 ppb, which was the highest concentration measured at the Mayville Wisconsin station during a mid-1990s EPA study. Wausau has agreed to use this single, conservative background value in the absence of daily data.

On November 29, 2010, DNR issued a letter that approved the CALPUFF BART protocol, as well as the emission rates to be used for the BART eligible emission units. DNR requested that the 2007 version of the protocol be updated to include the responses to EPA comments, and text that addresses current CALPUFF modeling guidance. In addition, the updated modeling results could be presented in the protocol (so that all information is located in one document).

This document is the updated modeling protocol and final modeling report for the refined Wausau “Subject to BART” analysis. This document does not duplicate the extensive information on visibility, CALPUFF, and EPA guidance on BART modeling procedures that are contained in the VISTAS protocol. Rather, this document describes how data specific to this setting will be used in the refined CALPUFF modeling. Section 1 of the protocol is this introduction. Section 2

briefly summarizes important concepts from the VISTAS protocol. Section 3 describes the specific CALMET data and procedures that were used to develop a refined, 4-km spaced meteorological grid for Wisconsin. Section 4 describes the CALPUFF input data and procedures, and Section 5 discusses the POSTUTIL and CALPOST processing, including the use of annual average natural background visibility. Section 6 presents the emission and stack data that were used for the BART eligible emission units, and Section 7 presents the modeling results. Appendices A and B present example CALMET and CALPUFF input file listings. A portable hard drive is also being submitted that contains all input data, control files, output files, and computer codes used in the analysis.

## **2.0 VISTAS BART Protocol**

The VISTAS member states have prepared a CALPUFF modeling protocol for BART determinations that is fully consistent with the EPA guidelines in 40 CFR Part 51 Appendix W and Appendix Y. The VISTAS protocol describes a two-step CALPUFF analysis methodology for BART exemption analyses. The initial CALPUFF analysis uses a “screening” or coarse resolution meteorological grid to determine if a particular source may be exempted from further BART analyses. The screening results are also used to determine which Class I areas should be included in any refined analysis. Assumptions for the initial screening analysis are conservative so that a source that contributes to visibility impairment is not exempted in error. If a source is shown to contribute to visibility impairment using the initial screening assessment, the source has the option to undertake refined CALPUFF modeling using finer meteorological grids to evaluate further whether it is subject to BART.

VISTAS developed both coarse (12-km grid resolution) and refined (4-km grid resolution) CALMET files. The fine grid CALMET files utilized both MM5 prognostic data, as well as meteorological observational data for the CALMET “Step 2” calculations. Overwater (buoy) data were used in addition to the standard hourly surface meteorological observations, precipitation

observations, and twice-daily upper air sounding data. The VISTAS website provides detailed documentation and supporting information on the refined CALMET processing, including the CALMET and CALPUFF input files that contain model configuration options and settings that have been optimized after consultation with stakeholders and consulting experts. These same CALMET and CALPUFF input files were used as the basis for the refined Wisconsin analyses.

The VISTAS protocol discusses several options for defining the natural background visibility. These include use of the annual average natural background extinction or the background extinction for the 20% best natural conditions days. Based on an email survey of the state agencies from the VISTAS member states, all of the VISTAS states allowed the use of annual average natural background extinction for refined CALPUFF analyses. This is also the guidance provided by EPA Region 5 staff during discussions with DNR and the Wisconsin Paper Council, and the method approved by DNR in their November 29, 2010 letter. Therefore, the refined Wausau CALPUFF modeling used the annual average natural background extinction for CALPOST processing.

The VISTAS BART guidance recommends that the threshold value used to define whether a source “contributes” to visibility impairment is a 0.5 dv change from natural background conditions. The 98<sup>th</sup> percentile 24-hr average predicted impact at the Class I area (equal to the 8<sup>th</sup> highest value) is to be compared to this contribution threshold value. According to clarification of the BART guidance received from EPA by the VISTAS workgroup, for a three-year simulation the modeling values to be compared with the threshold are the greatest of the three annual 8<sup>th</sup> highest values, or the 22<sup>nd</sup> highest value over all three years combined, whichever is greater.

### 3.0 CALMET Data and Procedures

EPA has updated the versions of the approved CALPUFF programs since the 2007 modeling protocol was developed, and has provided additional guidance on CALMET “switch” settings. The current approved versions are CALPUFF version 5.8 (level 070623), and CALMET version 5.8 (level 070623), and these versions were used for the updated refined BART CALPUFF analysis. On August 31, 2009, EPA published the memorandum “Clarification of EPA-FLM Recommended Settings for CALMET”. EPA stated that a 4 km fine CALMET grid size was acceptable, and provided recommended CALMET settings. The CALMET settings for the updated refined BART CALPUFF analysis meet all current EPA recommendations.

The first step in development of the refined Wisconsin analysis was to define the CALMET refined modeling domain and grid. As discussed in Section 1, DNR’s screening analysis indicated that there are three Class I areas to the north of Wisconsin, the Boundary Waters Canoe Area, Isle Royal National Park, and Seney Wildlife Area, where the screening analysis predicts some level of visibility impact from Wisconsin BART eligible sources. Therefore, the meteorological grid was developed to ensure coverage for these three Class I areas, as well as for the BART eligible source locations throughout Wisconsin. Standard IWAQM guidance also recommends that the CALMET grid extend a minimum of 50 km beyond any Class I or emission source location, so that puff recirculation can be properly accounted for.

The refined CALMET grid was defined as a 4-km spaced grid with 162 cells in the X-axis and 191 in the Y-axis. The grid projection was Lambert Conformal Conic, using the same projection parameters as in the DNR screening analysis (this allows the use of the same source X and Y coordinates as in the DNR analysis). The projection parameters include RLAT0 = 40N, RLON0 = 97W, XLAT1 = 33N, XLAT2 = 45N, XORIGKM = 247.0, YORIGKM = 240.0, and DATUM = NAS-C.

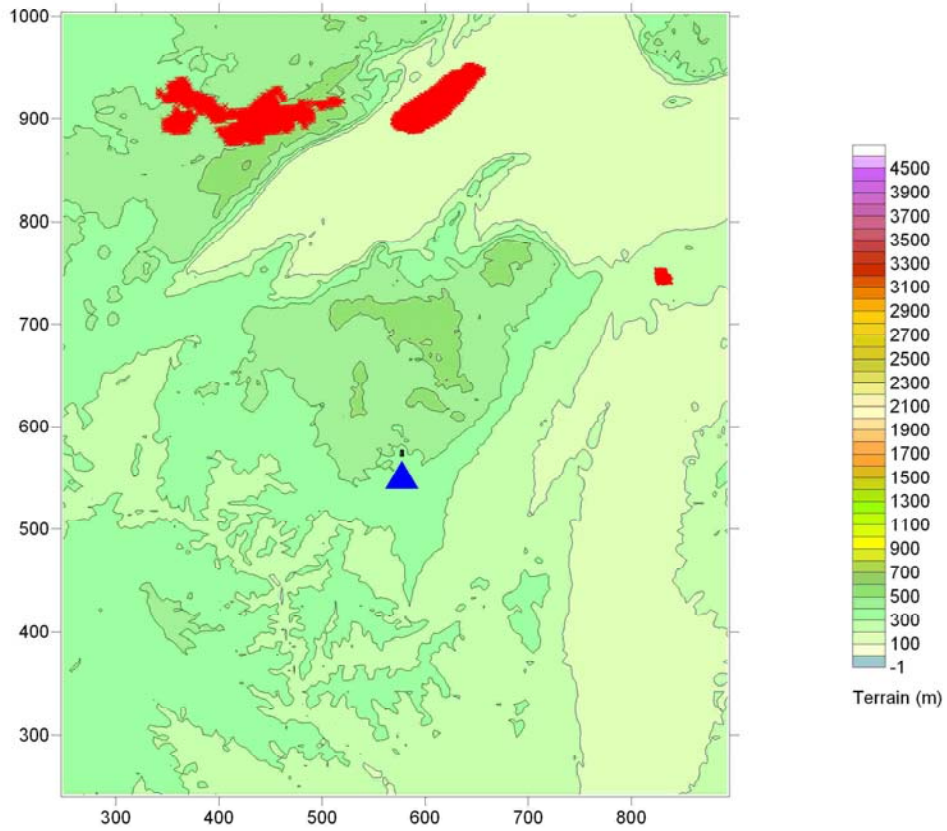
The meteorological data that was input to CALMET included the same three years (2002-2004) of CALMM5 data used by DNR, which was provided by LADCO. However, the refined CALMET runs also considered observation data from 22 surface stations, 16 precipitation stations, 5 overwater buoy stations, and 3 upper air stations. Figure 2 is a plot of the refined CALPUFF modeling domain. Figure 3 presents a plot of the locations of the observation stations in the modeling domain.

The CALMET program MAKEGEO uses USGS GTOPO30 digital terrain data (which includes coverage in Canada) and USGS Composite Theme Grid Land Use and Land Cover (LULC) data to create a gridded data file on geophysical parameters that is used by CALMET. The available USGS LULC coverage does not extend into the Canadian areas of the refined domain. These areas in Canada were assigned to the “forested land” category. Figure 4 presents a plot of the land use data for the modeling domain.

The VISTAS refined CALMET input files discussed in section 2 were used as the basis for the refined Wisconsin CALMET input files. The same CALMET configuration settings and options were utilized, as they are representative of Wisconsin terrain and land use patterns. All settings conform with current EPA and FLM guidance.

Appendix A presents a listing of a CALMET input file developed for the refined analysis.

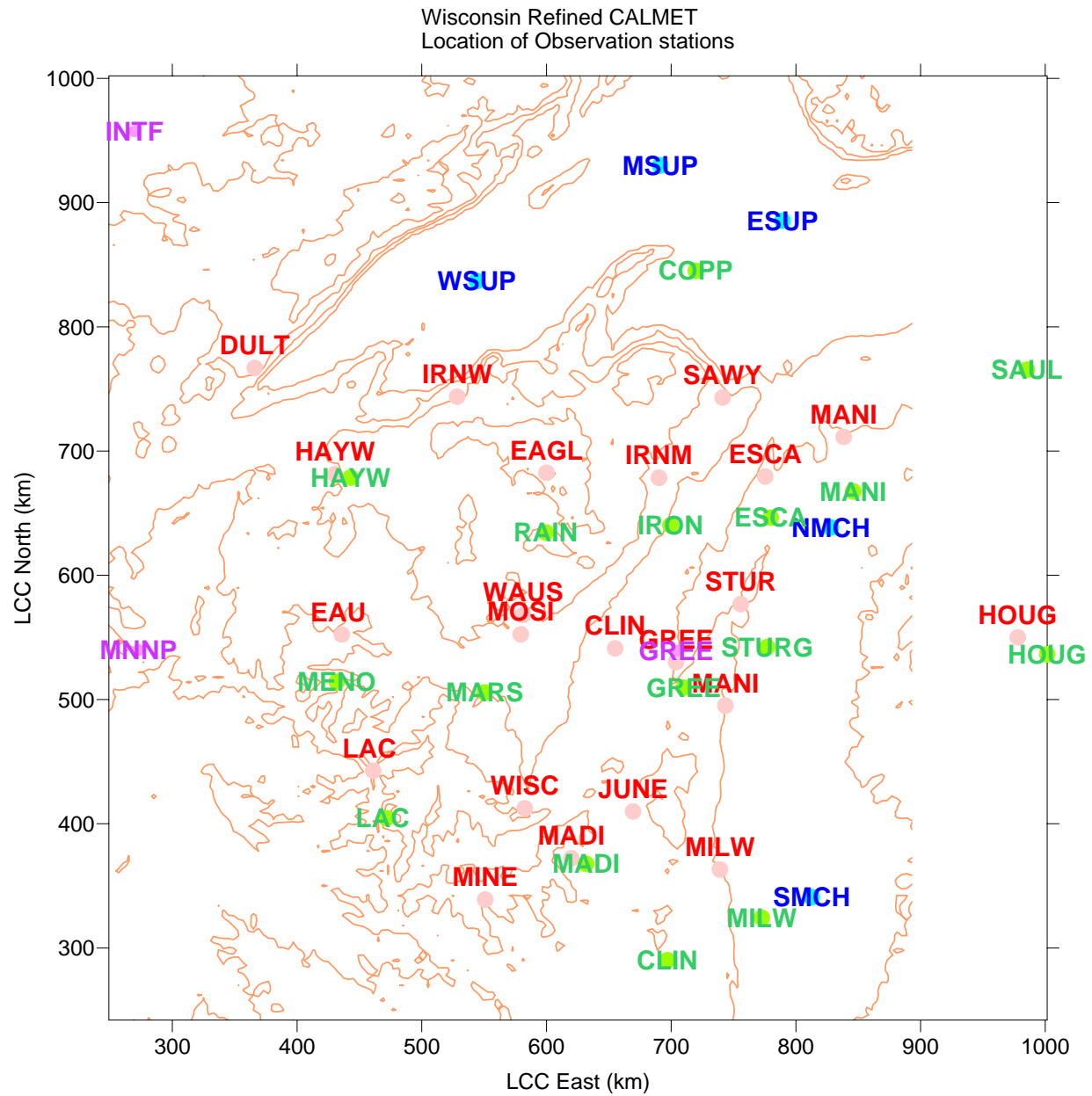
Figure 2 – Refined CALPUFF Modeling Domain and Locations of Class I Areas



NOTE: The Class I area receptors are in red, the Wausau Mill location is in blue, and the X and Y coordinates are the Lambert projection coordinates are in kms.

