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DEVELOPMENT AND VALIDATION OF STANDARDIZED INFRARED RADIATION MODEL (SIRRM)

GAS/PARTICLE RADIATION TRANSFER MODEL

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FOREWORD

This report was submitted by Photon Research Associates, Inc., La Jolla, CA 92037, under Contract No. F04611-78-C-0081 with the Air Force Rocket Propulsion Laboratory, Edwards AFB, CA 93523. The report documents the development of gas/particle radiative transfer model techniques as part of the Standardized InfraRed Radiation Model (SIRRM) during the time period from September 18, 1978 to June 30. 1979, under Air Force Project Task 305813BB.

The program manager and principal investigator is C. B. Ludwig and the co-investigators of this phase of the program covered in this report are G. N. Freeman and W. Malkmus (Photon Research Associates, Inc.) and M. Slack and R. Reed (Crumman Aerospace Corporation).

This report has been reviewed, and is approved for publication in accordance with the distribution statement on the cover, and on the DD Form 1473.

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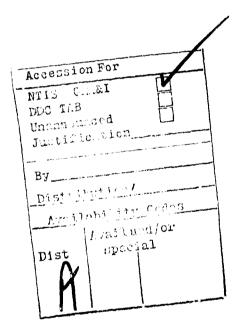
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PREFACE

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- Three-dimensional, non-uniform, axisymmetric medium
- Arbitrary particle mixtures and size distribution
- Line-by-line and band-model treatments
- Alternate levels of engineering approximation

These requirements are met by developing two basic approaches, a 6-flux and a N-flux approach. This distinction refers to the associated subdivision in angular space, six-orthogonal coordinates versus N nonorthogonal coordinates in azimuth and elevation, thus providing the alternatives of computational simplicity of the 6-flux model with generality of the N-flux model. Numerous parametric calculations were performed to study the effect of optical depth, scattering albedo, phase function and medium geometry on the target signature and computational requirements. \Comparisons with independent calculations such as the two-unmensional Monte Carlo method were made. The results show the different degrees of accuracy achievable with the various models as functions of the radiation field parameters. To the authors' knowledge, the N-flux model represents the first application of general finite-element techniques to emitting cylindrical media. It provides a powerful tool for the analysis of plume scattering problems of arbitrary complexity. The accuracy of the prediction is limited only by the computational resources of the user.

The solution of the equation of radiative transfer is derived on a monochromatic basis. For the application to the finite spectral intervals associated with the use of instruments with finite resolution, spectral averaging procedures are required. Thus, an extension of earlier band model formalisms for regular and random arrays of spectral lines was developed to include scattering as well as absorption in a medium in which radiative transfer is described by the two-flux method. Representations of the averaged transmissivity for the Elsasser model and the random model with exponential-tailed S⁻¹ line intensity distribution were derived and numerical results were obtained for large ranges of optical depths of absorbing and scattering media. The effects of inhomogeneities away from the line-of-sight are accounted for by a contribution to the source function which is calculated by a simplified six-flux technique.

A survey of available optical parameters (complex index of refraction as a function of wavelength and temperature) was made for particles of interest. The compilation by Aerospace was found useful as a starting base for a number of relevant species, namely carbon, aluminum oxide, magnesium oxide and zirconium oxide. Data for other species must yet be obtained. The complex indices of refraction for carbon and aluminum oxide were used, together with postulated size distributions, to calculate the volume extinction, scattering and absorption coefficients and the angular scattering phase function by the Mie Theory for homogeneous, spherical particles in the 1-25µm wavelength region. Based on these results, a tentative plan was developed for the organization of the Data File and its integration into the overall code architecture.

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INTRODUCTION

Validated models to predict rocket plume IR signatures must be capable of treating a wide range of propulsion and trajectory parameters as well as handling a number of sensor-related variables. While significant advances have been made in the state-of-knowledge of plume-related radiative transport phenomenologies during the last ten years, particularly in regard to advanced liquid propellants, prediction methods have not been developed to adequately treat the coupling of gas and particle radiative transfer in the plumes of solid propellant missiles. It is particularly important that models be developed to accurately treat plumes containing scattering particles while at the same time also yielding accurate results for conventional liquid propellant plumes.

The primary goals of the SIRRM program are (i) to develop an analytical methodology to treat coupled gas/particle radiative transport effects and (ii) to integrate this analytical capability together with the state-of-the-art treatments in hot gas/ atmospheric radiative transport (i.e., band model methodologies, band model parameters, line-by-line methodologies, spectral line parameters) into a comprehensive, user-oriented plume radiation model.

State-of-the-art plume signature predictions require four basic steps. First, the species concentrations and temperature in the exit plane are calculated through a combustion/nozzle code. Second, these exit plane parameters are input to a gasdynamic code which calculates the exhaust plume flow field. Third, the flow field serves as input to a radiation code that calculates the intrinsic plume signature. Finally, the apparent plume signature viewed by a remote observer sensor is obtained through attenuating the source intensity by the atmospheric absorption. When the source radiation is correlated with the atmospheric attenuation (radiating plume species are also present in the atmosphere as absorbers), the last two steps are combined into a single calculation.

The present study focuses on the third and fourth component, i.e., the calculation of the intrinsic and apparent plume signatures. The program is divided into four major tasks. Task I calls for the formulation of the SIRRM code methodology. In Task II, the code is to be developed and its capabilities demonstrated. The validation and application of SIRRM takes place in Task III. The documentation and code acceptance is done in Task IV.

Task I is concerned with the development of an integrated physical model which accounts for the essential spectroscopic and geometric effects governing single and multiple photon

element in this task is the theoretical description of coupling phenomena associated with the spectral and spatial redistribution of radiation by scattering. In order to provide a correct description, the model must account for the interaction of emitted, absorbed, and scattered radiation from line and continuum sources along multiple, intersecting paths. The task is divided into two subtasks. Subtask IA deals with the development of the coupled, multi-scattering model, while Subtask IB calls for the formulation of the overall methodology, i.e., the integration of the scattering model with the other parts of the code. This document is an interim technical report describing the development of the coupled multi-scattering model performed under Subtask IA.

Section 2 contains the description of our efforts in developing the multi-scattering model with cylindrical geometry, using both engineering approximations (six-flux) and exact solutions (N-flux). Diagnostic calculations with these models and comparisons of the results with those obtained by independent codes indicate that the required objectives have been met.

Since the present scattering model is developed for the monochromatic case, a parallel study was begun to treat the spectral averaging. The step forward to a band-model formulation coupling particle scattering with spectral line absorption is taken in Section 3, where the Elsasser and random model formulations

are developed for a two-flux scattering model and conceptual approaches for the six-flux and N-flux models are outlined.

The fundamental inputs into any scattering models are the optical parameters of the particles. These are discussed in Section 4, including the Mie theory, the indices of refraction and absorption, the scattering due to single particles and particle size distributions, and the handling of the Data File as it would be integrated into the overall code architecture.

SCATTERING MODEL

2.1 Introduction

The objective of Task IA is to formulate a scattering model that incorporates the following capabilities:

- Coupled emission, absorption and scattering treatment
- Multiple scattering with cylindrical geometry
- Three-dimensional, non-uniform, axisymmetric medium
- Arbitrary particle mixtures and size distributions
- Arbitrary angular scattering phase function
- Line-by-line and band-model treatments
- Alternate levels of engineering approximation

The approach to meet these objectives is described in this section. Our efforts included (i) the selection of physical models and solution methods for the gas/particle radiation transfer problems, (ii) the justification of selected models and methods by means of diagnostic calculations, and (iii) the final development of the selected models and methods for integration into SIRRM.

In Section 2.2, we provide the requisite technical back-ground, beginning with the equation of radiative transfer and its implicit and explicit solution schemes.

In Section 2.3, we review and assess the combined interdisciplinary experience of contributors in several fields (e.g., planetary and stellar atmospheric physics, radiant heat transfer, plume physics) where multiple scattering techniques have reached an advanced stage of development. Based on this assessment we select the most appropriate approach to the plume problem, and justify this selection on the basis of key physical requirements and model capabilities. The discussion presented in this section formed one of the essential elements of the Program Plan (Data Item 1).

The other essential element of the Program Plan was a set of recommendations for the final development of the scattering model to be incorporated into SIRRM. Since January 25, 1979, the recommendations were followed and the results of developing engineering approximations and the N-flux model are reported in Sections 2.4 and 2.5, respectively, together with numerical results of diagnostic and parametric calculations.

The comparisons of the computational results between the different models developed in the present program and with independent model calculations are made in Section 2.6. Conclusions about the models and the results are summarized in Section 2.7.

2.2 Scattering Theory

The transfer of radiant energy through matter that emits, absorbs, and scatters radiation was first described by astrophysicists at the beginning of this century. The governing

equation was derived by specializing the Boltzmann equation of kinetic theory to the case of photons. To this day, the original formulation remains essentially unchanged. However, the range of problems amenable to exact solution is quite small, and for most cases of interest an approximate treatment is required.

The formal solution to the equation of transfer provides the essential framework for development of a coupled gas/particle radiation model. The task is to obtain a practical mathematical solution based upon reasonable physical approximations to the spectral and spatial characteristics of the medium. In this section we review the available computational methods in order to identify promising approaches for the problem at hand.

A comprehensive discussion and development of the radiant transfer equation was given by Viskanta (1964). This equation may be written in the integro-differential form

$$\frac{dN_{\lambda}(\vec{s})}{ds} = -n\left(\sigma_{\lambda}^{(a)} + \sigma_{\lambda}^{(s)}\right)N_{\lambda}(\vec{s}) + n\sigma_{\lambda}^{(a)}N_{\lambda}^{0}(s) + \frac{n\sigma_{\lambda}^{(s)}}{4\pi} \int_{4\pi}^{(s)} P_{\lambda}(\vec{s}', \vec{s})N_{\lambda}(\vec{s}') d\Omega(\vec{s}') \tag{2-1}$$

where

 $N_{\lambda}(\vec{s})$ = spectral radiance at point s in direction \vec{s}

 $P_{\lambda}(\vec{s}',\vec{s})$ = phase function for scattering from \vec{s}' to \vec{s}

 $N_{\lambda}^{O}(s)$ = blackbody spectral radiance at point s

 $\sigma_{\lambda}^{(a)}, \sigma_{\lambda}^{(s)}$ = cross sections for absorption and scattering

n = number density

The terms on the right-hand side represent, respectively, the extinction (absorption plus scattering-out) and production (emission plus scattering-in) of radiation in the direction \vec{s} . The phase function is normalized such that

$$\frac{1}{4\pi} \int P_{\lambda}(\vec{s}', \vec{s}) d\Omega(\vec{s}') = 1 \qquad (2-2)$$

The absorption and scattering probabilities $(n\sigma_{\lambda}^{(a)})$ and $n\sigma_{\lambda}^{(s)}$ and the volumetric emission $(n\sigma_{\lambda}^{(a)}N_{\lambda}^{o})$ are defined to include implicitly the effect of multiple gas and particle species (i) and temperatures (T_i) through the relations

$$n\sigma_{\lambda}^{(a)} = \sum_{i} n_{i}\sigma_{\lambda i}^{(a)}$$

$$n\sigma_{\lambda}^{(s)} = \sum_{i} n_{i}\sigma_{\lambda i}^{(s)}$$

$$n\sigma_{\lambda}^{(a)}N_{\lambda}^{o} = \sum_{i} n_{i}\sigma_{\lambda i}^{(a)}N_{\lambda i}^{o}$$

$$n\sigma_{\lambda}^{(s)}P_{\lambda} = \sum_{i} n_{i}\sigma_{\lambda i}^{(s)}P_{\lambda i}$$

The effect of different particle sizes is also implicitly included.

The formal solution to the equation of transfer between points \mathbf{s}_{O} and \mathbf{s}_{O} may be written in the form

$$N_{\lambda}(\vec{s}) = N_{\lambda}(\vec{s}_{0})T(s_{0},s) + \int_{s_{0}}^{s} S_{\lambda}(s') \frac{dT(s',s)}{ds'} ds' \qquad (2-3)$$

where the source function (S_{λ}) at point s' in the direction s' and the transmittance (T_{λ}) from s' to s are given by

$$S_{\lambda}(\vec{s}') = (1-\omega_{\lambda})N_{\lambda}^{O} + \frac{\omega_{\lambda}}{4\pi} \int_{4\pi} P_{\lambda}(\vec{s}'', \vec{s}')N_{\lambda}(\vec{s}'') d\Omega(\vec{s}'')$$
 (2-4)

$$T_{\lambda}(s',s) = \exp\left[\tau_{\lambda}(s') - \tau_{\lambda}(s)\right]$$
 (2-5)

and where the optical depth for extinction and the albedo for single scattering are defined as

$$\tau_{\lambda}(s) = \int_{0}^{s} n(\sigma_{\lambda}^{(a)} + \sigma_{\lambda}^{(s)}) ds'$$

$$\omega_{\lambda} = \frac{\sigma_{\lambda}^{(s)}}{\sigma_{\lambda}^{(a)} + \sigma_{\lambda}^{(s)}}$$
(2-6)

This formal solution is not explicit in that the local radiance is seen to depend on the radiance distribution throughout the medium which is not known at the outset. Explicit solutions can be obtained only for special cases as noted below.

In the absence of scattering $(\sigma_{\lambda}^{(g)} = 0)$, the formal solution to the equation of transfer assumes the explicit form

$$N_{\lambda}(\vec{s}) = N_{\lambda}(\vec{s}_{0})T(s_{0},s) + \int_{s_{0}}^{s} N_{\lambda}^{0}(s') \frac{dT_{\lambda}(s's)}{ds'} ds'$$

where the transmittance between points s' and s is given by

T (s',s) = exp
$$\left[-\int_{s'}^{s} n\sigma_{\lambda}^{(a)} ds''\right]$$

The evaluation of this equation on a monochromatic or spectrally-averaged basis is the end object of numerous existing line-by-line or band model radiation codes.

The finite difference form of the formal solution to the equation of transfer may be written directly from the preceding integral equations. Thus, the radiance at position k in direction i is given by

$$N_{ik} = N_{i,k-1}^{T_{k-1,k}} + \frac{1}{2}(S_{ik} + S_{i,k-1}) (1-T_{k-1,k})$$

$$(i=1,2,\ldots,1,k=1,2,\ldots,K)$$
(2-7)

where the source function is given by

$$S_{ik} = (1-\omega_k)N_k^0 + \omega_k \sum_{j=1}^{I} P_{ji}N_{jk}$$
 (i=1,2,...,I;k=1,2,...K)

and where the transmittance and phase integral are given by

$$T_{k-1,k} = \exp \left[(\sigma_{k-1}^{(a)} + \sigma_{k-1}^{(s)}) n_{k-1} s_{k-1} - (\sigma_{k}^{(a)} + \sigma_{k}^{(s)}) n_{k} s_{k} \right]$$

$$(k=1,2,\ldots,K)$$

$$P_{ji} = \frac{1}{4\pi} \int_{\delta\Omega_{j}} P(s_{j}, s_{i}) d\Omega(s_{j})$$

$$(i, j=1, 2, ..., I)$$

In these equations, the subscripts i,j,k denote the direction \vec{s}, \vec{s}' and position s, respectively, and the subscript λ has been deleted for clarity. With appropriate boundary conditions on the radiance $N_{ik}(k=0,K;i=1,2...,I)$, which include the effects of any external sources, these equations constitute a closed set.

Solution techniques may be grouped into two fundamental classes: implicit or explicit. The implicit methods solve the coupled equations for N_{ik} and S_{ik} in an iterative manner starting from an assumed initial condition. A convenient procedure is to define an initial source function equal to the known thermal emission component $N_{ik}^{(1)}$ and corrected source function $S_{ik}^{(2)}$ directly. This procedure is then repeated to convergence according to the iteration formulae

$$N_{ik}^{(n)} = N_{i,k-1}^{(n)} T_{k-1,k} + \frac{1}{2} (S_{ik}^{(n)} + S_{i,k-1}^{(n)}) (1-T_{k-1,k})$$

$$S_{ik}^{(n+1)} = S_{ik}^{(n)} + \omega_k \sum_{j=1}^{I} P_{ji} N_{jk}^{(n)}$$

$$(1-T_{k-1,k}) = \sum_{i=1}^{I} (1-T_{k-1,k}) + \sum_{i=1}^{I$$

where the superscript (n) denotes the number of iterations or, equivalently, the order of scattering.