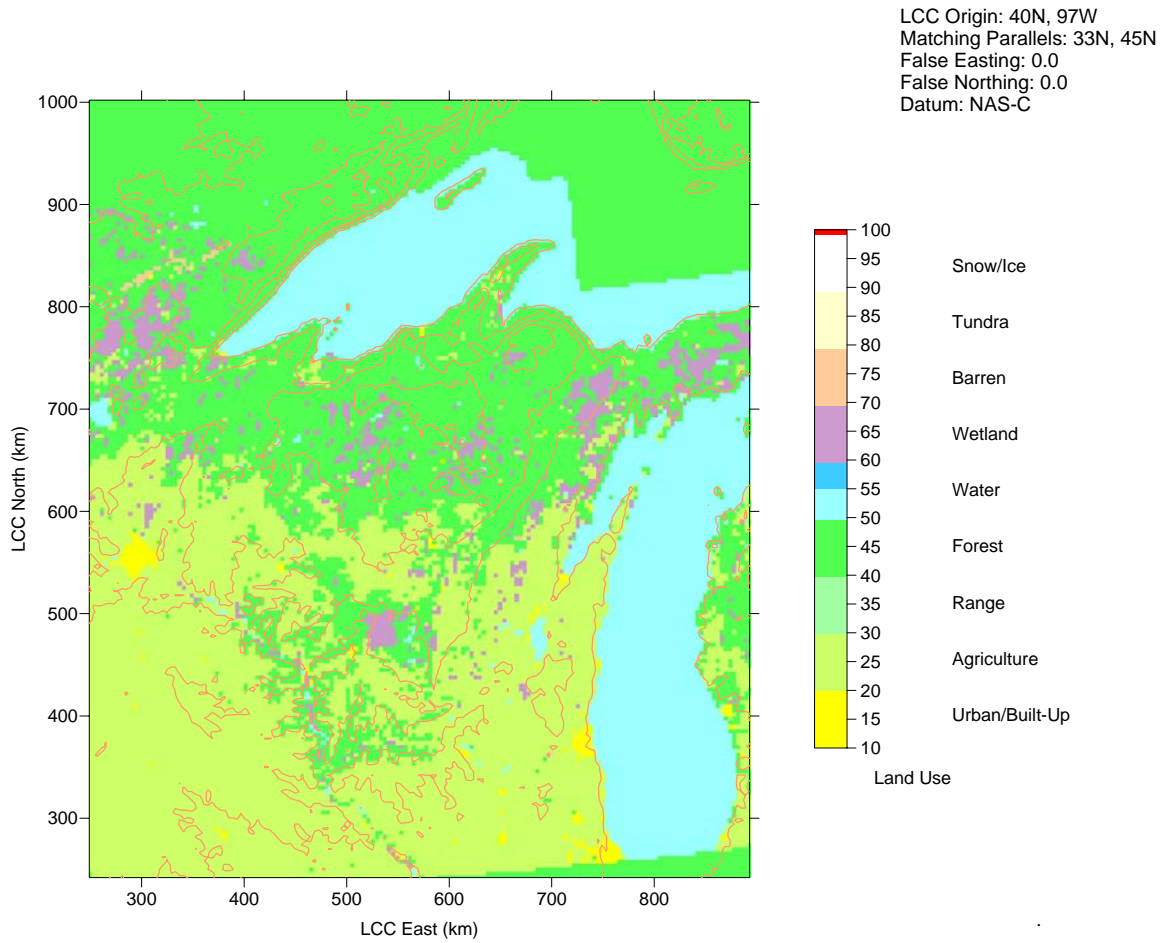


Figure 3 – Locations of Observational Stations in Refined CALPUFF Modeling Domain



NOTES: Surface stations are in red, overwater (buoy) in blue, precipitation in green, and upper air in purple.

Figure 4 – Land Use Data for Refined CALPUFF Modeling Domain



Some of the notable CALMET configuration settings included:

- CALMET vertical layers set to 10, with cell face heights at 20, 40, 80, 160, 320, 640, 1200, 2000, 3000, and 4000 meters,
- CALMET diagnostic options, including slope flow effects (ISLOPE = 1) with the radius of influence of terrain (TERRAD) set to 15 km,
- EPA recommended CALMET settings for the radius of influence parameters (R1=50, R2=100, RMAX1=100, RMAX2=200, and RMAX3=200).

## 4.0 CALPUFF Procedures

Both the DNR screening and VISTAS refined CALPUFF input files were used as the basis for the refined Wisconsin CALPUFF input files. All CALPUFF settings and options are consistent with EPA and FLM guidance, and the refined VISTAS protocol recommendations. The background ozone (O<sub>3</sub>) concentrations were used directly from the DNR CALPUFF input files. Instead of using monthly background ammonia concentrations from the DNR screening analyses, a single conservative background ammonia concentration of 3.9 ppb was used. Appendix B presents a listing of a CALPUFF input file developed for the analysis.

The DNR screening analysis considered three emitted pollutants, SO<sub>2</sub>, NO<sub>x</sub>, and PM<sub>2.5</sub>. The refined CALPUFF modeling will consider speciated particulate emissions, including coarse PM (PMC), fine PM equivalent to PM<sub>2.5</sub> (PMF), sulfate (SO<sub>4</sub>), nitrate (NO<sub>3</sub>), elemental carbon (EC), and secondary organic aerosols (SOA or OC). Section 6 of this protocol describes how the speciated particulate emissions were calculated.

## 5.0 POSTUTIL and CALPOST Procedures

In CALPUFF, the scavenging of the available ammonia for nitrate formation is computed on a puff-by-puff basis. Because of the possibility of multiple puff overlap, this approach may tend to overestimate the amount of ammonia available to form nitrate. In order to address the cumulative consumption of ammonia, the POSTUTIL program can be used to repartition nitric acid and nitrate. POSTUTIL calculates new NO<sub>3</sub>/HNO<sub>3</sub> equilibrium concentrations for the sample source in the Class I area using the background ammonia. As recommended in the VISTAS procedures for refined CALPUFF modeling, the POSTUTIL “MNITRATE=1” method was used for the refined Wisconsin CALPUFF modeling. Instead of using the monthly values for background ammonia concentrations from the DNR screening analyses, a single conservative background ammonia concentration of 3.9 ppb was used.

The modeling protocol submitted in 2007 proposed to use CALPOST Visibility Method 6 to calculate visibility impacts. However, since that time there have been revisions to CALPOST that allow the use of the updated IMPROVE extinction equation. CALPOST version 6.221 can implement the “Method 8 Mode 5” methodology to calculate visibility impacts, and the FLMs recommend its use in “Federal Land Managers’ Air Quality Related Values Work Group (FLAG) Phase I Report—Revised (2010)” (herein referred to as FLAG 2010). Therefore, this updated methodology will be used for the Wausau BART analysis. The new CALPUFF Graphical Users Interface automatically enters the appropriate values, based on the Class I area being selected, for natural background species and relative humidity adjustment factors using the data in Tables 6 through 9 of Section 3.3.5 of FLAG 2010 (which are based on “Guidance for Estimating Natural Visibility Conditions Under the Regional Haze Rule”, EPA-454/B-03-005, September 2003).

## 6.0 Emissions Modeled

EPA's BART modeling guidance state that the 24-hr maximum actual emissions rates for the period of the modeling simulation (2002-2004) should be used for the analysis. If this data is not available, then the short term "allowable" or "potential" emission rates should be used. Wausau provided DNR with source information that has been reviewed to determine the 24-hr maximum actual emissions rates.

There are a number of BART eligible emission units at the Wausau Mosinee mill. Stack S10 is a combined stack for the 3 boilers that are the primary source of the BART pollutants, and the recovery boiler (B21) vents to stack S11. In addition, although the BART eligibility is in question for some smaller emissions sources (the lime kiln, smelt tank, and slaker), these units were also included in the modeling to be conservative. Based on discussions between RMT (Wausau Paper's consultant) and DNR on December 7 and 9, 2010, it was agreed to use PTE emissions for all emission units, as listed in Table 1. Note that the PTE for S10 was adjusted to reflect fuel oil usage limitations as specified in the air permit. The PM emissions have been speciated using the information for coal and oil fired boilers developed by the National Park Service<sup>1</sup>.

---

<sup>1</sup> <http://www2.nature.nps.gov/air/permits/ect/index.cfm>

Table 1 – Wausau Emissions (in units of g/sec)

<b>Unit</b>	<b>SO2</b>	<b>NOx</b>	<b>PM</b>	<b>PMC</b>	<b>PMF</b>	<b>EC</b>	<b>SO4</b>	<b>SOA</b>
S10	143.1	28.1	11.68	1.16	2.39	0.09	6.43	1.61
S11	13.1	5.4	2.32	0.73	1.26	0.10	0.19	0.04
S12	0.4	0.0	0.43	0.00	0.43	0.00	0.00	0.00
S52	0.0	0.0	0.43	0.00	0.43	0.00	0.00	0.00
S53	1.4	0.9	0.62	0.32	0.19	0.01	0.10	0.00
Total	158.0	34.4	15.5	2.2	4.7	0.2	6.7	1.7

## 7.0 Modeling Results

CALPOST computes the daily maximum change in deciviews relative to the annual average natural background visibility, separately for each Class I area. For evaluating the source impact relative to the “Subject to BART” visibility impact threshold of 0.5 dv, the daily maximum changes in visibility are sorted to determine the 98th percentile (8<sup>th</sup> highest) value for each of three years, and the 98th percentile (22<sup>nd</sup> highest) value for the completed three year period. If all 98th percentile values are below the threshold of 0.5 dv based on the refined CALPUFF analysis, the source is exempt from the BART requirements.

Another equivalent metric commonly used for BART exemption modeling is to sum the number of days per year that are above the 0.5 dv threshold for each Class I area separately; if there are 7 or fewer days above the threshold for each year, then the 98<sup>th</sup> percentile value is also below the threshold. This is the metric approved by DNR for use in the BART modeling analysis.

Specifically, for each Class I area separately, if there are less than 8 days per year that the predicted visibility impacts are greater than 0.5 deciviews relative to the annual average natural background visibility, then the BART eligible emission units at the modeled facility will be determined not to be Subject to BART.

Table 2 presents the results from the refined CALPUFF BART exemption modeling. These results demonstrate that there are less than 8 days per year, at each individual Class I area, when the visibility impacts from the Wausau mill BART eligible emission unit are above 0.5 dv. Therefore, the **Wausau mill BART eligible sources are not Subject to BART** requirements.

Table 2 – Wausau CALPUFF BART Exemption Results Summary

<b>Class I area</b>	<b>Distance (km) from source to Class I area boundary</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2002</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2003</b>	<b># of days with impact &gt; 0.5 dv in Class I area: 2004</b>	<b># of days with impact &gt; 0.5 dv in Class I area for 3 year period</b>	<b>Maximum 98<sup>th</sup> Percentile impact over either annual or 3-yr periods</b>
Seney , MI	~ 320	0	0	1	1	0.20
Isle Royal, MN	~ 340	0	0	0	0	0.15
Boundary Waters, MN	~ 360	2	0	0	2	0.13



**Appendix A – Example CALMET Input File**

WI Refined CALMET - 4km Grid - Dec 2010 - with Updated EPA Settings

----- Run title (3 lines) -----  
CALMET MODEL CONTROL FILE  
-----

INPUT GROUP: 0 -- Input and Output File Names

Subgroup (a)

Default Name	Type	File Name
-----	----	-----
GEO.DAT	input	! GEODAT=geo.dat !
SURF.DAT	input	! SRFDAT=SURF02Z.DAT !
CLOUD.DAT	input	* CLDDAT= *
PRECIP.DAT	input	! PRCDAT=PRECIP02.DAT !
MM4.DAT	input	* MM4DAT=E:\MRPO-calmm5\2002\MM5.020102.3D.DAT *
MM4.DAT	input	! MM4DAT=E:\MRPO-calmm5\2002\020102.MM5 !
WT.DAT	input	* WTDAT= *
CALMET.LST	output	! METLST=cal020102.lst !
CALMET.DAT	output	! METDAT=E:\WI-Calmet\2002Output\cal020102.dat !
PACOUT.DAT	output	* PACDAT= *

All file names will be converted to lower case if LCFILES = T  
Otherwise, if LCFILES = F, file names will be converted to UPPER CASE  
T = lower case ! LCFILES = T !  
F = UPPER CASE

NUMBER OF UPPER AIR & OVERWATER STATIONS:

Number of upper air stations (NUSTA) No default ! NUSTA = 3 !  
Number of overwater met stations  
(NOWSTA) No default ! NOWSTA = 5 !

!END!

-----  
Subgroup (b)

---

-----  
Upper air files (one per station)  
-----

Default Name	Type	File Name
UP1.DAT	input	1 ! UPDAT=UPGB0204.DAT! !END!
UP2.DAT	input	2 ! UPDAT=UPMN0204.DAT! !END!
UP3.DAT	input	3 ! UPDAT=UPIN0204.DAT! !END!

-----  
Subgroup (c)  
-----

Overwater station files (one per station)

SEA1.DAT	input	1 ! SEADAT=4500102.DAT! !END!
SEA2.DAT	input	2 ! SEADAT=4500202.DAT! !END!
SEA3.DAT	input	3 ! SEADAT=4500402.DAT! !END!
SEA4.DAT	input	4 ! SEADAT=4500602.DAT! !END!
SEA5.DAT	input	5 ! SEADAT=4500702.DAT! !END!

-----  
Default Name Type File Name  
-----  
-----

-----  
Subgroup (d)  
-----

Other file names  
-----

Default Name	Type	File Name
DIAG.DAT	input	* DIADAT= *
PROG.DAT	input	* PRGDAT= *
TEST.PRT	output	* TSTPRT= *
TEST.OUT	output	* TSTOUT= *
TEST.KIN	output	* TSTKIN= *
TEST.FRD	output	* TSTFRD= *
TEST.SLP	output	* TSTSLP= *

-----  
NOTES: (1) File/path names can be up to 70 characters in length  
(2) Subgroups (a) and (d) must have ONE 'END' (surround by  
delimiters) at the end of the group  
(3) Subgroups (b) and (c) must have an 'END' (surround by

delimiters) at the end of EACH LINE

!END!

-----  
INPUT GROUP: 1 -- General run control parameters  
-----

Starting date: Year (IBYR) -- No default ! IBYR= 2002 !  
Month (IBMO) -- No default ! IBMO= 1 !  
Day (IBDY) -- No default ! IBDY= 2 !  
Hour (IBHR) -- No default ! IBHR= 0 !

Base time zone (IBTZ) -- No default ! IBTZ= 6 !  
PST = 08, MST = 07  
CST = 06, EST = 05

Length of run (hours) (IRLG) -- No default ! IRLG= 24 !

Run type (IRTYPE) -- Default: 1 ! IRTYPE= 1 !

0 = Computes wind fields only  
1 = Computes wind fields and micrometeorological variables  
(u\*, w\*, L, zi, etc.)  
(IRTYPE must be 1 to run CALPUFF or CALGRID)

Compute special data fields required  
by CALGRID (i.e., 3-D fields of W wind  
components and temperature)  
in addition to regular Default: T ! LCALGRD = T !  
fields ? (LCALGRD)  
(LCALGRD must be T to run CALGRID)

Flag to stop run after  
SETUP phase (ITEST) Default: 2 ! ITEST= 2 !  
(Used to allow checking  
of the model inputs, files, etc.)  
ITEST = 1 - STOPS program after SETUP phase  
ITEST = 2 - Continues with execution of  
COMPUTATIONAL phase after SETUP

Test options specified to see if  
they conform to regulatory

values? (MREG)                      No Default                      ! MREG = 1                      !