The explicit methods solve a single matrix equation for N_{ik} obtained by direct substitution of the equation for S_{ik} . After rearrangement of terms, the equation may be written in the form

$$\sum_{j=1}^{I} \left\{ N_{jk} (2\delta_{ji} T^{-1}_{k-1,k} - \omega_{k} P_{ji}) - N_{j,k-1} (2\delta_{ji} + \omega_{k-1} P_{ji}) \right\} =$$

$$(i=1,2,\ldots,I;k=1,2,\ldots,K)$$

$$(1-\omega_{k}) N_{k}^{0} + (1-\omega_{k-1}) N_{k-1}^{0}$$

$$(2-9)$$

where $\delta_{j,i}$ is the Kronecker delta.

Implicit and explicit methods offer different relative advantages depending upon the optical and geometric properties of the medium. These properties control the number of iteration steps (implicit methods) or the size of the inversion matrix (explicit methods) required to obtain a given level of accuracy.

2.3 Review, Assessment and Selection of Multiple Scattering Formulation

2.3.1 Review and assessment. In the presence of scattering, explicit solutions can be obtained only for suitable approximations to the radiation flux or to the

phase function. Approximations to the radiation flux include the optically thin limit (in which the medium is not dense enough to absorb or scatter radiant energy from other elements in the volume) and the optically thick limit (in which the medium is so dense that radiant energy from other elements does not penetrate). These limits are of interest only as test cases for a more general model of arbitrary optical opacity. Approximations to the phase function include the discreteordinate method, the moment method, and the method of expansion in orthogonal functions (Krook, 1955) from which the well-known two-stream approximations of Schuster-Schwarzschild and Milne-Eddington (Viskanta, 1964) may be derived as special cases. In addition to these explicit methods, various implicit solution techniques have been developed. These techniques typically construct a convergent solution by incremental steps in space or time.

A summary of available multiple scattering solution techniques is given in Table 2-1. These methods can be roughly grouped into four general classes: superposition, transformation, convergence, and matrix. Each of these is discussed and developed in detail in the survey papers by Hansen and Travis (1974) and Irvine (1975). The identifications are necessarily somewhat arbitrary in that some interrelated techniques are included and some distinguishable subsets are excluded. Tables 2-2 and 2-3 list selected key investigations according to solution technique and chronology, respectively. The former was compiled

TABLE 2-1. METHODS OF SOLUTION FOR MULTIPLE SCATTERING

MINE	Description	BENEFIT	DAMMPCE
	1. Supe	rposition Methods	
Layer addition	Superposition of layers of known transmission and reflection	Wide application for planetary atmospheres	Non-emitting madium; plans- parallel slab
Invariant imbedding	Superposition of optically thin layers	Non-homogeneous atmospheres	Long computations
Synthesis	Superposition of volume elements	Generality	Inefficient for symmetric media
	2. Trans	formation Methods	
X and Y functions	Mathematical transformation of variant imbedding	Single angle dependence	Reyleigh scattering
Spherical harmonica	Phase intogral replaced by sum of sphurical harmonics	General applicability	Large matrix
Eigenfunction	Phase integral replaced by sum of orthogonal eigenfunctions	Mathematically rigorous	Numerically prohibitive
	3. Cons	vergence Methods	
Iteration	Successive down- ward and upward integrations to convergence	Simple concept	Low opacity
Successive orders	Superposition of solutions for different orders of scattering	Simple concept	Low opacity; non-emitting medium
Monte Carlo	Individual photon histories traced and superposed	Simple concept; general applicability	Long computations
	ë. Ma	trix Hethods	
Discrete ordinate	Phase integral replaced by finite sum	General applicability	Large matrix
Pinite difference	Operates on differential equation	Componetional experience	Large matrix unless symmetric
Pinite • lement	Operates on integral equation	Reduces to two-stream limit	Large matrix uhiess symmetric

TABLE 2-2. REFERENCES FOR MULTIPLE SCATTERING METHODS

DORTEM	REFERENCES	
Layer Addition	van de Hulst, et al. (1963, 1968), Twomey, et al. (1966), Hansen (1969, 1971), Hovenier (1971), Lacis and Hansen (1974), Grant and Hunt (1969, 1971), Plass, et al. (1973), Preisendorfer (1965).	
Successive Orders	van de Hulst and Irvine (1962), Irvine (1964, 1965), Uesugi and Irvine (1970), Poon and Ueno (1974), Dave (1964), Hovenier (1971), Chou (1978).	
Invariant Imbedding	Uesugi and Irvine (1970), Bellium, et al. (1960, 1963, 1967, 1969), Wing (1962).	
Iteration	Herman, et al. (1963, 1965, 1970), Dave, et al. (1970), Crosbie and Linsenbardt (1975).	
X and Y Functions	van de Hulst (1970), Sobolev (1972, 1974), Carlstedt and Mullikin (1966), Pahor and Kuscer (1966), Busbridge (1960, 1967).	
Discrete Ordinate	Chandrasekhar (1950), Jefferies (1955), Chu and Churchill (1955), Lenoble (1956), Chin and Churchill (1965), Samuelson (1969), Liou (1973, 1974), Weinman and Guetter (1972), Whitney (1972, 1974), Kofink (1967), Love and Grosh (1965).	
Spherical Harmonics	Canosa and Penafiel (1973), Davison (1958), Case and Zweifel (1967), Deuze, et al. (1973), Dave, et al. (1974).	
Eigenfunctions	Case and Zweifel (1967), Kuscer and McCormick (1973, 1974), Kaper, et al. (1970), Mika (1961).	
Monte Carlo	Plass and Kattawar (1968, 1971), Panielson, et al (1968, 1969), Sanford and Pauls (1973), Collins, et al. (1965, 1972), Van Blerkom, et al. (1971, 1974), McKee and Cox (1974), Kattawar, et al. (1971, 1973), Marchuk and Mikhailov (1967).	
Diffusion	Heasley (1977), Barkstrom and Arduini (1978), Laung (1975, 1976)	
Finite Difference	Barkstrom (1976).	

TABLE 2-3. CHRONOLOGY OF MULTIPLE SCATTERING METHODS

YEAR	PEFERENCE	PHYSICAL SITUATION	PETHOD OF ANALYSIS
1955	Jefferies	Point and infinite line sources	Diffusion approximation
''''		in a semi-infinite medium	and and approximation
1955	Jefferles	A strip of collimated radiation incident on a semi-infinite medium	Discrete ordinates and Fourier transform
1955	Churchill	Point source outside an aniso- tropically acattering finite layer	Six-flux erproximation
1956	Giovanelli	An infinite line source on the surface or within a semi-infinite medium and a thick finite layer	Approach of Jefferies
.959	Chandrasekhar	An infinitely narrow collinated beam incident on a semi-infinite medium	loveriance principles
1963	Bollman, Falaba, and Uene	An infinitely narrow collimated beam incident on a finite laver with properties varying with depth	Invariant imbedding
1963	Eellman, Kalaba, and Ueno	Collinated radiation incident on an inhomorpheous animotropically scattering finite layer	Invariant imbedding
1964	Smith	A uniferm strip of cullimated radiation incident on a semi-infinite medium	Assumed cosine-varying in- cident radiation to reduce the two-dimensional in- tegral equation for the source function to a one- dimensional form
1954	Malkevich, Monin, and Rozenberg	Collimated radiation incident on a finite layer with nonuniform reflection from lower surface	Horizontal variation of downward radiation neglected
1965	Chin and Churchill	An arbitrary, cylindrical source in an anisotropically scattering finite slab	Six-flux approximation
1966	Dolin	Narrow collimated beam incident on a semi-infinite anisotropi- cally scattering medium	Small angle transport equa- tion used for unscattered radiation, diffusion equa- tion used for scattered radiation
1967	Weinman and Swarztrauber	Collimated radiation incident on a nonabsorbing finite layer with horizontally non-uniform scattering coefficient	Giovanelli's approximation used to molve transport equation
1967	Smith and Hunt	A uniform strip of collimated radiation incident on a semi- finite layer	Same as Smith 1964
1967	Hunt	Collimated radiation incident normally on an anisotropically scattering finite medium	Expression of scattering function as series of Legendre polynomials
1967	Bobco	An isothermal semi-infinite simb	Diffusion equation solved by iterative technique
1967	Drobyshevich	Collimated radiation incident on an anisotropically scattering finite layer with nonuniform reflection from lover surface	Forizontal variation of downward radiation neglected
1968	Remanova	An infinitely nerrow collimated bear normally incident on a semi-infinite anisotromically scattering medium	Three-dimensional moments used to solve the trans- port equation for large depths
1968	Pomanova	An infinitely narrow collimated bear normally incident on an an- isotropically scattering finite layer	Frall apple approximation to the transport equation solved using spatial roments

TABLE 2-3. (CONTINUED)

TRAN	PEPERENCE	PHYSICAL SITUATION	PETHOD OF AMALYSIS
1968	Punk	Collimated woint source normally incident on an anisotropically scattering finite layer	Monte Carlo technique
1568	Funt	Collimated Pessel function varying radiation incident on a finite layer	Integra) equation for the source function formulated using Green's function
1968	Stockham and Love	An isothermal finite cylindrical medium: both isotropic and ani- sotropic scattering	Yonte Carlo technique
1969 -	Punt	Collimated radiation normally incident on a finite layer	Integral equation for source function trans- formed into singular equation for the re- solvent kernel using properties of a Green's function
1769	Bravo-Zhivatovskiy, Dolin, Luchinin, and Savel'yev	Nerrow collimated bear normally incident on a semi-infinite an-isotropically scattering redium	Uses molution to transport equation found in Dolin 1966
1969	Solubitably and Tantosnev	An infinitely narrow collinated beam at the center of an aniso-tropically scattering apherical medium	Monte Carlo technique
1970	Love and Turner	An imothermal semi-infinite alab	Fonte Carlo technique
1970	Scofield and	A strip of diffuse radiation in- cident on an eritting seri-in-	Monte Carlo technique
		finite medium; both isotropis and anisotropic scattering	
1971	Rybick1	An infinitely narrow collimated beam incident on a finite layer and semi-infinite medium	Fourier transform and in- variance techniques
1971	Romanova	An infinitely narrow collimated beam normally incident on a non-absorbing finite layer	Three-dimensional invariance principles used to obtain equations for spatial moments of intensity
:971	Romenova	An infinitely narrow collimated beam normally incident on an an- isotropically acattering finite layer	Ipatial moments used to solve small angle approximation to the transport equation
1971	Kochetkov	An infinitely narrow collimated been normally incident on an an- isotropically scattering finite layer	Pethod of spherical harmonica
1972	Dushmakova, Zoge, and Katsev	Diffuse and unidirectional acurces inside an anisotropical-ly scattering finite layer and semi-infinite medium	Pankel transformation used to solve diffusion equation
1973	Romanovs	An infinitely narrow collimated beam normally incident on a finite layer with weak absorption	Intensity represented by trans- verse noments and invariance principles used
1973	Punk	A narrow collimated beam and noncollimated source in an in- finite anisotropically scatter- ing medium	Monte Carlo technique
1973	Gordon and Fnittel	An infinitely narrow collimated bear normally incident on an an- instropically scattering finite layer	Iterative ray tracing, no re- turn backscattering and small angle approximations used
1974	Crusbie and Koewing	A step and finite strip of col- limated and diffuse radiation incident on a finite layer	Assumption of cosine wary- ing incident radiation re- duces problem to one-di- mensional problem; Ambarzumian's method
1974	Beckett, Foster, Mutson and Moss	A collimated cylindrical bram normally incident on a finite layer	Hunt's interral equations solved by iteration

from previous review papers together with our own review; the latter was taken directly from Crosbie and Linsenbardt (1976). Neither list is comprehensive, but both serve to illustrate the depth and scope of previous work. Most of these studies applied to non-emitting plane-parallel atmospheres. Only a limited number of the studies included the effects of thermal emission, inhomogeneous media, non-isotropic scattering, and non-planar geometry which are important in the present application. None of the studies treated all these effects simultaneously. Despite these deficiencies, this previous work provides a firm base for the selection and development of a multiple scattering model for two-phase plumes.

Existing rocket plume radiation codes which treat particle radiation were also reviewed and are presented in Table 2.4,

The early codes of Fontenot (1965), Gulrajani (1964), Hunt (1966) and the Aeronutronic group (Carson, 1965) were all developed to predict base heating. The more recent plume codes of Wilson (1973), Rieger (1974), Vanderbilt and Slack (1976), and the ARAP group (Fishburne, 1977) were developed to predict plume IR signatures. These codes approximate particle radiation effects by either of two methods: Pseudo-gas or one-dimensional beam. The pseudo-gas approximation includes particle emission/absorption effects but neglects

TABLE 2-4. EXISTING ROCKET PLUME RADIATION CODES INCLUDING PARTICLES

PLUME CODES	METHOD	COMMENTS
Fontenot (1965)	Inverse wavelength method	Simplistic approach; single particle size neglects scattering. Estimates base-heating for constant radial properties.
Gulrajani (1964) Hunt (1966)	Neutron-scattering analogy method	Based on Morizumi and Carpenters (64) analysis. Five particle sizes. Includes 1-D multiple scattering. Neglects anisotropic scattering effects. Assumes $Q_t \rightarrow 2$. Base-heating estimates.
Aeronutronic (Carlson, 1965)	One-dimensional beam approximation	Plume codes to analyze Saturn ullage retro and strap-on exhaust plumes. 1-D anisotropic scattering approximation.
Wilson and Hahn (1973)	One-dimensional beam approximation	Uncoupled gas and particles. Adds separately computed particle radiation and gas radiation. Narrow inband resolution. Treats 5 particle sizes.
Rieger (1974)	Pseudo-gas approxi- mation	Computes coupled particle and gas emission and absorption.
Vanderbilt & Slack (1976)	One-dimensional beam approximation	Couples multiple scattering and gaseous emission/absorption along beams. Spectral resolution 5 cm $^{-1}$. Treats 5 particle sizes. Treats gases as non-scattering particles yet retains correct solution of radiative transfer equation for gases alone and particles alone. Spectral and inband output coupled to LOWTRAN codes. Dave code for Q_s , Q_a
ROCRD Code (ARAP)	Pseudo -gas approximation	Includes particle emission, but neglects scattering. Assumes Beer's law.

scattering. The one-dimensional-beam approximation includes particle scattering effects along the beam but neglects scattering into the beam from other directions. Of the two methods. the one-dimensional beam provides a better physical description of particle radiation effects with no appreciable penalty in computational complexity. Wilson used the one-dimensional beam approximation to compute the particulate radiation field in an uncoupled manner for subsequent addition to the gaseous radiation field; the validity of this procedure is restricted to low optical depths. Vanderbilt and Slack used the one-dimensional beam approximation to compute the coupled particulate and gaseous radiation fields, thereby allowing the gas to absorb the particulate continuum and the particles to scatter and absorb the gaseous radiation. The code developed by Vanderbilt and Slack is the best of the existing codes for predicting spectral IR signatures from missile plumes. However, the lack of coupling between adjacent 1-D beams means that it does not meet the SIRRM code requirements, in general.

One dimensional beam methods have received extensive application and limited verification in the area of radiative heat transfer but only limited application and no verification in the area of radiative signatures. This distinction is important because of a fundamental difference between the two phenomena. On the one hand, the radiative heat transfer to an adjacent point is the sum of contributions from non-parallel beams. On the other hand, the radiative signature to a remote point is the sum of contributions from parallel beams. The

latter effect is much more sensitive to errors in the directional properties of the radiation field. Thus, the SIRRM code must include a three-dimensional scattering model in order to achieve the required directional accuracies.

The state of the s

2.3.2 Selection and justification. As noted previously, the SIRRM scattering model for two-phase plumes must include the following key physical effects: thermal emission, inhomogeneous media, non-isotropic scattering, and non-planar geometry. In addition, the SIRRM scattering model must possess the following computational attributes: accuracy, efficiency, and generality. These general attributes have rather specific meanings for the present application as noted below:

Accuracy The chosen technique should incorporate sufficient spatial resolution to yield an accurate mathematical solution. The technique should demonstrate adequate agreement with the predictions of alternative methods for suitable test cases.