0 = NO checks are made

1 = Technical options must conform to USEPA guidance

          IMIXH    -1            Maul-Carson convective mixing height  
                                  over land; OCD mixing height overwater  
          ICOARE    0            OCD deltaT method for overwater fluxes  
          THRESHL  0.0           Threshold buoyancy flux over land needed  
                                  to sustain convective mixing height grow

!END!

-----

INPUT GROUP: 2 -- Map Projection and Grid control parameters

-----

Projection for all (X,Y):

-----

Map projection

(PMAP)                                      Default: UTM                      ! PMAP = LCC                      !

UTM : Universal Transverse Mercator  
TTM : Tangential Transverse Mercator  
LCC : Lambert Conformal Conic  
PS : Polar Stereographic  
EM : Equatorial Mercator  
LAZA : Lambert Azimuthal Equal Area

False Easting and Northing (km) at the projection origin

(Used only if PMAP= TTM, LCC, or LAZA)

(FEAST)                                      Default=0.0                      ! FEAST = 0.000                      !  
(FNORTH)                                     Default=0.0                      ! FNORTH = 0.000                      !

UTM zone (1 to 60)

(Used only if PMAP=UTM)

(IUTMZN)                                     No Default                      ! IUTMZN = -999                      !

Hemisphere for UTM projection?

(Used only if PMAP=UTM)

(UTMHEM)                                     Default: N                      ! UTMHEM = N                      !

N : Northern hemisphere projection  
S : Southern hemisphere projection

Latitude and Longitude (decimal degrees) of projection origin

(Used only if PMAP= TTM, LCC, PS, EM, or LAZA)

(RLAT0)                      No Default              ! RLAT0 = 40N    !

(RLON0)                      No Default              ! RLON0 = 97W    !

TTM : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience

LCC : RLON0 identifies central (true N/S) meridian of projection  
      RLAT0 selected for convenience

PS : RLON0 identifies central (grid N/S) meridian of projection  
      RLAT0 selected for convenience

EM : RLON0 identifies central meridian of projection  
      RLAT0 is REPLACED by 0.0N (Equator)

LAZA: RLON0 identifies longitude of tangent-point of mapping plane  
      RLAT0 identifies latitude of tangent-point of mapping plane

Matching parallel(s) of latitude (decimal degrees) for projection

(Used only if PMAP= LCC or PS)

(XLAT1)                      No Default              ! XLAT1 = 33N    !

(XLAT2)                      No Default              ! XLAT2 = 45N    !

LCC : Projection cone slices through Earth's surface at XLAT1 and XLAT2

PS : Projection plane slices through Earth at XLAT1  
      (XLAT2 is not used)

-----

Note: Latitudes and longitudes should be positive, and include a  
      letter N,S,E, or W indicating north or south latitude, and  
      east or west longitude. For example,  
      35.9 N Latitude = 35.9N  
      118.7 E Longitude = 118.7E

Datum-region

-----

The Datum-Region for the coordinates is identified by a character  
string. Many mapping products currently available use the model of the  
Earth known as the World Geodetic System 1984 (WGS-84). Other local  
models may be in use, and their selection in CALMET will make its output  
consistent with local mapping products. The list of Datum-Regions with  
official transformation parameters is provided by the National Imagery and  
Mapping Agency (NIMA).

NIMA Datum - Regions(Examples)

-----  
WGS-84 WGS-84 Reference Ellipsoid and Geoid, Global coverage (WGS84)  
NAS-C NORTH AMERICAN 1927 Clarke 1866 Spheroid, MEAN FOR CONUS (NAD27)  
NAR-C NORTH AMERICAN 1983 GRS 80 Spheroid, MEAN FOR CONUS (NAD83)  
NWS-84 NWS 6370KM Radius, Sphere  
ESR-S ESRI REFERENCE 6371KM Radius, Sphere

Datum-region for output coordinates

(DATUM) Default: WGS-84 ! DATUM = NAS-C !

Horizontal grid definition:

-----  
Rectangular grid defined for projection PMAP,  
with X the Easting and Y the Northing coordinate

No. X grid cells (NX) No default ! NX = 162 !  
No. Y grid cells (NY) No default ! NY = 191 !

Grid spacing (DGRIDKM) No default ! DGRIDKM = 4. !  
Units: km

Reference grid coordinate of  
SOUTHWEST corner of grid cell (1,1)

X coordinate (XORIGKM) No default ! XORIGKM = 247.000 !  
Y coordinate (YORIGKM) No default ! YORIGKM = 240.000 !  
Units: km

Vertical grid definition:

-----  
No. of vertical layers (NZ) No default ! NZ = 10 !

Cell face heights in arbitrary  
vertical grid (ZFACE(NZ+1)) No defaults  
Units: m

! ZFACE = 0.,20.,40.,80.,160.,320.,640.,1200.,2000.,3000.,4000. !

!END!

-----  
INPUT GROUP: 3 -- Output Options  
-----

DISK OUTPUT OPTION

Save met. fields in an unformatted  
output file ? (LSAVE) Default: T ! LSAVE = T !  
(F = Do not save, T = Save)

Type of unformatted output file:  
(IFORMO) Default: 1 ! IFORMO = 1 !

- 1 = CALPUFF/CALGRID type file (CALMET.DAT)
- 2 = MESOPUFF-II type file (PACOUT.DAT)

LINE PRINTER OUTPUT OPTIONS:

Print met. fields ? (LPRINT) Default: F ! LPRINT = T !  
(F = Do not print, T = Print)  
(NOTE: parameters below control which  
met. variables are printed)

Print interval  
(IPRINF) in hours Default: 1 ! IPRINF = 1 !  
(Meteorological fields are printed  
every 1 hours)

Specify which layers of U, V wind component  
to print (IUVOU(NZ)) -- NOTE: NZ values must be entered  
(0=Do not print, 1=Print)  
(used only if LPRINT=T) Defaults: NZ\*0  
! IUVOU = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !  
-----

Specify which levels of the W wind component to print  
(NOTE: W defined at TOP cell face -- 16 values)  
(IWOUT(NZ)) -- NOTE: NZ values must be entered

---



(0=Do not print, 1=Print)  
(used only if LPRINT=T & LCALGRD=T)  
-----  
Defaults: NZ\*0  
! IWOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Specify which levels of the 3-D temperature field to print  
(ITOUT(NZ)) -- NOTE: NZ values must be entered  
(0=Do not print, 1=Print)  
(used only if LPRINT=T & LCALGRD=T)  
-----  
Defaults: NZ\*0  
! ITOUT = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Specify which meteorological fields  
to print  
(used only if LPRINT=T) Defaults: 0 (all variables)  
-----

Variable	Print ?	
	(0 = do not print,	
	1 = print)	
-----	-----	
! STABILITY =	0	! - PGT stability class
! USTAR =	0	! - Friction velocity
! MONIN =	0	! - Monin-Obukhov length
! MIXHT =	0	! - Mixing height
! WSTAR =	0	! - Convective velocity scale
! PRECIP =	0	! - Precipitation rate
! SENSHEAT =	0	! - Sensible heat flux
! CONVZI =	0	! - Convective mixing ht.

Testing and debug print options for micrometeorological module

Print input meteorological data and  
internal variables (LDB) Default: F ! LDB = F !  
(F = Do not print, T = print)  
(NOTE: this option produces large amounts of output)

First time step for which debug data

are printed (NN1)                      Default: 1              ! NN1 = 1 !

Last time step for which debug data  
are printed (NN2)                      Default: 1              ! NN2 = 2 !

Testing and debug print options for wind field module  
(all of the following print options control output to  
wind field module's output files: TEST.PRT, TEST.OUT,  
TEST.KIN, TEST.FRD, and TEST.SLP)

Control variable for writing the test/debug  
wind fields to disk files (IOUTD)  
(0=Do not write, 1=write)              Default: 0              ! IOUTD = 0 !

Number of levels, starting at the surface,  
to print (NZPRN2)                      Default: 1              ! NZPRN2 = 1 !

Print the INTERPOLATED wind components ?  
(IPR0) (0=no, 1=yes)                      Default: 0              ! IPR0 = 0 !

Print the TERRAIN ADJUSTED surface wind  
components ?  
(IPR1) (0=no, 1=yes)                      Default: 0              ! IPR1 = 0 !

Print the SMOOTHED wind components and  
the INITIAL DIVERGENCE fields ?  
(IPR2) (0=no, 1=yes)                      Default: 0              ! IPR2 = 0 !

Print the FINAL wind speed and direction  
fields ?  
(IPR3) (0=no, 1=yes)                      Default: 0              ! IPR3 = 0 !

Print the FINAL DIVERGENCE fields ?  
(IPR4) (0=no, 1=yes)                      Default: 0              ! IPR4 = 0 !

Print the winds after KINEMATIC effects  
are added ?  
(IPR5) (0=no, 1=yes)                      Default: 0              ! IPR5 = 0 !

Print the winds after the FROUDE NUMBER  
adjustment is made ?  
(IPR6) (0=no, 1=yes)                      Default: 0              ! IPR6 = 0 !

Print the winds after SLOPE FLOWS  
are added ?

(IPR7) (0=no, 1=yes)                    Default: 0            ! IPR7 = 0 !

Print the FINAL wind field components ?

(IPR8) (0=no, 1=yes)                    Default: 0            ! IPR8 = 0 !

!END!

-----  
INPUT GROUP: 4 -- Meteorological data options  
-----

NO OBSERVATION MODE                    (NOOBS) Default: 0            ! NOOBS = 0 !

0 = Use surface, overwater, and upper air stations

1 = Use surface and overwater stations (no upper air observations)  
    Use MM5 for upper air data

2 = No surface, overwater, or upper air observations

    Use MM5 for surface, overwater, and upper air data

NUMBER OF SURFACE & PRECIP. METEOROLOGICAL STATIONS

Number of surface stations    (NSSTA) No default            ! NSSTA = 22 !

Number of precipitation stations

(NPSTA=-1: flag for use of MM5 precip data)

(NPSTA) No default            ! NPSTA = 16 !

CLOUD DATA OPTIONS

Gridded cloud fields:

(ICLOUD) Default: 0            ! ICLOUD = 0 !

ICLOUD = 0 - Gridded clouds not used

ICLOUD = 1 - Gridded CLOUD.DAT generated as OUTPUT

ICLOUD = 2 - Gridded CLOUD.DAT read as INPUT

ICLOUD = 3 - Gridded cloud cover from Prognostic Rel. Humidity

FILE FORMATS

Surface meteorological data file format

(IFORMS) Default: 2            ! IFORMS = 2 !

(1 = unformatted (e.g., SMERGE output))

(2 = formatted    (free-formatted user input))

Precipitation data file format

(IFORMP) Default: 2 ! IFORMP = 2 !

(1 = unformatted (e.g., PMERGE output))

(2 = formatted (free-formatted user input))

Cloud data file format

(IFORMC) Default: 2 ! IFORMC = 2 !

(1 = unformatted - CALMET unformatted output)

(2 = formatted - free-formatted CALMET output or user input)

!END!

-----  
INPUT GROUP: 5 -- Wind Field Options and Parameters  
-----

WIND FIELD MODEL OPTIONS

Model selection variable (IWFCOD) Default: 1 ! IWFCOD = 1 !

0 = Objective analysis only

1 = Diagnostic wind module

Compute Froude number adjustment

effects ? (IFRADJ) Default: 1 ! IFRADJ = 1 !

(0 = NO, 1 = YES)

Compute kinematic effects ? (IKINE) Default: 0 ! IKINE = 0 !

(0 = NO, 1 = YES)

Use O'Brien procedure for adjustment

of the vertical velocity ? (IOBR) Default: 0 ! IOBR = 0 !

(0 = NO, 1 = YES)

Compute slope flow effects ? (ISLOPE) Default: 1 ! ISLOPE = 1 !

(0 = NO, 1 = YES)

Extrapolate surface wind observations

to upper layers ? (IEXTRP) Default: -4 ! IEXTRP = -4 !

(1 = no extrapolation is done,

2 = power law extrapolation used,

3 = user input multiplicative factors

for layers 2 - NZ used (see FEXTRP array)  
4 = similarity theory used  
-1, -2, -3, -4 = same as above except layer 1 data  
at upper air stations are ignored

Extrapolate surface winds even

if calm? (ICALM) Default: 0 ! ICALM = 0 !  
(0 = NO, 1 = YES)

Layer-dependent biases modifying the weights of  
surface and upper air stations (BIAS(NZ))

-1<=BIAS<=1

Negative BIAS reduces the weight of upper air stations

(e.g. BIAS=-0.1 reduces the weight of upper air stations  
by 10%; BIAS= -1, reduces their weight by 100 %)

Positive BIAS reduces the weight of surface stations

(e.g. BIAS= 0.2 reduces the weight of surface stations  
by 20%; BIAS=1 reduces their weight by 100%)

Zero BIAS leaves weights unchanged (1/R\*\*2 interpolation)

Default: NZ\*0

! BIAS = 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 , 0 !

Minimum distance from nearest upper air station  
to surface station for which extrapolation  
of surface winds at surface station will be allowed  
(RMIN2: Set to -1 for IEXTRP = 4 or other situations  
where all surface stations should be extrapolated)

Default: 4. ! RMIN2 = -1.0 !

Use gridded prognostic wind field model  
output fields as input to the diagnostic

wind field model (IPROG) Default: 0 ! IPROG = 14 !  
(0 = No, [IWFCOD = 0 or 1])

1 = Yes, use CSUMM prog. winds as Step 1 field, [IWFCOD = 0]

2 = Yes, use CSUMM prog. winds as initial guess field [IWFCOD = 1]

3 = Yes, use winds from MM4.DAT file as Step 1 field [IWFCOD = 0]

4 = Yes, use winds from MM4.DAT file as initial guess field [IWFCOD = 1]

5 = Yes, use winds from MM4.DAT file as observations [IWFCOD = 1]

13 = Yes, use winds from MM5.DAT file as Step 1 field [IWFCOD = 0]

14 = Yes, use winds from MM5.DAT file as initial guess field [IWFCOD = 1]

15 = Yes, use winds from MM5.DAT file as observations [IWFCOD = 1]

Timestep (hours) of the prognostic

model input data (ISTEPPG) Default: 1 ! ISTEPPG = 1 !

---



(Used only if IPROG = 1) Units: km  
-----

Maximum acceptable divergence in the  
divergence minimization procedure  
(DIVLIM)

Default: 5.E-6 ! DIVLIM= 5.0E-06 !