Generality The chosen technique should incorporate sufficient spatial dimensions to represent the real physical problem. The technique should treat all ranges and variations of optical depth, scattering albedo, and phase function such that no limit is imposed on the nature of gas or particle properties, plume properties, or observation geometry.

Efficiency The chosen technique should yield computed results in machine times which are comparable to current gaseous emission codes for an equivalent number of aspect angles. Approximate techniques which incorporate physically acceptable assumptions should be evaluated in order to determine the most efficient technique for different classes of problems.

The available methods were screened according to these criteria.

The dual requirement for engineering efficiency and scientific accuracy, together with the need for computational generality, and flexibiltiy, led to the selection of straightforward, finite difference procedures of variable spatial and angular resolution applied to locally cylindrical plume geometries. Two basic approaches (herein identified as "6-flux" and "N-flux") were distinguished and various alternatives were developed within each classification. The semantic distinction refers to the associated subdivision of angular space (6 orthogonal coordinates versus N non-orthogonal coordinates in azimuth and elevation). However, the fundamental distinction lies with the generality of the N-flux model versus the computational simplicity of the 6-flux model.

Within the 6 flux model class, two independent sub-class models were developed. The first model employs suitable approximations for the axial and tangential radiances in order to

reduce the solution to a one-dimensional radial integration similar to existing two-stream models (Vanderbilt-Slack, 1976). The second model employs a transformation of the multiple-scattering source function in order to reduce the solution to a coupled set of radial, tangential, and axial integrations, each of which is functionally similar to existing two-stream models. The two six-flux models are described and compared in Section 2.4.

The N-flux model class includes a generalized set of optional finite-element grid geometries including independent specification of the angular resolution in the azimuth and elevation planes. The model solves the coupled three-dimensional equations of transport using a generalized matrix network to describe the radiance distribution and source function at each grid point. The model can be exercised as either an engineering or scientific tool by simple variations in the spatial and angular grid network to achieve low or high resolution. Analytical details are presented in Section 2.5.

Each of the models was exercised parametrically over the full physical range to illustrate the effect of optical depth, scattering albedo, and phase function on the local and integrated target signature characteristics. These calculations provided independent tests of the physical credibility and computational efficiency of the separate models as well as comparative tests of the accuracy and self-consistency of the composite model set.

Additional verification was accomplished by performing independent calculations with the NASA plane-parallel slab code (Barkstrom, 1976) modified to include self-emission and with the Grumman Monte Carlo code. Results of these calculations are summarized in Section 2.6.

2.4 Six-Flux Engineering Approaches

The basis for the engineering approaches developed under

Task IA is the approximation of the multiple-scattering source

function by a six-element orthogonal angular quadrature. In

this approximation, the general differential equation of transport

$$\frac{dN(\tau,\Omega)}{d\tau} = -N(\tau,\Omega) + \left|1-\omega(\tau)\right| N^{O}(\tau) + \frac{\omega(\tau)}{4\pi} \int_{4\pi}^{\Phi} P(\tau,\Omega,\Omega')N(\tau,\Omega')d\Omega' \quad (2-10)$$

is written for each orthogonal direction (i = 1,2,3) in the specific form

$$\pm \frac{dN_{i}^{\pm}}{d\tau_{i}} = -N_{i}^{\pm} + (1-\omega)N^{O} + \omega \left[fN_{i}^{\pm} + bN_{i}^{\mp} + s \sum_{i'\neq i} (N_{i'}^{+} + N_{i'}^{-}) \right]$$
 (2-11)

where the terms on the right-hand side represent the respective contributions of extinction, emission, and multiple scattering to the radiance gradient and where the terms f,b,s represent the forward, backward, and sideward components of the phase function (f+b+4s=1). This quadrature provides the minimum angular

resolution necessary to account for three-dimensional multiple-scattering effects. For comparison to previous approaches, the standard two-element, one-dimensional approximation is obtained as a special case (s=0).

The orthogonal coordinates are chosen to coincide with the natural cylindrical coordinates (r,θ,z) of the axisymmetric plume. Figure 2-1 illustrates the grid network formed by the intersections of these coordinates with the boundaries of the finite elements (radial and axial). The six-flux equations are solved as a coupled set along these mutually-perpendicular integration paths. Radiances for oblique lines-of-sight are computed in a subsequent step based upon the known six-flux solution.

Within the six-flux model class, two different levels of engineering approximation were employed. The first method yields an approximate explicit solution based upon elimination of the tangential (θ) equations of transport in favor of simple closure relations. The second method yields an exact implicit solution based upon iteration or inversion of the fully-coupled multi-dimensional equation set. Both methods were originally developed for infinite cylindrical geometries (negligible axial gradients). The second method was subsequently extended to finite cylindrical geometries in order to assess the influence of plume axial gradients on the radiance at side aspect. The following paragraphs provide analytical details.

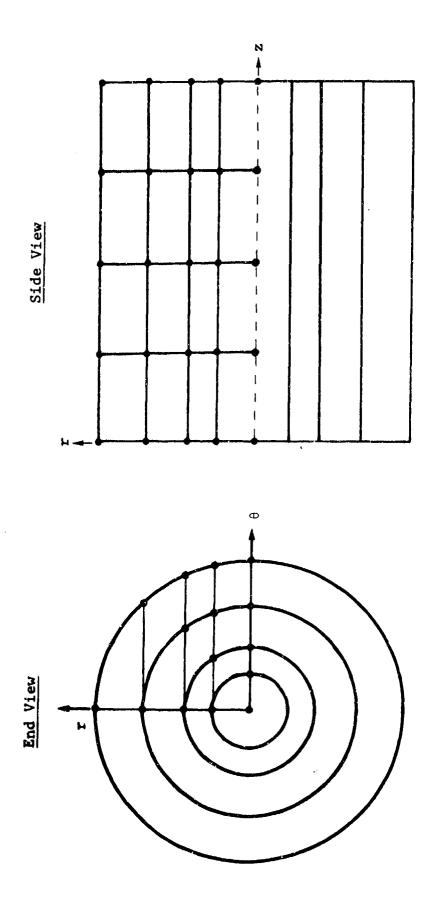


Figure 2-1. Six-Flux Model Coordinate System.

2.4.1 Method I. This method is based upon a transformation of the three-dimensional equations of radiative transport to equivalent one-dimensional form for infinite planar or cylindrical media. The resulting equation is integrated along a piecewise-uniform path perpendicular to the surface to yield an explicit solution for the normal and tangential radiances throughout the medium. Radiances at intermediate angles are determined by analytical interpolation. No restriction is placed upon the range of physical properties considered (e.g. optical depth, scattering albedo, phase function) or upon the variation of those properties within the medium.

The ability to treat planar and cylindrical geometries in a self-consistent manner provides an important connection between the available literature (mostly planar) and the plume problem (nearly cylindrical). As a result of this connection, the utility of existing plane parallel codes (e.g., Barkstrom, 1976) for defining baseline parametric trends is greatly enhanced. In addition, solutions for planar media which are infinite in one or two directions (i.e., ribbon or slab, respectively) provide convenient bounds on the emitted radiation from equivalent cylindrical media.

In the rectangular r,θ,z coordinate system of Figure 2-1, the equation of transfer reduces to the six simultaneous equations

$$\frac{dN_{r}^{\pm}}{d\tau_{r}} = -N_{r}^{\pm} + (1-\omega)N^{O} + \omega \left[fN_{r}^{\pm} + bN_{r}^{\mp} + s(N_{\theta}^{+} + N_{\theta}^{-} + N_{z}^{+} + N_{z}^{-})\right]$$

$$\pm \frac{dN_{\theta}^{\pm}}{d\tau_{\theta}} = -N_{\theta}^{\pm} + (1-\omega)N^{O} + \omega \left[fN_{\theta}^{\pm} + bN_{\theta}^{\mp} + s(N_{r}^{+} + N_{r}^{-} + N_{z}^{+} + N_{z}^{-})\right]$$

$$\pm \frac{dN_{z}^{\pm}}{d\tau_{z}} = -N_{z}^{\pm} + (1-\omega)N^{O} + \omega \left[fN_{z}^{\pm} + bN_{z}^{\mp} + s(N_{r}^{+} + N_{r}^{-} + N_{\theta}^{+} + N_{\theta}^{-})\right]$$
(2-12)

These six equations are further reduced to two ($\pm r$) by means of suitable analytical approximations for the perpendicular radiances ($\pm \theta$, $\pm z$). The z-equation is simplified through the assumption of negligible axial gradients

$$\left|\frac{dN_z}{d\tau_z}\right| \ll N_z^{\pm}$$

which reduces the equation to simple algebraic form. The θ -equation is eliminated through the assumption of approximate proportionality factors

$$N_{\theta}^{\pm} \approx \Phi_{p} N_{z}^{\pm}$$
 (plane)
 $N_{\theta}^{\pm} \approx \Phi_{c} \{ (N_{r}^{+} + N_{r}^{-}) \text{ (cylinder)} \}$

which are piecewise-constant over a radial integration step. The range $0<\Phi_p<1$ defines the limits of thin and thick optical paths in the θ -direction. In general, the factor Φ is approximated

by the corresponding radiance ratio for a non-scattering medium.