Maximum number of iterations in the  
divergence min. procedure (NITER)

Default: 50 ! NITER = 50 !

Number of passes in the smoothing  
procedure (NSMTH(NZ))

NOTE: NZ values must be entered

Default: 2,(mxnz-1)\*4 ! NSMTH =

2 , 4 , 4 , 4 , 4 , 4 , 4 , 4 , 4 , 4 , 4 !

Maximum number of stations used in  
each layer for the interpolation of  
data to a grid point (NINTR2(NZ))

NOTE: NZ values must be entered

Default: 99. ! NINTR2 =

99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 , 99 !

Critical Froude number (CRITFN)

Default: 1.0 ! CRITFN = 1. !

Empirical factor controlling the  
influence of kinematic effects

(ALPHA)

Default: 0.1 ! ALPHA = 0.1 !

Multiplicative scaling factor for  
extrapolation of surface observations

to upper layers (FEXTR2(NZ))

Default: NZ\*0.0

! FEXTR2 = 0., 0., 0., 0., 0., 0., 0., 0., 0., 0. !

(Used only if IEXTRP = 3 or -3)

#### BARRIER INFORMATION

Number of barriers to interpolation  
of the wind fields (NBAR)

Default: 0 ! NBAR = 0 !

THE FOLLOWING 4 VARIABLES ARE INCLUDED

ONLY IF NBAR > 0

NOTE: NBAR values must be entered

No defaults

for each variable

Units: km

X coordinate of BEGINNING  
of each barrier (XBBAR(NBAR))       ! XBBAR = 0. !  
Y coordinate of BEGINNING  
of each barrier (YBBAR(NBAR))       ! YBBAR = 0. !  
  
X coordinate of ENDING  
of each barrier (XEBAR(NBAR))       ! XEBAR = 0. !  
Y coordinate of ENDING  
of each barrier (YEBAR(NBAR))       ! YEBAR = 0. !

DIAGNOSTIC MODULE DATA INPUT OPTIONS

Surface temperature (IDIOPT1)       Default: 0       ! IDIOPT1 = 0 !  
0 = Compute internally from  
    hourly surface observations  
1 = Read preprocessed values from  
    a data file (DIAG.DAT)

Surface met. station to use for  
the surface temperature (ISURFT)   No default       ! ISURFT = 1 !  
(Must be a value from 1 to NSSTA)  
(Used only if IDIOPT1 = 0)

-----

Domain-averaged temperature lapse  
rate (IDIOPT2)                      Default: 0       ! IDIOPT2 = 0 !  
0 = Compute internally from  
    twice-daily upper air observations  
1 = Read hourly preprocessed values  
    from a data file (DIAG.DAT)

Upper air station to use for  
the domain-scale lapse rate (IUPT) No default       ! IUPT = 1 !  
(Must be a value from 1 to NUSTA)  
(Used only if IDIOPT2 = 0)

-----

Depth through which the domain-scale  
lapse rate is computed (ZUPT)       Default: 200.   ! ZUPT = 200. !  
(Used only if IDIOPT2 = 0)       Units: meters

-----



Domain-averaged wind components

(IDIOPT3) Default: 0 ! IDIOPT3 = 0 !

0 = Compute internally from  
twice-daily upper air observations  
1 = Read hourly preprocessed values  
a data file (DIAG.DAT)

Upper air station to use for

the domain-scale winds (IUPWND) Default: -1 ! IUPWND = -1 !

(Must be a value from -1 to NUSTA)

(Used only if IDIOPT3 = 0)

-----

Bottom and top of layer through

which the domain-scale winds

are computed

(ZUPWND(1), ZUPWND(2)) Defaults: 1., 1000. ! ZUPWND= 1., 1000. !

(Used only if IDIOPT3 = 0) Units: meters

-----

Observed surface wind components

for wind field module (IDIOPT4) Default: 0 ! IDIOPT4 = 0 !

0 = Read WS, WD from a surface  
data file (SURF.DAT)

1 = Read hourly preprocessed U, V from  
a data file (DIAG.DAT)

Observed upper air wind components

for wind field module (IDIOPT5) Default: 0 ! IDIOPT5 = 0 !

0 = Read WS, WD from an upper  
air data file (UP1.DAT, UP2.DAT, etc.)

1 = Read hourly preprocessed U, V from  
a data file (DIAG.DAT)

LAKE BREEZE INFORMATION

Use Lake Breeze Module (LLBREZE)

Default: F ! LLBREZE = F !

Number of lake breeze regions (NBOX) ! NBOX = 0 !

X Grid line 1 defining the region of interest

! XG1 = 0. !

X Grid line 2 defining the region of interest

```

! XG2 = 0. !
Y Grid line 1 defining the region of interest
! YG1 = 0. !
Y Grid line 2 defining the region of interest
! YG2 = 0. !

X Point defining the coastline (Straight line)
(XBCST) (KM) Default: none ! XBCST = 0. !

Y Point defining the coastline (Straight line)
(YBCST) (KM) Default: none ! YBCST = 0. !

X Point defining the coastline (Straight line)
(XECST) (KM) Default: none ! XECST = 0. !

Y Point defining the coastline (Straight line)
(YECST) (KM) Default: none ! YECST = 0. !

Number of stations in the region Default: none ! NLB = 0 !
(Surface stations + upper air stations)

Station ID's in the region (METBXID(NLB))
(Surface stations first, then upper air stations)
! METBXID = 0 !

!END!

-----

INPUT GROUP: 6 -- Mixing Height, Temperature and Precipitation Parameters
-----

EMPIRICAL MIXING HEIGHT CONSTANTS

Neutral, mechanical equation
(CONSTB) Default: 1.41 ! CONSTB = 1.41 !
Convective mixing ht. equation
(CONSTE) Default: 0.15 ! CONSTE = 0.15 !
Stable mixing ht. equation
(CONSTN) Default: 2400. ! CONSTN = 2400.!
Overwater mixing ht. equation
(CONSTW) Default: 0.16 ! CONSTW = 0.16 !
```

---

Absolute value of Coriolis  
parameter (FCORIOI)                      Default: 1.E-4   ! FCORIOI = 1.0E-04!  
Units: (1/s)

SPATIAL AVERAGING OF MIXING HEIGHTS

Conduct spatial averaging  
(IAVEZI) (0=no, 1=yes)                    Default: 1       ! IAVEZI = 1 !

Max. search radius in averaging  
process (MNMDAV)                          Default: 1       ! MNMDAV = 1 !  
Units: Grid  
      cells

Half-angle of upwind looking cone  
for averaging (HAFANG)                    Default: 30.     ! HAFANG = 30. !  
Units: deg.

Layer of winds used in upwind  
averaging (ILEVZI)                        Default: 1       ! ILEVZI = 1 !  
(must be between 1 and NZ)

CONVECTIVE MIXING HEIGHT OPTIONS:

Method to compute the convective  
mixing height(IMIXH)                      Default: 1       ! IMIXH = -1 !  
    1: Maul-Carson for land and water cells  
   -1: Maul-Carson for land cells only -  
      OCD mixing height overwater  
    2: Batchvarova and Gryning for land and water cells  
   -2: Batchvarova and Gryning for land cells only  
      OCD mixing height overwater

Threshold buoyancy flux required to  
sustain convective mixing height growth  
overland (THRESHL)                        Default: 0.05   ! THRESHL = 0.00 !  
(expressed as a heat flux                units: W/m3  
per meter of boundary layer)

Threshold buoyancy flux required to  
sustain convective mixing height growth  
overwater (THRESHW)                       Default: 0.05   ! THRESHW = 0.05 !  
(expressed as a heat flux                units: W/m3  
per meter of boundary layer)

Option for overwater lapse rates used  
in convective mixing height growth

(ITWPROG) Default: 0 ! ITWPROG = 0 !  
0 : use SEA.DAT lapse rates and deltaT (or assume neutral  
conditions if missing)  
1 : use prognostic lapse rates (only if IPROG>2)  
and SEA.DAT deltaT (or neutral if missing)  
2 : use prognostic lapse rates and prognostic delta T  
(only if iprog>12 and 3D.DAT version# 2.0 or higher)

Land Use category ocean in 3D.DAT datasets

(ILUOC3D) Default: 16 ! ILUOC3D = 16 !  
Note: if 3D.DAT from MM5 version 3.0, iluoc3d = 16  
if MM4.DAT, typically iluoc3d = 7

#### OTHER MIXING HEIGHT VARIABLES

Minimum potential temperature lapse  
rate in the stable layer above the  
current convective mixing ht. Default: 0.001 ! DPTMIN = 0.001 !  
(DPTMIN) Units: deg. K/m

Depth of layer above current conv.  
mixing height through which lapse  
rate is computed (DZZI) Default: 200. ! DZZI = 200. !  
Units: meters

Minimum overland mixing height Default: 50. ! ZIMIN = 50. !  
(ZIMIN) Units: meters

Maximum overland mixing height Default: 3000. ! ZIMAX = 3000. !  
(ZIMAX) Units: meters

Minimum overwater mixing height Default: 50. ! ZIMINW = 50. !  
(ZIMINW) -- (Not used if observed  
overwater mixing hts. are used) Units: meters

Maximum overwater mixing height Default: 3000. ! ZIMAXW = 3000. !  
(ZIMAXW) -- (Not used if observed  
overwater mixing hts. are used) Units: meters

#### OVERWATER SURFACE FLUXES METHOD and PARAMETERS

(ICOARE) Default: 10 ! ICOARE = 0 !  
0: original deltaT method (OCD)  
10: COARE with no wave parameterization (jwave=0, Charnock)  
11: COARE with wave option jwave=1 (Oost et al.)  
and default wave properties  
-11: COARE with wave option jwave=1 (Oost et al.)  
and observed wave properties (must be in SEA.DAT files)  
12: COARE with wave option 2 (Taylor and Yelland)  
and default wave properties

-12: COARE with wave option 2 (Taylor and Yelland)  
and observed wave properties (must be in SEA.DAT files)

Coastal/Shallow water length scale (DSHELF)  
(for modified z0 in shallow water)  
( COARE fluxes only)

Default : 0. ! DSHELF = 0. !  
units: km

COARE warm layer computation (IWARM) ! IWARM = 0 !  
1: on - 0: off (must be off if SST measured with  
IR radiometer) Default: 0

COARE cool skin layer computation (ICOOL) ! ICOOL = 0 !  
1: on - 0: off (must be off if SST measured with  
IR radiometer) Default: 0

#### TEMPERATURE PARAMETERS

3D temperature from observations or  
from prognostic data? (ITPROG) Default:0 !ITPROG = 0 !

0 = Use Surface and upper air stations  
(only if NOOBS = 0)  
1 = Use Surface stations (no upper air observations)  
Use MM5 for upper air data  
(only if NOOBS = 0,1)  
2 = No surface or upper air observations  
Use MM5 for surface and upper air data  
(only if NOOBS = 0,1,2)

Interpolation type  
(1 = 1/R ; 2 = 1/R\*\*2) Default:1 ! IRAD = 1 !

Radius of influence for temperature  
interpolation (TRADKM) Default: 500. ! TRADKM = 500. !  
Units: km

Maximum Number of stations to include  
in temperature interpolation (NUMTS) Default: 5 ! NUMTS = 5 !

Conduct spatial averaging of temp-  
eratures (IAVET) (0=no, 1=yes) Default: 1 ! IAVET = 1 !



**Refined BART Modeling Protocol for  
Wausau Mosinee Mill**

```

! SS3 = 'LA C'      26430      461.012      442.705      6      10  !
! SS4 = 'EAU '     26435      435.857      552.352      6      10  !
! SS5 = 'GREE'    26450      703.911      530.179      6      10  !
! SS6 = 'MANI'    26455      743.477      495.113      6      10  !
! SS7 = 'STUR'    26458      755.730      576.760      6      10  !
! SS8 = 'WAUS'    26463      581.202      568.247      6      10  !
! SS9 = 'MOSI'    26465      579.307      552.489      6      10  !
! SS10 = 'CLIN'   26502      655.019      541.168      6      10  !
! SS11 = 'WISC'   26503      582.532      412.498      6      10  !
! SS12 = 'EAGL'   26504      599.980      682.545      6      10  !
! SS13 = 'MINE'   26507      550.817      338.941      6      10  !
! SS14 = 'HAYW'   26508      430.347      681.295      6      10  !
! SS15 = 'JUNE'   26509      669.429      409.931      6      10  !
! SS16 = 'dulu'   27450      366.011      766.977      6      10  !
! SS17 = 'MANI'   25408      838.315      711.433      5      10  !
! SS18 = 'SAWY'   26284      741.237      743.078      5      10  !
! SS19 = 'HOUG'   26380      977.905      549.977      5      10  !
! SS20 = 'ESCA'   26480      775.417      679.550      5      10  !
! SS21 = 'IRNM'   27437      690.168      678.511      5      10  !
! SS22 = 'IRNW'   27445      528.314      743.857      5      10  !

```

```

-----
1
  Four character string for station name
  (MUST START IN COLUMN 9)

```

```

2
  Five digit integer for station ID

```

!END!

```

-----
INPUT GROUP: 8 -- Upper air meteorological station parameters
-----

```

```

UPPER AIR STATION VARIABLES
(One record per station -- 0 records in all)

```

```

      1      2
      Name    ID      X coord.  Y coord.  Time zone
              (km)    (km)
-----
! US1 = 'GREE' 14898  703.911  539.179  6  !

```

```
! US2  ='MNNP'   94983   272.681   539.843   6   !
! US3  ='INTF'   14918   269.489   957.416   6   !
```

-----

```
1
  Four character string for station name
  (MUST START IN COLUMN 9)
```

```
2
  Five digit integer for station ID
```

!END!

-----

INPUT GROUP: 9 -- Precipitation station parameters

-----

```
PRECIPITATION STATION VARIABLES
(One record per station -- 0 records in all)
(NOT INCLUDED IF NPSTA = 0)
```

	1	2		
	Name	Station Code	X coord. (km)	Y coord. (km)
! PS1	'COPP'	201780	719.681	845.179 !
! PS2	'ESCA'	202626	779.850	646.454 !
! PS3	'HOUG'	203936	1001.500	536.427 !
! PS4	'IRON'	204090	699.421	640.283 !
! PS5	'MANI'	205073	845.933	667.499 !
! PS6	'SAUL'	207366	985.172	765.448 !
! PS7	'CLIN'	471667	696.889	290.151 !
! PS8	'GREE'	473269	709.933	509.579 !
! PS9	'HAYW'	473511	442.947	678.728 !
! PS10	'LAC '	474370	471.515	404.495 !
! PS11	'MADI'	474961	631.719	367.656 !
! PS12	'MARS'	475120	550.348	505.662 !
! PS13	'MENO'	475335	431.807	514.277 !
! PS14	'MILW'	475479	772.567	324.344 !
! PS15	'RAIN'	476939	599.369	634.460 !
! PS16	'STURG'	478267	776.964	542.174 !



-----  
1

Four character string for station name  
(MUST START IN COLUMN 9)

2

Six digit station code composed of state  
code (first 2 digits) and station ID (last  
4 digits)

!END!

**Appendix B – Example CALPUFF Input File**

Example BART Modeling  
2010 Updated BART Modeling for Wausau, WI -  
Using Refined 4km CALMET Data File and SENE, ISLE, BOWA Class I Receptors

----- Run title (3 lines) -----

CALPUFF MODEL CONTROL FILE  
-----

INPUT GROUP: 0 -- Input and Output File Names

-----  
Default Name    Type            File Name  
-----  
CALMET.DAT    input            \* METDAT =            \*  
          or  
ISCMET.DAT    input            \* ISCDAT =            \*  
          or  
PLMMET.DAT    input            \* PLMDAT =            \*  
          or  
PROFILE.DAT    input            \* PRFDAT =            \*  
SURFACE.DAT    input            \* SFCDAT =            \*  
RESTARTB.DAT  input            \* RSTARTB=            \*  
-----  
CALPUFF.LST    output            ! PUFLST =WausuaRef02.lst !  
CONC.DAT        output            ! CONDAT =wausuaRef02.dat !  
DFLX.DAT        output            \* DFDAT =            \*  
WFLX.DAT        output            \* WFDAT =            \*  
VISB.DAT        output            \* VISDAT =wausua02DNR.VIB \*  
RESTARTE.DAT    output            \* RSTARTE=            \*

Emission Files  
-----

PTEMARB.DAT    input            \* PTDAT =            \*  
VOLEMARB.DAT    input            \* VOLDAT =            \*  
BAEMARB.DAT    input            \* ARDAT =            \*  
LNEMARB.DAT    input            \* LNDAT =            \*

Other Files  
-----

OZONE.DAT        input            \* OZDAT =            \*  
VD.DAT            input            \* VDDAT =            \*  
CHEM.DAT        input            \* CHEMDAT=            \*  
H2O2.DAT        input            \* H2O2DAT=            \*  
HILL.DAT        input            \* HILDAT=            \*  
HILLRCT.DAT     input            \* RCTDAT=            \*  
COASTLN.DAT     input            \* CSTDAT=            \*  
FLUXBDY.DAT     input            \* BDYDAT=            \*  
BCON.DAT        input            \* BCNDAT=            \*  
DEBUG.DAT        output            \* DEBUG =            \*  
MASSFLX.DAT     output            \* FLXDAT=            \*  
MASSBAL.DAT     output            \* BALDAT=            \*  
FOG.DAT          output            \* FOGDAT=            \*

-----  
All file names will be converted to lower case if LCFILES = T  
Otherwise, if LCFILES = F, file names will be converted to UPPER CASE  
T = lower case            ! LCFILES = T !  
F = UPPER CASE

NOTE: (1) file/path names can be up to 70 characters in length

Provision for multiple input files  
-----

**Refined BART Modeling Protocol for  
Wausau Mosinee Mill**

---

```

Number of CALMET.DAT files for run (NMETDAT)
          Default: 1          ! NMETDAT = 364 !

Number of PTEMARB.DAT files for run (NPTDAT)
          Default: 0          ! NPTDAT = 0 !

Number of BAEMARB.DAT files for run (NARDAT)
          Default: 0          ! NARDAT = 0 !

Number of VOLEMARB.DAT files for run (NVOLDAT)
          Default: 0          ! NVOLDAT = 0 !

```

!END!

```

-----
Subgroup (0a)
-----

```

The following CALMET.DAT filenames are processed in sequence if NMETDAT>1

Default Name	Type	File Name		
-----	----	-----		
none	input	! METDAT=e:\WI-Calmet\2002output\cal020102.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020103.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020104.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020105.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020106.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020107.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020108.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020109.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020110.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020111.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020112.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020113.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020114.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020115.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020116.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020117.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020118.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020119.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020120.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020121.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020122.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020123.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020124.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020125.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020126.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020127.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020128.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020129.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020130.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020131.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020201.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020202.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020203.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020204.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020205.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020206.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020207.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020208.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020209.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020210.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020211.DAT	!	!END!
none	input	! METDAT=e:\WI-Calmet\2002output\cal020212.DAT	!	!END!













INPUT GROUP: 1 -- General run control parameters

-----

Option to run all periods found  
in the met. file (METRUN) Default: 0 ! METRUN = 0 !

METRUN = 0 - Run period explicitly defined below  
METRUN = 1 - Run all periods in met. file

Starting date: Year (IBYR) -- No default ! IBYR = 2002 !  
(used only if Month (IBMO) -- No default ! IBMO = 1 !  
METRUN = 0) Day (IBDY) -- No default ! IBDY = 02 !  
Hour (IBHR) -- No default ! IBHR = 1 !

Base time zone (XBTZ) -- No default ! XBTZ = 6.0 !  
PST = 8., MST = 7.  
CST = 6., EST = 5.

Length of run (hours) (IRLG) -- No default ! IRLG = 8736 !

Number of chemical species (NSPEC)  
Default: 5 ! NSPEC = 9 !

Number of chemical species  
to be emitted (NSE) Default: 3 ! NSE = 7 !

Flag to stop run after  
SETUP phase (ITEST) Default: 2 ! ITEST = 2 !  
(Used to allow checking  
of the model inputs, files, etc.)  
ITEST = 1 - STOPS program after SETUP phase  
ITEST = 2 - Continues with execution of program  
after SETUP

Restart Configuration:

Control flag (MRESTART) Default: 0 ! MRESTART = 0 !

0 = Do not read or write a restart file  
1 = Read a restart file at the beginning of  
the run  
2 = Write a restart file during run  
3 = Read a restart file at beginning of run  
and write a restart file during run

Number of periods in Restart  
output cycle (NRESPD) Default: 0 ! NRESPD = 0 !

0 = File written only at last period  
>0 = File updated every NRESPD periods

Meteorological Data Format (METFM)  
Default: 1 ! METFM = 1 !

METFM = 1 - CALMET binary file (CALMET.MET)  
METFM = 2 - ISC ASCII file (ISCMET.MET)  
METFM = 3 - AUSPLUME ASCII file (PLMMET.MET)  
METFM = 4 - CTDM plus tower file (PROFILE.DAT) and  
surface parameters file (SURFACE.DAT)

PG sigma-y is adjusted by the factor (AVET/PGTIME)\*\*0.2  
Averaging Time (minutes) (AVET)  
Default: 60.0 ! AVET = 60. !

PG Averaging Time (minutes) (PGTIME)  
Default: 60.0 ! PGTIME = 60. !

!END!

-----  
INPUT GROUP: 2 -- Technical options  
-----

Vertical distribution used in the  
near field (MGAUSS)                   Default: 1       ! MGAUSS = 1   !  
  0 = uniform  
  1 = Gaussian

Terrain adjustment method  
(MCTADJ)                               Default: 3       ! MCTADJ = 3   !  
  0 = no adjustment  
  1 = ISC-type of terrain adjustment  
  2 = simple, CALPUFF-type of terrain  
      adjustment  
  3 = partial plume path adjustment

Subgrid-scale complex terrain  
flag (MCTSG)                           Default: 0       ! MCTSG = 0   !  
  0 = not modeled  
  1 = modeled

Near-field puffs modeled as  
elongated 0 (MSLUG)                   Default: 0       ! MSLUG = 0   !  
  0 = no  
  1 = yes (slug model used)

Transitional plume rise modeled ?  
(MTRANS)                              Default: 1       ! MTRANS = 1   !  
  0 = no (i.e., final rise only)  
  1 = yes (i.e., transitional rise computed)

Stack tip downwash? (MTIP)           Default: 1       ! MTIP = 1   !  
  0 = no (i.e., no stack tip downwash)  
  1 = yes (i.e., use stack tip downwash)

Method used to simulate building  
downwash? (MBDW)                      Default: 1       ! MBDW = 1   !  
  1 = ISC method  
  2 = PRIME method

Vertical wind shear modeled above  
stack top? (MSHEAR)                   Default: 0       ! MSHEAR = 0   !  
  0 = no (i.e., vertical wind shear not modeled)  
  1 = yes (i.e., vertical wind shear modeled)

Puff splitting allowed? (MSPLIT)      Default: 0       ! MSPLIT = 0   !  
  0 = no (i.e., puffs not split)  
  1 = yes (i.e., puffs are split)

Chemical mechanism flag (MCHEM)       Default: 1       ! MCHEM = 1   !  
  0 = chemical transformation not  
      modeled  
  1 = transformation rates computed  
      internally (MESOPUFF II scheme)  
  2 = user-specified transformation  
      rates used  
  3 = transformation rates computed

internally (RIVAD/ARM3 scheme)  
4 = secondary organic aerosol formation  
   computed (MESOPUFF II scheme for OH)

Aqueous phase transformation flag (MAQCHEM)  
(Used only if MCHEM = 1, or 3)      Default: 0      ! MAQCHEM = 0   !  
0 = aqueous phase transformation  
   not modeled  
1 = transformation rates adjusted  
   for aqueous phase reactions

Wet removal modeled ? (MWET)            Default: 1      ! MWET = 1   !  
0 = no  
1 = yes

Dry deposition modeled ? (MDRY)         Default: 1      ! MDRY = 1   !  
0 = no  
1 = yes  
(dry deposition method specified  
  for each species in Input Group 3)

Method used to compute dispersion  
coefficients (MDISP)                    Default: 3      ! MDISP = 3   !  
  
1 = dispersion coefficients computed from measured values  
   of turbulence, sigma v, sigma w  
2 = dispersion coefficients from internally calculated  
   sigma v, sigma w using micrometeorological variables  
   (u\*, w\*, L, etc.)  
3 = PG dispersion coefficients for RURAL areas (computed using  
   the ISCST multi-segment approximation) and MP coefficients in  
   urban areas  
4 = same as 3 except PG coefficients computed using  
   the MESOPUFF II eqns.  
5 = CTDM sigmas used for stable and neutral conditions.  
   For unstable conditions, sigmas are computed as in  
   MDISP = 3, described above. MDISP = 5 assumes that  
   measured values are read

Sigma-v/sigma-theta, sigma-w measurements used? (MTURBVW)  
(Used only if MDISP = 1 or 5)         Default: 3      ! MTURBVW = 3   !  
1 = use sigma-v or sigma-theta measurements  
   from PROFILE.DAT to compute sigma-y  
   (valid for METFM = 1, 2, 3, 4)  
2 = use sigma-w measurements  
   from PROFILE.DAT to compute sigma-z  
   (valid for METFM = 1, 2, 3, 4)  
3 = use both sigma-(v/theta) and sigma-w  
   from PROFILE.DAT to compute sigma-y and sigma-z  
   (valid for METFM = 1, 2, 3, 4)  
4 = use sigma-theta measurements  
   from PLMMET.DAT to compute sigma-y  
   (valid only if METFM = 3)

Back-up method used to compute dispersion  
when measured turbulence data are  
missing (MDISP2)                        Default: 3      ! MDISP2 = 3   !  
(used only if MDISP = 1 or 5)  
2 = dispersion coefficients from internally calculated  
   sigma v, sigma w using micrometeorological variables  
   (u\*, w\*, L, etc.)  
3 = PG dispersion coefficients for RURAL areas (computed using  
   the ISCST multi-segment approximation) and MP coefficients in  
   urban areas  
4 = same as 3 except PG coefficients computed using

---

the MESOPUFF II eqns.

PG sigma-y,z adj. for roughness?      Default: 0      ! MROUGH = 0 !  
(MROUGH)  
0 = no  
1 = yes

Partial plume penetration of      Default: 1      ! MPARTL = 1 !  
elevated inversion?  
(MPARTL)  
0 = no  
1 = yes

Strength of temperature inversion      Default: 0      ! MTINV = 0 !  
provided in PROFILE.DAT extended records?  
(MTINV)  
0 = no (computed from measured/default gradients)  
1 = yes

PDF used for dispersion under convective conditions?  
Default: 0      ! MPDF = 0 !  
(MPDF)  
0 = no  
1 = yes

Sub-Grid TIBL module used for shore line?  
Default: 0      ! MSGTIBL = 0 !  
(MSGTIBL)  
0 = no  
1 = yes

Boundary conditions (concentration) modeled?  
Default: 0      ! MBCON = 0 !  
(MBCON)  
0 = no  
1 = yes, using formatted BCON.DAT file  
2 = yes, using unformatted CONC.DAT file

Analyses of fogging and icing impacts due to emissions from  
arrays of mechanically-forced cooling towers can be performed  
using CALPUFF in conjunction with a cooling tower emissions  
processor (CTEMISS) and its associated postprocessors. Hourly  
emissions of water vapor and temperature from each cooling tower  
cell are computed for the current cell configuration and ambient  
conditions by CTEMISS. CALPUFF models the dispersion of these  
emissions and provides cloud information in a specialized format  
for further analysis. Output to FOG.DAT is provided in either  
'plume mode' or 'receptor mode' format.

Configure for FOG Model output?      Default: 0      ! MFOG = 0 !  
(MFOG)  
0 = no  
1 = yes - report results in PLUME Mode format  
2 = yes - report results in RECEPTOR Mode format

TEST options specified to see if  
they conform to regulatory  
values? (MREG)      Default: 1      ! MREG = 0 !  
0 = NO checks are made  
1 = Technical options must conform to USEPA  
Long Range Transport (LRT) guidance

---

```

METFM      1 or 2
AVET       60. (min)
PGTIME     60. (min)
MGAUSS     1
MCTADJ     3
MTRANS     1
MTIP       1
MCHEM      1 or 3 (if modeling SOx, NOx)
MWET       1
MDRY       1
MDISP      2 or 3
MPDF       0 if MDISP=3
           1 if MDISP=2
MROUGH     0
MPARTL     1
SYTDEP     550. (m)
MHFTSZ     0

```

!END!