For a uniform medium, this ratio is given simply by the emissivity ratio

$$\Phi_{p} = 1-\exp(-\tau_{\theta})$$

$$\Phi_{c} = \left[1-\exp(-\tau_{\theta})\right] / \left[1-\exp(-\tau)\right]$$

where τ is the total optical depth for a cylinder of diameter 2R and τ_{θ} is the half optical depth measured perpendicular to the line-of-sight. For a uniform cylinder $\tau_{\theta} = \tau [1-(r/R)^2]^{\frac{1}{2}}$. Substitution of these approximations into the axial equation of transport yields the explicit formulae

$$N_{z}^{\pm} = \left[(1-\omega)N^{O} + \omega s \left(N_{r}^{+} + N_{r}^{-}\right) \right] \left[1 - \omega \left(f + b + 2\Phi_{p} s\right) \right]^{-1} \qquad \text{(planar)}$$

$$N_{z}^{\pm} = \left[(1-\omega)N^{O} + \omega s \left(1 + \Phi_{c}\right) \left(N_{r}^{+} + N_{r}^{-}\right) \right] \left[1 - \omega \left(f + b\right) \right]^{-1} \qquad \text{(cylindrical)}$$

With these approximations, the six-flux radial equation of transport may be written in the functional form of the two-flux equation with modified coefficients. That is,

$$\pm \frac{dN_r^{\pm}}{d\tau_r} = -\alpha N_r^{\pm} + \beta N_r^{\mp} + \gamma N^0 \qquad (2-14)$$

where the piecewise-constant coefficients are given by

$$\frac{\text{planar}}{\alpha = 1 - \omega(\text{f+s}\chi)} \qquad \alpha = 1 - \omega\{\text{f+s}[\chi + \Phi_{\mathbf{c}}(1 + \chi)]\}$$

$$\beta = \omega(\text{b+s}\chi) \qquad \beta = \omega\{\text{b+s}[\chi + \Phi_{\mathbf{c}}(1 + \chi)]\}$$

$$\gamma = (1 + \chi)(1 - \omega) \qquad \gamma = (1 + \chi)(1 - \omega)$$

$$\chi = 2s\omega(1 + \Phi_{\mathbf{p}})[1 - \omega(\text{f+b+2}\Phi_{\mathbf{p}}s)]^{-1} \qquad \chi = 2s\omega[1 - \omega(\text{f+b})]^{-1}$$

Integration over the uniform region from station (k-1) to k yields explicit expressions for the outward and inward radiance components at the outer and inner surfaces, respectively.

where the subscript r has been deleted for clarity. These equations contain the boundary conditions (N $_k^-$ and N $_{k-1}^+$), the local source function (N $_k^0$), and the local transport coefficients (reflectivity, transmissivity, and emissivity) defined by the relations

$$\rho = \frac{1-\xi^2}{1-\xi^2\eta^2} \eta$$

$$\xi = \frac{\beta}{\alpha+\lambda}$$

$$t = \frac{1-\eta^2}{1-\xi^2\eta^2} \xi$$

$$\eta = \exp\left[-\lambda(\tau_k - \tau_{k-1})\right]$$

$$\xi = \frac{\gamma}{\alpha-\beta} (1-t-\beta)$$

$$\lambda = \left[(\alpha+\beta)(\alpha-\beta)\right]^{\frac{1}{2}}$$

The axial radiances at the outer and inner surfaces then follow directly.

The formulation for a uniform planar medium is complete at this point. For a non-uniform or cylindrical medium, the formulation is completed by superposition of elemental solutions for K piecewise-uniform subregions (k=1,2,...,K) subject to the overall boundary conditions

$$N_{K}^{-} = 0$$
 (no external sources)
 $N_{O}^{+} = N_{O}^{-}$ (axial symmetry)

The radiance emerging from the K-th (outer) element is determined from Eq. 2-15 by successive Gaussian elimination beginning at the center (k = 0). The general recursion relation (developed by induction) is given by

where

$$A_{k} = \rho_{k} + t_{k} C_{k} A_{k-1} ; \qquad A_{o} = 1$$

$$B_{k} = \rho_{k} N_{k}^{o} + C_{k} B_{k-1} ; \qquad B_{o} = 0$$

$$C_{k} = t_{k} (1 - \rho_{k} A_{k-1})^{-1}$$

$$\rho_{k} = \epsilon_{k} (1 + C_{k} A_{k-1})$$

The surface radiance $(N_K^+ = B_K^-)$ is determined explicitly from the sequential summation.

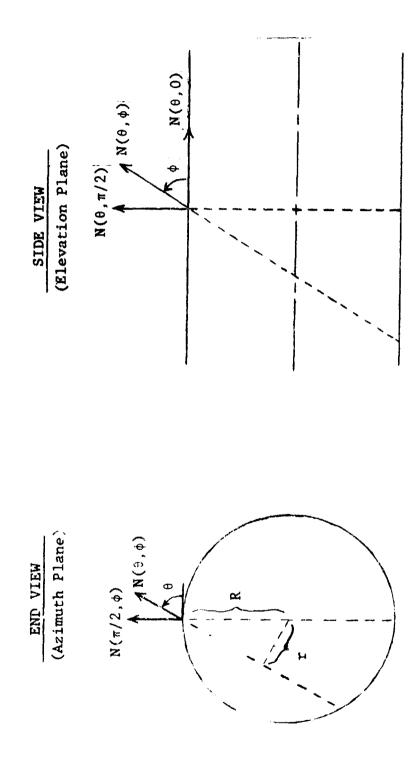
The preceding analysis yields the normal and tangential radiance components for planar and cylindrical media. Between these limits, the variation of radiance with elevation angle (ϕ) or azimuth angle (θ) (standard spherical coordinates) must be determined by a subsequent calculation. The following paragraphs describe a simple approximate procedure based on analytical interpolation. An alternative exact procedure based on the known six-flux source function distribution is described under Method II.

The approximate interpolation formulae are based on weighted sums of the local-to-normal and local-to-tangential radiance ratios for a non-scattering medium. From Figure 2-2

$$N(\theta, \phi) \approx \left[\frac{N(\theta, \phi)}{N(\pi/2, \pi/2)}\right]_{\omega=0} N(\pi/2, \pi/2) \sin^2 \phi$$

$$+ \left[\frac{N(\theta, \phi)}{N(\pi/2, 0)} \right]_{\omega=0} N(\pi/2, 0) \cos^2 \phi$$

where $N(\pi/2,\pi/2)$ and $N(\pi/2,0)$ are the normal and tangential radiances predicted by the six-flux model. The θ and ϕ variations are conveniently separated by rewriting $N(\theta,\phi)$ as the product



Estimation of Three-Dimensional Radiance Distribution from Computed Radiances at Normal and Tangent Aspect. Figure 2-2.

$$N(\theta,\phi) = \left[\frac{N(\theta,\phi)}{(N(\pi/2,\phi))}\right]_{\omega=0} \cdot \left\{ \left[\frac{N(\pi/2,\phi)}{N(\pi/2,\pi/2)}\right]_{\omega=0} N(\pi/2,\pi/2) \sin^2\phi \right\}$$

$$+\left[\frac{N(\pi/2,\phi)}{N(\pi/2,0)}\right]_{\omega=0}N(\pi/2,0)\cos^2\phi$$

This separation facilitates the integration over θ to obtain the local station radiation.

$$J'(\phi) = 2\sin\phi \int_{0}^{r} N(\theta, \phi) dr$$

$$= 2R\sin\phi \int_{0}^{\theta} N(\theta, \phi) \cos\theta d\theta$$

$$= 2R\sin\phi N(\pi/2, \phi) \int_{0}^{\theta} \left[\frac{N(\theta, \phi)}{N(\pi/2, \phi)} \right]_{\omega=0}^{\cos\theta d\theta}$$

For uniform media the radiances for $\omega=0$ are simply proportional to the corresponding emissivities; i.e.,

$$N(\theta, \phi) \sim \varepsilon (\tau \sin\theta / \sin\phi)$$

where τ is the optical depth based on the cylinder diameter. In this case, the radiance is given by the simplified formula

$$N(\theta,\phi) = \varepsilon(\tau \sin\theta/\sin\phi) \left| N(\pi/2,\pi/2) \sin^2\theta/\varepsilon(\tau) + N(\pi/2,0) \cos^2\phi \right|$$

and the station radiation is well-approximated by the empirical formula (see Figure 2-3)

$$J'(\phi) = 2R\sin\phi\varepsilon(\tau/\sin\phi)[N(\pi/2,\pi/2)\sin^2\phi/\varepsilon(\tau)+N(\pi/2,0)\cos^2\phi] \cdot$$

$$\varepsilon(.546\tau/\sin\phi)/\varepsilon(.695\tau/\sin\phi)$$

The non-dimensional ratio $J'(\phi)/2R\sin\phi N(\pi/2,\phi)$ is observed to approach $\pi/4$ or 1 in the optically thin or thick limits, respectively. These interpolation formulae suffice to define the entire radiation field at a given axial station in terms of the computed normal and tangential radiances.

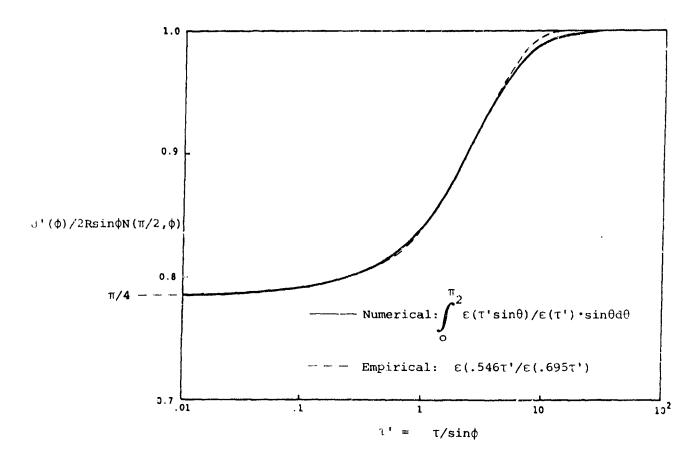


Figure 2-3. Estimation of Station Radiation for Uniform Cylinder from Computer Centerline Radiance.

2.4.2 Method II. This method is based upon a transformation of the three-dimensional equations of transport to equivalent isotropic form for infinite and finite cylindrical media. The resulting equations are integrated along and perpendicular to the radial direction to yield an implicit solution for the isotropic (scalar) source function distribution. Numerical results are obtained by iteration using the method of successive scatterings. Solutions for arbitrary aspect are then computed directly in terms of the known source function by means of a two-stream integration along the observer line-of-sight. As for Method I, no restriction is placed upon the range or variation of physical properties within the medium.

The general six-flux equation of transport (see Eq. 2-11)

$$\pm \frac{dN_{i}^{\pm}}{d\tau_{i}} = -N_{i}^{\pm} + (1-\omega)N^{O} + \omega \left[fN_{i}^{\pm} + bN_{i}^{\mp} + s \sum_{i'\neq i} (N_{i'}^{+} + N_{i'}^{-}) \right] \quad (i=1,2,3)$$

is transformed by simple rearrangement of terms to equivalent isotropic form

$$\pm \frac{dN_{i}^{\pm}}{d\tau_{i}} = -\alpha N_{i}^{\pm} + \beta N_{i}^{\mp} + \gamma (N^{0} + N') \qquad (i=1,2,3)$$
 (2-16)

where the piecewise-constant coefficients are given by

$$\alpha = 1 - \omega(f - s)$$

$$\beta = \omega(b-s)$$

$$\gamma = (1 - \omega)$$

and where the source function for multiple scattering is given by

$$N' = \frac{\omega s}{(1-\omega)} \sum_{i=-1}^{3} (N_{i}^{+}, + N_{i}^{-})$$

The key feature of the transformed equation set is that the source function now contains only the side-scattering component of the phase function such that N' is a scalar quantity. This important simplification is obtained for any arbitrary phase function which can be approximated by a six-angle orthogonal quadrature.

For a known source function N' (as from a previous iteration) the transformed equation of transport is functionally equivalent to Method I and can be integrated as before to yield

$$N_{ik}^{+} = \rho_{ik}N_{ik}^{-} + t_{ik}N_{ik-1}^{+} + \epsilon_{ik}(N_{k}^{0} + N_{k}^{1})$$

$$(i=1,2,3)$$

$$N_{ik-1}^{-} = t_{ik}N_{ik}^{-} + \rho_{ik}N_{ik-1}^{+} + \epsilon_{ik}(N_{k}^{0} + N_{k}^{1})$$
(2-17)

where ρ_{ik} , t_{ik} , ϵ_{ik} are the reflectivity, transmissivity, and emissivity of the element k in the direction i and N_{ik}^{\pm} , N_{ik-1}^{\pm} are the surface radiances. The coefficients ρ , t, ϵ are evaluated as before (see Method I) in terms of the coefficients α , β , γ , where the latter have new definitions as given above. The requisite geometrical relations are given in Section 2.5. For an array of K elements (k=1, 2,...K), the boundary conditions are

$$N_{io}^{+} = N_{io}^{-}$$
 (axial symmetry)
 $N_{iK}^{-} = 0$ (no external sources)

In the direction of integration, these equations form a system of 2K equations for the 2K radiances $N_{\bf k}^{\pm}$ defined over

a K-point linear mesh. An additional 2K equations of similar form are required for each perpendicular direction in order to define the source function at each grid point k. For each path, the transport equation yields a one-dimensional solution matrix which is a diagonal sub-block of the full three-dimensional solution matrix. The individual paths are coupled through the scalar source function at intersecting points such that the complete ensemble of paths is required to obtain closure. The complete matrix is solved by iteration using the single-scattering solution (multiple-scattering source function equal to zero) as the initial guess. For the cases investigated to date (see Section 2.6), two to eight iterations (depending on the product $\omega \cdot \tau$) were sufficient to achieve source function convergence within two percent.

The importance of the isotropic source function transformation can be appreciated by reference to the tangential (0) integration paths in Figure 2-1. Except for the s-axis, the intersections of these paths with the cylindrical shells are observed to occur at oblique angles which do not coincide with the six-angle quadrature of the phase function. This situation presents no difficulty when the source function is isotropic. However, a non-uniform angular quadrature would be required to represent non-isotropic scattering into the line-of-sight. Such an approach is somewhat more complex than the present method, but offers the potential for greater accuracy through higher resolution. This capability is provided by the multiple-flux model described in the following section.

2.5 Multiple-Flux Exact Approach

The basis for the exact approach developed under Task IA is the approximation of the multiple-scattering source function by an angular quadrature of arbitrary resolution. The accuracy of the predicted radiation field increases as the resolution of the angular quadrature increases. Exact solutions are obtained by increasing the resolution in successive calculations until convergence is achieved.

The arbitrary angular resolution of the multiple-flux (henceforth called "N-flux") model affords considerable flexibility in the analysis of multiple-scattering problems. The resolution in the elevation and azimuth planes is independently specified so that individual optima can be established. An important feature of the formulation is the introduction of non-uniform azimuthal resolution which increases as the radius increases in order to achieve uniform accuracy from the isotropic center to the anisotropic edge. An alternative formulation based on uniform azimuthal resolution was considered early in the model development effort, but was abandoned due to its greater complexity and lesser accuracy compared to the present approach.