-----  
INPUT GROUP: 3a, 3b -- Species list  
-----

-----  
Subgroup (3a)  
-----

The following species are modeled:

The following species are modeled:

```

! CSPEC =      SO2 !      !END!
! CSPEC =      SO4 !      !END!
! CSPEC =      NOX !      !END!
! CSPEC =      HNO3 !     !END!
! CSPEC =      NO3 !      !END!
! CSPEC =      PMC !      !END!
! CSPEC =      PMF !      !END!
! CSPEC =      EC !       !END!
! CSPEC =      SOA !      !END!

```

SPECIES NAME (Limit: 12 Characters in length)	MODELED (0=NO, 1=YES)	EMITTED (0=NO, 1=YES)	Dry DEPOSITED (0=NO, 1=COMPUTED-GAS 2=COMPUTED-PARTICLE 3=USER-SPECIFIED)	OUTPUT GROUP NUMBER (0=NONE, 1=1st CGRUP, 2=2nd CGRUP, 3= etc.)
! SO2 =	1,	1,	1,	0 !
! SO4 =	1,	1,	2,	0 !
! NOX =	1,	1,	1,	0 !
! HNO3 =	1,	0,	1,	0 !
! NO3 =	1,	0,	2,	0 !
! PMC =	1,	1,	2,	0 !
! PMF =	1,	1,	2,	0 !
! EC =	1,	1,	2,	0 !
! SOA =	1,	1,	2,	0 !

!END!



(XLAT2)                      No Default              ! XLAT2 = 45N !

LCC : Projection cone slices through Earth's surface at XLAT1 and XLAT2  
PS : Projection plane slices through Earth at XLAT1  
(XLAT2 is not used)

-----  
Note: Latitudes and longitudes should be positive, and include a  
letter N,S,E, or W indicating north or south latitude, and  
east or west longitude. For example,  
35.9 N Latitude = 35.9N  
118.7 E Longitude = 118.7E

Datum-region

-----  
The Datum-Region for the coordinates is identified by a character  
string. Many mapping products currently available use the model of the  
Earth known as the World Geodetic System 1984 (WGS-84). Other local  
models may be in use, and their selection in CALMET will make its output  
consistent with local mapping products. The list of Datum-Regions with  
official transformation parameters is provided by the National Imagery and  
Mapping Agency (NIMA).

NIMA Datum - Regions(Examples)

-----  
WGS-84    WGS-84 Reference Ellipsoid and Geoid, Global coverage (WGS84)  
NAS-C    NORTH AMERICAN 1927 Clarke 1866 Spheroid, MEAN FOR CONUS (NAD27)  
NAR-C    NORTH AMERICAN 1983 GRS 80 Spheroid, MEAN FOR CONUS (NAD83)  
NWS-84    NWS 6370KM Radius, Sphere  
ESR-S    ESRI REFERENCE 6371KM Radius, Sphere

Datum-region for output coordinates

(DATUM)                      Default: WGS-84              ! DATUM = NAS-C !

METEOROLOGICAL Grid:

Rectangular grid defined for projection PMAP,  
with X the Easting and Y the Northing coordinate

No. X grid cells (NX)              No default              ! NX = 162 !  
No. Y grid cells (NY)              No default              ! NY = 191 !  
No. vertical layers (NZ)            No default              ! NZ = 10 !  
  
Grid spacing (DGRIDKM)            No default              ! DGRIDKM = 4. !  
Units: km

Cell face heights in arbitrary

vertical grid (ZFACE(NZ+1))        No defaults  
Units: m  
! ZFACE = 0.,20.,40.,80.,160.,320.,640.,1200.,2000.,3000.,4000. !

Reference Coordinates  
of SOUTHWEST corner of  
grid cell(1, 1):

X coordinate (XORIGKM)            No default              ! XORIGKM = 247. !  
Y coordinate (YORIGKM)            No default              ! YORIGKM = 240. !  
Units: km

COMPUTATIONAL Grid:

---



The computational grid is identical to or a subset of the MET. grid.  
 The lower left (LL) corner of the computational grid is at grid point  
 (IBCOMP, JBCOMP) of the MET. grid. The upper right (UR) corner of the  
 computational grid is at grid point (IECOMP, JECOMP) of the MET. grid.  
 The grid spacing of the computational grid is the same as the MET. grid.

X index of LL corner (IBCOMP) (1 <= IBCOMP <= NX)	No default	! IBCOMP = 1 !
Y index of LL corner (JBCOMP) (1 <= JBCOMP <= NY)	No default	! JBCOMP = 1 !
X index of UR corner (IECOMP) (1 <= IECOMP <= NX)	No default	! IECOMP = 162 !
Y index of UR corner (JECOMP) (1 <= JECOMP <= NY)	No default	! JECOMP = 191 !

SAMPLING Grid (GRIDDED RECEPTORS):

The lower left (LL) corner of the sampling grid is at grid point  
 (IBSAMP, JBSAMP) of the MET. grid. The upper right (UR) corner of the  
 sampling grid is at grid point (IESAMP, JESAMP) of the MET. grid.  
 The sampling grid must be identical to or a subset of the computational  
 grid. It may be a nested grid inside the computational grid.  
 The grid spacing of the sampling grid is DGRIDKM/MESHNDN.

Logical flag indicating if gridded receptors are used (LSAMP) (T=yes, F=no)	Default: T	! LSAMP = F !
X index of LL corner (IBSAMP) (IBCOMP <= IBSAMP <= IECOMP)	No default	! IBSAMP = 1 !
Y index of LL corner (JBSAMP) (JBCOMP <= JBSAMP <= JECOMP)	No default	! JBSAMP = 1 !
X index of UR corner (IESAMP) (IBCOMP <= IESAMP <= IECOMP)	No default	! IESAMP = 162 !
Y index of UR corner (JESAMP) (JBCOMP <= JESAMP <= JECOMP)	No default	! JESAMP = 191 !
Nesting factor of the sampling grid (MESHNDN) (MESHNDN is an integer >= 1)	Default: 1	! MESHNDN = 1 !

!END!

-----  
 INPUT GROUP: 5 -- Output Options  
 -----

FILE	* DEFAULT VALUE	* VALUE THIS RUN
----	-----	-----

**Refined BART Modeling Protocol for  
Wausau Mosinee Mill**

---

```
Concentrations (ICON)           1           ! ICON = 1  !
Dry Fluxes (IDRY)               1           ! IDRY = 0  !
Wet Fluxes (IWET)               1           ! IWET = 0  !
Relative Humidity (IVIS)        1           ! IVIS = 0  !
  (relative humidity file is
   required for visibility
   analysis)
Use data compression option in output file?
(LCOMPRES)                       Default: T           ! LCOMPRES = T !
```

\*

0 = Do not create file, 1 = create file

DIAGNOSTIC MASS FLUX OUTPUT OPTIONS:

```
Mass flux across specified boundaries
for selected species reported hourly?
(IMFLX)                           Default: 0           ! IMFLX = 0  !
  0 = no
  1 = yes (FLUXBDY.DAT and MASSFLX.DAT filenames
           are specified in Input Group 0)
```

```
Mass balance for each species
reported hourly?
(IMBAL)                           Default: 0           ! IMBAL = 0  !
  0 = no
  1 = yes (MASSBAL.DAT filename is
           specified in Input Group 0)
```

LINE PRINTER OUTPUT OPTIONS:

```
Print concentrations (ICPRT)      Default: 0           ! ICPRT = 0  !
Print dry fluxes (IDPRT)         Default: 0           ! IDPRT = 0  !
Print wet fluxes (IWPRT)        Default: 0           ! IWPRT = 0  !
(0 = Do not print, 1 = Print)
```

```
Concentration print interval
(ICFRQ) in hours                 Default: 1           ! ICFRQ = 1  !
Dry flux print interval
(IDFRQ) in hours                 Default: 1           ! IDFRQ = 1  !
Wet flux print interval
(IWFRQ) in hours                 Default: 1           ! IWFRQ = 1  !
```

```
Units for Line Printer Output
(IPRTU)                          Default: 1           ! IPRTU = 3  !
      for                          for
      Concentration                Deposition
  1 =      g/m**3                  g/m**2/s
  2 =      mg/m**3                 mg/m**2/s
  3 =      ug/m**3                 ug/m**2/s
  4 =      ng/m**3                 ng/m**2/s
  5 =      Odour Units
```

```
Messages tracking progress of run
written to the screen ?
(IMESG)                          Default: 2           ! IMESG = 2  !
  0 = no
  1 = yes (advection step, puff ID)
  2 = yes (YYYYJJJHH, # old puffs, # emitted puffs)
```

SPECIES (or GROUP for combined species) LIST FOR OUTPUT OPTIONS

**Refined BART Modeling Protocol for  
Wausau Mosinee Mill**

```

----- CONCENTRATIONS ----- ----- DRY FLUXES ----- ----- WET FLUXES -----
-- MASS FLUX --
SPECIES
/GROUP      PRINTED?  SAVED ON DISK?  PRINTED?  SAVED ON DISK?  PRINTED?  SAVED ON DISK?
SAVED ON DISK?
-----
!          SO2 =    0,      1,      0,      0,      0,      0,
0 !
!          SO4 =    0,      1,      0,      0,      0,      0,
0 !
!          NOX =    0,      1,      0,      0,      0,      0,
0 !
!          HNO3 =   0,      1,      0,      0,      0,      0,
0 !
!          NO3 =    0,      1,      0,      0,      0,      0,
0 !
!          PMC =    0,      1,      0,      0,      0,      0,
0 !
!          PMF =    0,      1,      0,      0,      0,      0,
0 !
!          EC =     0,      1,      0,      0,      0,      0,
0 !
!          SOA =    0,      1,      0,      0,      0,      0,
0 !

```

OPTIONS FOR PRINTING "DEBUG" QUANTITIES (much output)

```

Logical for debug output
(LDEBUG)                                Default: F      ! LDEBUG = F !

First puff to track
(IPFDEB)                                Default: 1      ! IPFDEB = 1 !

Number of puffs to track
(NPFDEB)                                Default: 1      ! NPFDEB = 1 !

Met. period to start output
(NN1)                                    Default: 1      ! NN1 = 1 !

Met. period to end output
(NN2)                                    Default: 10     ! NN2 = 10 !

!END!

```

INPUT GROUP: 6a, 6b, & 6c -- Subgrid scale complex terrain inputs

-----  
Subgroup (6a)  
-----

```

Number of terrain features (NHILL)      Default: 0      ! NHILL = 0 !

Number of special complex terrain
receptors (NCTREC)                      Default: 0      ! NCTREC = 0 !

Terrain and CTSG Receptor data for
CTSG hills input in CTDM format ?
(MHILL)                                  No Default     ! MHILL = 2 !
1 = Hill and Receptor data created

```

by CTDM processors & read from  
HILL.DAT and HILLRCT.DAT files  
2 = Hill data created by OPTHILL &  
input below in Subgroup (6b);  
Receptor data in Subgroup (6c)

Factor to convert horizontal dimensions Default: 1.0 ! XHILL2M = 1. !  
to meters (MHILL=1)

Factor to convert vertical dimensions Default: 1.0 ! ZHILL2M = 1. !  
to meters (MHILL=1)

X-origin of CTDM system relative to No Default ! XCTDMKM = 0.0E00 !  
CALPUFF coordinate system, in Kilometers (MHILL=1)

Y-origin of CTDM system relative to No Default ! YCTDMKM = 0.0E00 !  
CALPUFF coordinate system, in Kilometers (MHILL=1)

! END !

-----  
Subgroup (6b)  
-----

1 \*\*  
HILL information

HILL AMAX1 NO. (m)	XC AMAX2 (km) (m)	YC (km)	THETAH (deg.)	ZGRID (m)	RELIEF (m)	EXPO 1 (m)	EXPO 2 (m)	SCALE 1 (m)	SCALE 2 (m)
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----
----	----								

-----  
Subgroup (6c)  
-----

COMPLEX TERRAIN RECEPTOR INFORMATION

XRCT (km)	YRCT (km)	ZRCT (m)	XHH
-----	-----	-----	----

-----  
1

Description of Complex Terrain Variables:

XC, YC = Coordinates of center of hill  
THETAH = Orientation of major axis of hill (clockwise from  
North)  
ZGRID = Height of the 0 of the grid above mean sea  
level  
RELIEF = Height of the crest of the hill above the grid elevation  
EXPO 1 = Hill-shape exponent for the major axis  
EXPO 2 = Hill-shape exponent for the major axis  
SCALE 1 = Horizontal length scale along the major axis  
SCALE 2 = Horizontal length scale along the minor axis  
AMAX = Maximum allowed axis length for the major axis  
BMAX = Maximum allowed axis length for the major axis

XRCT, YRCT = Coordinates of the complex terrain receptors  
ZRCT = Height of the ground (MSL) at the complex terrain  
Receptor

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---

XHH = Hill number associated with each complex terrain receptor  
(NOTE: MUST BE ENTERED AS A REAL NUMBER)

\*\*

NOTE: DATA for each hill and CTSG receptor are treated as a separate  
input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUP: 7 -- Chemical parameters for dry deposition of gases  
-----

SPECIES LAW COEFFICIENT NAME (dimensionless)	DIFFUSIVITY (cm**2/s)	ALPHA STAR	REACTIVITY	MESOPHYLL RESISTANCE (s/cm)	HENRY'S
! SO2 =	0.1509,	1000.,	8.,	0.,	
0.04 !					
! NOX =	0.1656,	1.,	8.,	5.,	
3.5 !					
! HNO3 =	0.1628,	1.,	18.,	0.,	
0.00000008 !					
!END!					

-----  
INPUT GROUP: 8 -- Size parameters for dry deposition of particles  
-----

For SINGLE SPECIES, the mean and standard deviation are used to  
compute a deposition velocity for NINT (see group 9) size-ranges,  
and these are then averaged to obtain a mean deposition velocity.

For GROUPED SPECIES, the size distribution should be explicitly  
specified (by the 'species' in the group), and the standard deviation  
for each should be entered as 0. The model will then use the  
deposition velocity for the stated mean diameter.

SPECIES NAME	GEOMETRIC MASS MEAN DIAMETER (microns)	GEOMETRIC STANDARD DEVIATION (microns)
! SO4 =	0.48,	2. !
! NO3 =	0.48,	2. !
! PMC =	2.0,	2.0 !
! PMF =	0.48,	2. !
! EC =	0.48,	2. !
! SOA =	0.48,	2. !

!END!

-----  
INPUT GROUP: 9 -- Miscellaneous dry deposition parameters  
-----

```

Reference cuticle resistance (s/cm)
(RCUTR)                      Default: 30    ! RCUTR = 30.0 !
Reference ground resistance (s/cm)
(RGR)                        Default: 10    ! RGR = 10.0 !
Reference pollutant reactivity
(REACTR)                     Default: 8     ! REACTR = 8.0 !

Number of particle-size intervals used to
evaluate effective particle deposition velocity
(NINT)                        Default: 9     ! NINT = 9 !

Vegetation state in unirrigated areas
(IVEG)                        Default: 1     ! IVEG = 1 !
  IVEG=1 for active and unstressed vegetation
  IVEG=2 for active and stressed vegetation
  IVEG=3 for inactive vegetation

```

!END!

-----

INPUT GROUP: 10 -- Wet Deposition Parameters

-----

Scavenging Coefficient -- Units: (sec)\*\*(-1)

Pollutant	Liquid Precip.	Frozen Precip.
! SO2 =	3.0E-05,	0.0E00 !
! SO4 =	1.0E-04,	3.0E-05 !
! NOX =	0.0E00,	0.0E00 !
! HNO3 =	6.0E-05,	0.0E00 !
! NO3 =	1.0E-04,	3.0E-05 !

!END!

-----

INPUT GROUP: 11 -- Chemistry Parameters

-----

```

Ozone data input option (MOZ)   Default: 1           ! MOZ = 0 !
(Used only if MCHEM = 1, 3, or 4)
  0 = use a monthly background ozone value
  1 = read hourly ozone concentrations from
    the OZONE.DAT data file

```

```

Monthly ozone concentrations
(Used only if MCHEM = 1, 3, or 4 and
MOZ = 0 or MOZ = 1 and all hourly O3 data missing)
(BCKO3) in ppb                    Default: 12*80.

```

```

! BCKO3 = 31.00, 31.00, 31.00, 37.00, 37.00, 37.00, 33.00, 33.00, 33.00, 27.00, 27.00, 27.00
!

```

```

Monthly ammonia concentrations
(Used only if MCHEM = 1, or 3)
(BCKNH3) in ppb                   Default: 12*10.

```

! BCKNH3 = 12\*3.9 !

Nighttime SO2 loss rate (RNITE1)  
in percent/hour                   Default: 0.2                   ! RNITE1 = .2 !

Nighttime NOx loss rate (RNITE2)  
in percent/hour                   Default: 2.0                   ! RNITE2 = 2.0 !

Nighttime HNO3 formation rate (RNITE3)  
in percent/hour                   Default: 2.0                   ! RNITE3 = 2.0 !

H2O2 data input option (MH2O2)   Default: 1                   ! MH2O2 = 1 !  
(Used only if MAQCHEM = 1)  
  0 = use a monthly background H2O2 value  
  1 = read hourly H2O2 concentrations from  
      the H2O2.DAT data file

Monthly H2O2 concentrations  
(Used only if MQACHEM = 1 and  
  MH2O2 = 0 or MH2O2 = 1 and all hourly H2O2 data missing)  
(BCKH2O2) in ppb                   Default: 12\*1.  
\* BCKH2O2 = 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00 \*  
! BCKH2O2 = 0.50, 0.50, 0.50, 2.30, 2.30, 2.30, 3.50, 3.50, 3.50, 0.80, 0.80, 0.80 !

--- Data for SECONDARY ORGANIC AEROSOL (SOA) Option  
(used only if MCHEM = 4)

The SOA module uses monthly values of:  
  Fine particulate concentration in ug/m^3 (BCKPMF)  
  Organic fraction of fine particulate       (OFRAC)  
  VOC / NOX ratio (after reaction)       (VCNX)  
to characterize the air mass when computing  
the formation of SOA from VOC emissions.  
Typical values for several distinct air mass types are:

Month	1	2	3	4	5	6	7	8	9	10	11	12
	Jan	Feb	Mar	Apr	May	Jun	Jul	Aug	Sep	Oct	Nov	Dec
<b>Clean Continental</b>												
BCKPMF	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.	1.
OFRAC	.15	.15	.20	.20	.20	.20	.20	.20	.20	.20	.20	.15
VCNX	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.
<b>Clean Marine (surface)</b>												
BCKPMF	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5	.5
OFRAC	.25	.25	.30	.30	.30	.30	.30	.30	.30	.30	.30	.25
VCNX	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.	50.
<b>Urban - low biogenic (controls present)</b>												
BCKPMF	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.	30.
OFRAC	.20	.20	.25	.25	.25	.25	.25	.25	.20	.20	.20	.20
VCNX	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.	4.
<b>Urban - high biogenic (controls present)</b>												
BCKPMF	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.	60.
OFRAC	.25	.25	.30	.30	.30	.55	.55	.55	.35	.35	.35	.25
VCNX	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.
<b>Regional Plume</b>												
BCKPMF	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.	20.
OFRAC	.20	.20	.25	.35	.25	.40	.40	.40	.30	.30	.30	.20
VCNX	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.	15.
<b>Urban - no controls present</b>												
BCKPMF	100.	100.	100.	100.	100.	100.	100.	100.	100.	100.	100.	100.

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---

```
OFRAC  .30  .30  .35  .35  .35  .55  .55  .55  .35  .35  .35  .30
VCNX   2.   2.   2.   2.   2.   2.   2.   2.   2.   2.   2.   2.
```

Default: Clean Continental

```
! BCKPMF = 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00, 1.00 !
! OFRAC  = 0.15, 0.15, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.20, 0.15 !
! VCNX   = 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00, 50.00,
50.00 !
```

!END!

-----

INPUT GROUP: 12 -- Misc. Dispersion and Computational Parameters

-----

Horizontal size of puff (m) beyond which  
time-dependent dispersion equations (Heffter)  
are used to determine sigma-y and  
sigma-z (SYTDEP) Default: 550. ! SYTDEP = 5.5E02 !

Switch for using Heffter equation for sigma z  
as above (0 = Not use Heffter; 1 = use Heffter  
(MHFTSZ) Default: 0 ! MHFTSZ = 0 !

Stability class used to determine plume  
growth rates for puffs above the boundary  
layer (JSUP) Default: 5 ! JSUP = 5 !

Vertical dispersion constant for stable  
conditions (k1 in Eqn. 2.7-3) (CONK1) Default: 0.01 ! CONK1 = .01 !

Vertical dispersion constant for neutral/  
unstable conditions (k2 in Eqn. 2.7-4)  
(CONK2) Default: 0.1 ! CONK2 = .1 !

Factor for determining Transition-point from  
Schulman-Scire to Huber-Snyder Building Downwash  
scheme (SS used for Hs < Hb + TBD \* HL)  
(TBD) Default: 0.5 ! TBD = .5 !  
TBD < 0 ==> always use Huber-Snyder  
TBD = 1.5 ==> always use Schulman-Scire  
TBD = 0.5 ==> ISC Transition-point

Range of land use categories for which  
urban dispersion is assumed  
(IURB1, IURB2) Default: 10 ! IURB1 = 10 !  
19 ! IURB2 = 19 !

Site characterization parameters for single-point Met data files -----  
(needed for METFM = 2,3,4)

Land use category for modeling domain  
(ILANDUIN) Default: 20 ! ILANDUIN = 20 !

Roughness length (m) for modeling domain  
(Z0IN) Default: 0.25 ! Z0IN = .25 !

Leaf area index for modeling domain  
(XLAIIN) Default: 3.0 ! XLAIIN = 3.0 !

Elevation above sea level (m)

---



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---

```

(ELEVIN)                                Default: 0.0    ! ELEVIN = .0 !

Latitude (degrees) for met location
(XLATIN)                                Default: -999. ! XLATIN = -999.0 !

Longitude (degrees) for met location
(XLONIN)                                Default: -999. ! XLONIN = -999.0 !

Specialized information for interpreting single-point Met data files -----

Anemometer height (m) (Used only if METFM = 2,3)
(ANEMHT)                                Default: 10.   ! ANEMHT = 10.0 !

Form of lateral turbulence data in PROFILE.DAT file
(Used only if METFM = 4 or MTURBVW = 1 or 3)
(ISIGMAV)                               Default: 1     ! ISIGMAV = 1 !
    0 = read sigma-theta
    1 = read sigma-v

Choice of mixing heights (Used only if METFM = 4)
(IMIXCTDM)                              Default: 0     ! IMIXCTDM = 0 !
    0 = read PREDICTED mixing heights
    1 = read OBSERVED mixing heights

Maximum length of a slug (met. grid units)
(XMXLEN)                                Default: 1.0   ! XMXLEN = 1.0 !

Maximum travel distance of a puff/slug (in
grid units) during one sampling step
(XSAMPLN)                               Default: 1.0   ! XSAMPLN = 1.0 !

Maximum Number of slugs/puffs release from
one source during one time step
(MXNEW)                                  Default: 99    ! MXNEW = 99 !

Maximum Number of sampling steps for
one puff/slug during one time step
(MXSAM)                                  Default: 99    ! MXSAM = 99 !

Number of iterations used when computing
the transport wind for a sampling step
that includes gradual rise (for CALMET
and PROFILE winds)
(NCOUNT)                               Default: 2     ! NCOUNT = 2 !

Minimum sigma y for a new puff/slug (m)
(SYMIN)                                  Default: 1.0   ! SYMIN = 1.0 !

Minimum sigma z for a new puff/slug (m)
(SZMIN)                                  Default: 1.0   ! SZMIN = 1.0 !

Default minimum turbulence velocities
sigma-v and sigma-w for each
stability class (m/s)
(SVMIN(6) and SWMIN(6))                 Default SVMIN : .50, .50, .50, .50, .50, .50
                                           Default SWMIN : .20, .12, .08, .06, .03, .016

                                           Stability Class :  A    B    C    D    E    F
                                           ---  ---  ---  ---  ---  ---
                                           ! SVMIN = 0.500, 0.500, 0.500, 0.500, 0.500, 0.500!
                                           ! SWMIN = 0.200, 0.120, 0.080, 0.060, 0.030, 0.016!

Divergence criterion for dw/dz across puff
used to initiate adjustment for horizontal
convergence (1/s)

```

```

Partial adjustment starts at CDIV(1), and
full adjustment is reached at CDIV(2)
(CDIV(2))                                Default: 0.0,0.0 ! CDIV = .0, .0 !

Minimum wind speed (m/s) allowed for
non-calm conditions. Also used as minimum
speed returned when using power-law
extrapolation toward surface
(WSCALM)                                Default: 0.5 ! WSCALM = .5 !

Maximum mixing height (m)
(XMAXZI)                                Default: 3000. ! XMAXZI = 3000.0 !

Minimum mixing height (m)
(XMINZI)                                Default: 50. ! XMINZI = 50.0 !

Default wind speed classes --
5 upper bounds (m/s) are entered;
the 6th class has no upper limit
(WSCAT(5))                                Default :
ISC RURAL : 1.54, 3.09, 5.14, 8.23, 10.8 (10.8+)

Wind Speed Class : 1      2      3      4      5
                   ---    ---    ---    ---    ---
                   ! WSCAT = 1.54, 3.09, 5.14, 8.23, 10.80 !

Default wind speed profile power-law
exponents for stabilities 1-6
(PLX0(6))                                Default : ISC RURAL values
ISC RURAL : .07, .07, .10, .15, .35, .55
ISC URBAN : .15, .15, .20, .25, .30, .30

Stability Class :  A      B      C      D      E      F
                   ---    ---    ---    ---    ---    ---
                   ! PLX0 = 0.07, 0.07, 0.10, 0.15, 0.35, 0.55 !

Default potential temperature gradient
for stable classes E, F (degK/m)
(PTG0(2))                                Default: 0.020, 0.035
                   ! PTG0 = 0.020, 0.035 !

Default plume path coefficients for
each stability class (used when option
for partial plume height terrain adjustment
is selected -- MCTADJ=3)
(PPC(6))                                Stability Class :  A      B      C      D      E      F
Default PPC : .50, .50, .50, .50, .35, .35
                   ---    ---    ---    ---    ---    ---
                   ! PPC = 0.50, 0.50, 0.50, 0.50, 0.35, 0.35 !

Slug-to-puff transition criterion factor
equal to sigma-y/length of slug
(SL2PF)                                Default: 10. ! SL2PF = 10.0 !

Puff-splitting control variables -----

VERTICAL SPLIT
-----

Number of puffs that result every time a puff
is split - nsplit=2 means that 1 puff splits
into 2
(NSPLIT)                                Default: 3 ! NSPLIT = 3 !

Time(s) of a day when split puffs are eligible to

```

be split once again; this is typically set once  
per day, around sunset before nocturnal shear develops.  
24 values: 0 is midnight (00:00) and 23 is 11 PM (23:00)  
0=do not re-split 1=eligible for re-split  
(IRESPLIT(24)) Default: Hour 17 = 1  
! IRESPLIT = 0,0 !

Split is allowed only if last hour's mixing  
height (m) exceeds a minimum value  
(ZISPLIT) Default: 100. ! ZISPLIT = 100.0 !

Split is allowed only if ratio of last hour's  
mixing ht to the maximum mixing ht experienced  
by the puff is less than a maximum value (this  
postpones a split until a nocturnal layer develops)  
(ROLDMAX) Default: 0.25 ! ROLDMAX = 0.25 !

HORIZONTAL SPLIT  
-----

Number of puffs that result every time a puff  
is split - nsplith=5 means that 1 puff splits  
into 5  
(NSPLITH) Default: 5 ! NSPLITH = 5 !

Minimum sigma-y (Grid Cells Units) of puff  
before it may be split  
(SYSPLITH) Default: 1.0 ! SYSPLITH = 1.0 !

Minimum puff elongation rate (SYSPLITH/hr) due to  
wind shear, before it may be split  
(SHSPLITH) Default: 2. ! SHSPLITH = 2.0 !

Minimum concentration (g/m<sup>3</sup>) of each  
species in puff before it may be split  
Enter array of NSPEC values; if a single value is  
entered, it will be used for ALL species  
(CNSPLITH) Default: 1.0E-07 ! CNSPLITH = 1.0E-07 !

Integration control variables -----

Fractional convergence criterion for numerical SLUG  
sampling integration  
(EPSSLUG) Default: 1.0e-04 ! EPSSLUG = 1.0E-04 !

Fractional convergence criterion for numerical AREA  
source integration  
(EPSAREA) Default: 1.0e-06 ! EPSAREA = 1.0E-06 !

Trajectory step-length (m) used for numerical rise  
integration  
(DSRISE) Default: 1.0 ! DSRISE = 1.0 !

Boundary Condition (BC) Puff control variables -----

Minimum height (m) to which BC puffs are mixed as they are emitted  
(MBCON=2 ONLY). Actual height is reset to the current mixing height  
at the release point if greater than this minimum.  
(HTMINBC) Default: 500. ! HTMINBC = 500.0 !

Search radius (in BC segment lengths) about a receptor for sampling  
nearest BC puff. BC puffs are emitted with a spacing of one segment  
length, so the search radius should be greater than 1.  
(RSAMPBC) Default: 4. ! RSAMPBC = 10.0 !

Near-Surface depletion adjustment to concentration profile used when  
sampling BC puffs?  
(MDEPBC) Default: 1 ! MDEPBC = 1 !  
0 = Concentration is NOT adjusted for depletion  
1 = Adjust Concentration for depletion

!END!

-----  
INPUT GROUPS: 13a, 13b, 13c, 13d -- Point source parameters  
-----

-----  
Subgroup (13a)  
-----

Number of point sources with  
parameters provided below (NPT1) No default ! NPT1 = 5 !

Units used for point source  
emissions below (IPTU) Default: 1 ! IPTU = 1 !

1 = g/s  
2 = kg/hr  
3 = lb/hr  
4 = tons/yr  
5 = Odour Unit \* m\*\*3/s (vol. flux of odour compound)  
6 = Odour Unit \* m\*\*3/min  
7 = metric tons/yr

Number of source-species  
combinations with variable  
emissions scaling factors  
provided below in (13d) (NSPT1) Default: 0 ! NSPT1 = 0 !

Number of point sources with  
variable emission parameters  
provided in external file (NPT2) No default ! NPT2 = 0 !

(If NPT2 > 0, these point  
source emissions are read from  
the file: PTEMARB.DAT)

!END!

-----  
Subgroup (13b)  
-----

The following species are modeled:

CSPEC = SO2  
CSPEC = SO4  
CSPEC = NOX  
CSPEC = HNO3  
CSPEC = NO3  
CSPEC = PMC  
CSPEC = PMF  
CSPEC = EC  
CSPEC = SOA

a

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POINT SOURCE: CONSTANT DATA

```

-----
Source      X      Y      Stack   Base   Stack   Exit   Exit   Bldg.  Emission
No.        Coordinate Coordinate Height Elevation Diameter Vel.  Temp.  Dwash  Rates
          (km)      (km)      (m)      (m)      (m)      (m/s) (deg. K)          SO2 SO4 NOX HNO3
NO3 PMC PMF EC SOA
-----
1 ! SRCNAM = S10 !
1 ! X = 577.39257, 553.25636, 64.9, 352.7, 2.6 , 14.6, 449, .0, 143.1, 6.43, 28.1, 0, 0, 1.16,
2.39, 0.09, 1.61 !
1 ! FMFAC =      1.0 ! !END!
2 ! SRCNAM = S11 !
2 ! X = 577.37101, 553.15432, 36 , 352.7, 1.64, 25.7, 457, .0, 13.12, 0.19, 5.35, 0, 0, 0.73,
1.26, 0.10, 0.04 !
2 ! FMFAC =      1.0 ! !END!
3 ! SRCNAM = S12 !
3 ! X = 577.37292, 553.18169, 41.8, 352.7, 0.8 , 11.4, 347, .0, 0.40 , 0 , 0 , 0, 0, 0.00,
0.43, 0.00, 0.00 !
3 ! FMFAC =      1.0 ! !END!
4 ! SRCNAM = S52 !
4 ! X = 577.32104, 553.10738, 23.4, 352.7, 0.51, 9.4 , 313, .0, 0 0 , 0 , 0 , 0, 0, 0.43,
0.00, 0.00, 0.00 !
4 ! FMFAC =      1.0 ! !END!
5 ! SRCNAM = S53 !
5 ! X = 577.31194, 553.12545, 25.9, 352.7, 0.98, 8.8 , 347, .0, 1.35 , 0.10, 0.92, 0, 0, 0.32,
0.19, 0.01, 0.00 !
5 ! FMFAC =      1.0 ! !END!
-----

```

a

Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.

```

SRCNAM is a 12-character name for a source
       (No default)
X      is an array holding the source data listed by the column headings
       (No default)
SIGYZI is an array holding the initial sigma-y and sigma-z (m)
       (Default: 0.,0.)
FMFAC  is a vertical momentum flux factor (0. or 1.0) used to represent
       the effect of rain-caps or other physical configurations that
       reduce momentum rise associated with the actual exit velocity.
       (Default: 1.0 -- full momentum used)

```

b

0. = No building downwash modeled, 1. = downwash modeled  
NOTE: must be entered as a REAL number (i.e., with decimal point)

c

An emission rate must be entered for every pollutant modeled.  
Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by IPTU (e.g. 1 for g/s).

-----  
Subgroup (13c)  
-----

BUILDING DIMENSION DATA FOR SOURCES SUBJECT TO DOWNWASH

```

-----
Source      a
No.        Effective building height, width, length and X/Y offset (in meters)
-----

```

every 10 degrees. LENGTH, XBADJ, and YBADJ are only needed for  
MBDW=2 (PRIME downwash option)

-----

-----

a  
Building height, width, length, and X/Y offset from the source are treated  
as a separate input subgroup for each source and therefore must end with  
an input group terminator. The X/Y offset is the position, relative to the  
stack, of the center of the upwind face of the projected building, with the  
x-axis pointing along the flow direction.

-----

Subgroup (13d)

-----

a  
POINT SOURCE: VARIABLE EMISSIONS DATA

-----

Use this subgroup to describe temporal variations in the emission  
rates given in 13b. Factors entered multiply the rates in 13b.  
Skip sources here that have constant emissions. For more elaborate  
variation in source parameters, use PTEMARB.DAT and NPT2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

0 =	Constant
1 =	Diurnal cycle (24 scaling factors: hours 1-24)
2 =	Monthly cycle (12 scaling factors: months 1-12)
3 =	Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)
4 =	Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12
5 =	Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----

a  
Data for each species are treated as a separate input subgroup  
and therefore must end with an input group terminator.

-----

INPUT GROUPS: 14a, 14b, 14c, 14d -- Area source parameters

-----

-----

Subgroup (14a)

-----

Number of polygon area sources with  
parameters specified below (NAR1) No default ! NAR1 = 0 !

Units used for area source  
emissions below (IARU) Default: 1 ! IARU = 1 !

---

```

1 =      g/m**2/s
2 =      kg/m**2/hr
3 =      lb/m**2/hr
4 =      tons/m**2/yr
5 =      Odour Unit * m/s (vol. flux/m**2 of odour compound)
6 =      Odour Unit * m/min
7 =      metric tons/m**2/yr

```

Number of source-species combinations with variable emissions scaling factors provided below in (14d) (NSAR1) Default: 0 ! NSAR1 = 0 !

Number of buoyant polygon area sources with variable location and emission parameters (NAR2) No default ! NAR2 = 0 !  
(If NAR2 > 0, ALL parameter data for these sources are read from the file: BAEMARB.DAT)

!END!

-----  
Subgroup (14b)  
-----

```

                                     a
      AREA SOURCE: CONSTANT DATA
      -----
Source          Effect.   Base      Initial   Emission
No.            Height    Elevation Sigma z    Rates
              (m)        (m)        (m)
-----

```

-----  
a  
Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.  
b  
An emission rate must be entered for every pollutant modeled. Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by IARU (e.g. 1 for g/m\*\*2/s).

-----  
Subgroup (14c)  
-----

```

      COORDINATES (km) FOR EACH VERTEX(4) OF EACH POLYGON
      -----
Source          Ordered list of X followed by list of Y, grouped by source
No.
-----

```

-----  
a  
Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
Subgroup (14d)  
-----

a

AREA SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 14b. Factors entered multiply the rates in 14b. Skip sources here that have constant emissions. For more elaborate variation in source parameters, use BAEMARB.DAT and NAR2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

- 0 = Constant
- 1 = Diurnal cycle (24 scaling factors: hours 1-24)
- 2 = Monthly cycle (12 scaling factors: months 1-12)
- 3 = Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)
- 4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)
- 5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

-----  
a  
Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 15a, 15b, 15c -- Line source parameters  
-----

-----  
Subgroup (15a)  
-----

Number of buoyant line sources with variable location and emission parameters (NLN2) No default ! NLN2 = 0 !

(If NLN2 > 0, ALL parameter data for these sources are read from the file: LNEMARB.DAT)

Number of buoyant line sources (NLINES) No default ! NLINES = 0 !

Units used for line source emissions below (ILNU) Default: 1 ! ILNU = 1 !

- 1 = g/s
- 2 = kg/hr
- 3 = lb/hr
- 4 = tons/yr
- 5 = Odour Unit \* m\*\*3/s (vol. flux of odour compound)
- 6 = Odour Unit \* m\*\*3/min
- 7 = metric tons/yr

Number of source-species combinations with variable emissions scaling factors provided below in (15c) (NSLN1) Default: 0 ! NSLN1 = 0 !



Maximum number of segments used to model  
each line (MXNSEG) Default: 7 ! MXNSEG = 7 !

The following variables are required only if NLINES > 0. They are  
used in the buoyant line source plume rise calculations.

Number of distances at which  
transitional rise is computed Default: 6 ! NLRISE = 6 !

Average building length (XL) No default ! XL = .0 !  
(in meters)

Average building height (HBL) No default ! HBL = .0 !  
(in meters)

Average building width (WBL) No default ! WBL = .0 !  
(in meters)

Average line source width (WML) No default ! WML = .0 !  
(in meters)

Average separation between buildings (DXL) No default ! DXL = .0 !  
(in meters)

Average buoyancy parameter (FPRIMEL) No default ! FPRIMEL = .0 !  
(in m\*\*4/s\*\*3)

!END!

-----  
Subgroup (15b)  
-----

BUOYANT LINE SOURCE: CONSTANT DATA  
-----

Source No.	Beg. X Coordinate (km)	Beg. Y Coordinate (km)	End. X Coordinate (km)	End. Y Coordinate (km)	Release Height (m)	Base Elevation (m)	Emission Rates
---------------	------------------------------	------------------------------	------------------------------	------------------------------	--------------------------	--------------------------	-------------------

a  
Data for each source are treated as a separate input subgroup  
and therefore must end with an input group terminator.

b  
An emission rate must be entered for every pollutant modeled.  
Enter emission rate of zero for secondary pollutants that are  
modeled, but not emitted. Units are specified by ILNTU  
(e.g. 1 for g/s).

-----  
Subgroup (15c)  
-----

BUOYANT LINE SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission  
rates given in 15b. Factors entered multiply the rates in 15b.  
Skip sources here that have constant emissions.

IVARY determines the type of variation, and is source-specific:

```
(IVARY)                                Default: 0
 0 = Constant
 1 = Diurnal cycle (24 scaling factors: hours 1-24)
 2 = Monthly cycle (12 scaling factors: months 1-12)
 3 = Hour & Season (4 groups of 24 hourly scaling factors,
    where first group is DEC-JAN-FEB)
 4 = Speed & Stab. (6 groups of 6 scaling factors, where
    first group is Stability Class A,
    and the speed classes have upper
    bounds (m/s) defined in Group 12
 5 = Temperature (12 scaling factors, where temperature
    classes have upper bounds (C) of:
    0, 5, 10, 15, 20, 25, 30, 35, 40,
    45, 50, 50+)
```

-----

a  
Data for each species are treated as a separate input subgroup  
and therefore must end with an input group terminator.

-----

INPUT GROUPS: 16a, 16b, 16c -- Volume source parameters

-----

-----  
Subgroup (16a)  
-----

Number of volume sources with  
parameters provided in 16b,c (NVL1)      No default !    NVL1 = 0 !

Units used for volume source  
emissions below in 16b      (IVLU)      Default: 1 !    IVLU = 1 !

```
1 = g/s
2 = kg/hr
3 = lb/hr
4 = tons/yr
5 = Odour Unit * m**3/s (vol. flux of odour compound)
6 = Odour Unit * m**3/min
7 = metric tons/yr
```

Number of source-species  
combinations with variable  
emissions scaling factors  
provided below in (16c)      (NSVL1)      Default: 0 !    NSVL1 = 0 !

Number of volume sources with  
variable location and emission  
parameters      (NVL2)      No default !    NVL2 = 0 !

(If NVL2 > 0, ALL parameter data for  
these sources are read from the VOLEMARB.DAT file(s) )

!END!

-----  
Subgroup (16b)  
-----

a  
VOLUME SOURCE: CONSTANT DATA

```

-----
      X           Y           Effect.   Base      Initial   Initial   Emission
Coordinate  Coordinate  Height    Elevation  Sigma y    Sigma z    Rates
      (km)       (km)       (m)       (m)       (m)       (m)
-----

```

a  
Data for each source are treated as a separate input subgroup and therefore must end with an input group terminator.

b  
An emission rate must be entered for every pollutant modeled. Enter emission rate of zero for secondary pollutants that are modeled, but not emitted. Units are specified by IVLU (e.g. 1 for g/s).

-----  
Subgroup (16c)  
-----

a  
VOLUME SOURCE: VARIABLE EMISSIONS DATA  
-----

Use this subgroup to describe temporal variations in the emission rates given in 16b. Factors entered multiply the rates in 16b. Skip sources here that have constant emissions. For more elaborate variation in source parameters, use VOLEMARB.DAT and NVL2 > 0.

IVARY determines the type of variation, and is source-specific:  
(IVARY) Default: 0

- 0 = Constant
- 1 = Diurnal cycle (24 scaling factors: hours 1-24)
- 2 = Monthly cycle (12 scaling factors: months 1-12)
- 3 = Hour & Season (4 groups of 24 hourly scaling factors, where first group is DEC-JAN-FEB)
- 4 = Speed & Stab. (6 groups of 6 scaling factors, where first group is Stability Class A, and the speed classes have upper bounds (m/s) defined in Group 12)
- 5 = Temperature (12 scaling factors, where temperature classes have upper bounds (C) of: 0, 5, 10, 15, 20, 25, 30, 35, 40, 45, 50, 50+)

a  
Data for each species are treated as a separate input subgroup and therefore must end with an input group terminator.

-----  
INPUT GROUPS: 17a & 17b -- Non-gridded (discrete) receptor information  
-----

-----  
Subgroup (17a)  
-----

Number of non-gridded receptors (NREC) No default ! NREC = 1995 !

!END!

Note - three Class I areas  
Seney SENE = 173 receptors  
Isle Royal ISLE = 966 receptors  
Boundary Waters BOWA = 856 receptors  
total number of receptors is 1995

**ALL RECEPTOR COORDINATES FOLLOW IN INPUT FILE**