Figure 2-4 shows end-view and side-view projections of representative lines-of-sight through an assembly of infinite cylindrical shells in which the source function for multiple scattering is assumed constant. In the present development,

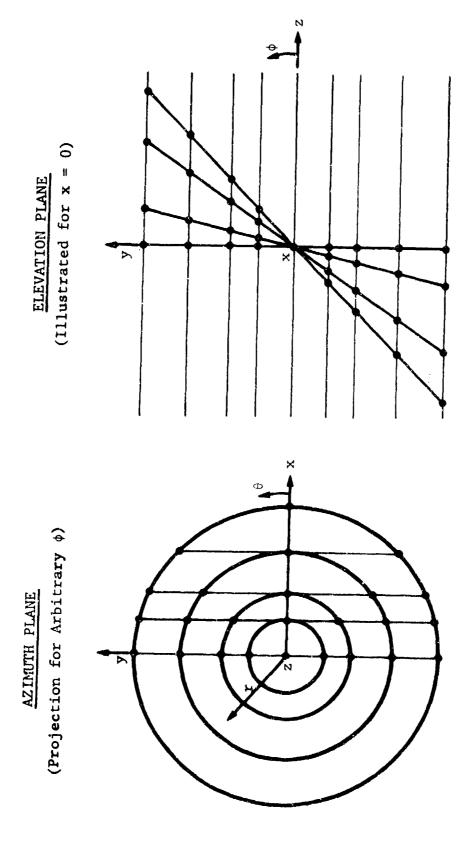


Figure 2-4. Cylindrical Shells and Lines of Sight

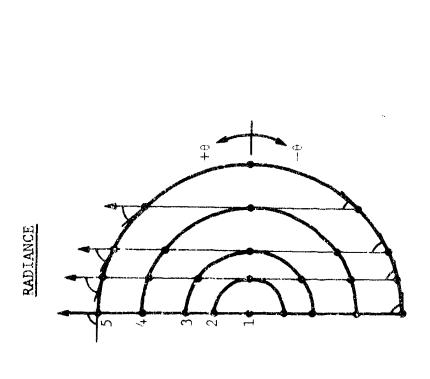
the source function for emission is also assumed constant over the same region, but this is not a general requirement. unknowns are the radiances at the intersections of the cylindrical shells with the parallel lines-of-sight. The constraints are the transport equations for the intervening paths together with the boundary conditions for each line of sight. Together they constitute a determinate linear set. In the figure, the rectangular z, y, x coordinates denote the plume axis, the line-of-sight direction, and the line-of-sight spacing, respectively. The spherical r, Θ , ϕ coordinates denote the radius, azimuth, and elevation of the local radiance vectors. The integration paths are determined by prescribed values of x and ϕ where the x's are chosen to correspond to the radii of the cylindrical shells. The integrations are performed over half of the azimuth plane $(-\pi/2 < \Theta < \pi/2)$ and half the elevation plane $(0 < \phi < \pi/2)$ or onefourth of the radiating volume (π steradians). The remaining volume is described by symmetry relations through simple reflections in Θ and Φ .

Figure 2-5 shows the relationship between the radiances and the source functions at different radial stations which provides the coupling mechanism for multiple lines-of-sight. This relationship is based on the assumption of infinite axial symmetry such that the properties of the medium vary only with radius. Positive or negative radiance vectors are constructed from line-of-sight radiances directed outward or inward, respectively. In the elevation

Figure 2-5a. Radiances and Source Functions



SOURCE FUNCTION



2

Figure 2-5b. Radiances and Source Functions

plane, the source function is approximated by an angular quadrature of tixed resolution given by the prescribed values of aspect angle. In the azimuth plane, the source function is approximated by an angular quadrature of variable resolution which increases in proportion to the prescribed number of cylindrical shells. The resolution increases from the center (where the radiation field is isotropic by symmetry) to the edge (where the radiation field is most anisotropic) such that the accuracy of the quadrature is maintained independent of radius.

Figure 2-6 shows the indexing system for the matrix elements in the azimuth plane. The indices i and k designate the spacing (x) of the line-of-sight and the radius (r) of the cylindrical shell. The index j (which does not affect the azimuth plane projection) designates the elevation angle (ϕ) of the line-of-sight relative to the cylinder axis. The i,k matrix elements form an asymmetric square array in which the ordering of the subscripts indicates the direction of the path (outward or inward).

Figure 2-7 shows the geometric relations for each matrix element in the azimuth plane. The inclination of the radiance vector relative to the radial direction is given by the complement of the angle θ_{ik} or θ_{ki} where

$$\theta_{ik} = -\theta_{ki} = \cos^{-1}\left(\frac{r_i}{r_k}\right)$$
 ; $\theta_{1k} = \pi/2$.

The line-of-sight distance s_{ijk} or s_{kji} measured from the midpoint (x-axis) to the point k is given by

Figure 2-6. Matrix Element Indexing in Azimuth Plane

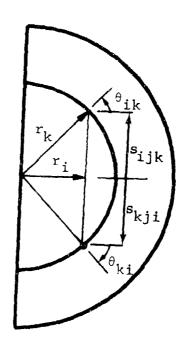


Figure 2-7. Geometrical Relations in Azimuth Plane

$$s_{ijk} = -s_{kji} = r_k \frac{\sin \theta_{ik}}{\sin \phi_j}$$

For uniform properties (ρ_k, κ_k) between points k and k+1, the transmission is given by

$$T_{ijk} = \exp\left[-\rho_k \kappa_k \left(s_{ijk+1} - s_{ijk}\right)\right]$$

With these definitions, the equation of transport between points k and k+1 in the lower quadrant and points k-1 and k in the upper quadrant may be written as

$$N_{kji} = N_{k+lji}T_{ijk} + \left[(1 - \omega_{k})N_{k}^{\circ} + \frac{\omega_{k}}{2} (S_{kji} + S_{k+lji}) \right] (1-T_{ijk})$$

$$N_{ijk} = N_{ijk-l}T_{ijk-l} + \left[(1 - \omega_{k-l})N_{k-l}^{\circ} + \frac{\omega_{k-l}}{2} (S_{kji} + S_{k-lji}) \right] (1 - T_{ijk-l})$$

$$(2-18)$$

source function S accounts for multiple scattering from all directions into the line of sight.

$$S_{ijk} = \sum_{i'=1}^{k} \sum_{j'=1}^{J} (P_{ii'kjj'}, N_{i'j'k} + P_{iki'jj'}, N_{kj'i'}) \Omega_{i'j'k} + P_{oj}N_{ok}\Omega_{o})$$

$$S_{kji} = \sum_{i'=1}^{k} \sum_{j'=1}^{J} P_{i'kij'j}N_{i'j'k} + P_{ki'ij'j}N_{kj'i'} \Omega_{i'j'k} + P_{oj}N_{ok}\Omega_{o})$$
(2-19)

where

k = number of discrete azimuth angles at radius r_k

J = total number of discrete elevation angles

P_{ii'kj'j} phase function for scattering from direction θ_{1'k}, φ_j, and its symmetric counterparts into direction θ_{ik}, φ_j

 Ω_{ijk} = solid angle for radiance vector N_{ijk} or N_{kji}

o = subscript for axial component $(\phi_0 = 0)$.

The axial component of the source function (subscript zero) is excluded from the double sum for reasons of symmetry. Symmetry with respect to z yields the following expression for exial radiance

$$N_{ok} = (1 - \omega_k) N_k^o + \omega_k S_{ck}$$

Symmetry with respect to θ yields the following expression for axial source function

$$S_{ok} = \sum_{j=1}^{J} P_{jo} \sum_{i=1}^{k} (N_{ijk} + N_{kji}) \Omega_{ijk} + P_{oo} N_{ok} \Omega_{o}$$

Substitution of S_{ok} into N_{ok} and N_{ok} into N_{ijk} and N_{kji} eliminates N_{ok} from the radiance matrix. These values are computed independently subsequent to solution of the matrix.

The solid angles are evaluated in terms of the discrete spherical coordinates θ_{ik} , ϕ_j shown in Figure 2-8. The solid angle boundaries are assumed to bisect the angular region between adjacent radiance vectors such that, for an interior element,

$$\Omega_{ijk} = -\left(\frac{\theta_{i+1,k} - \theta_{i-1,k}}{2}\right) \left[\cos\left(\frac{\phi_{j} + \phi_{j-1}}{2}\right) - \cos\left(\frac{\phi_{j} + \phi_{j+1}}{2}\right)\right] \qquad (i \neq 1,k ; j \neq J)$$

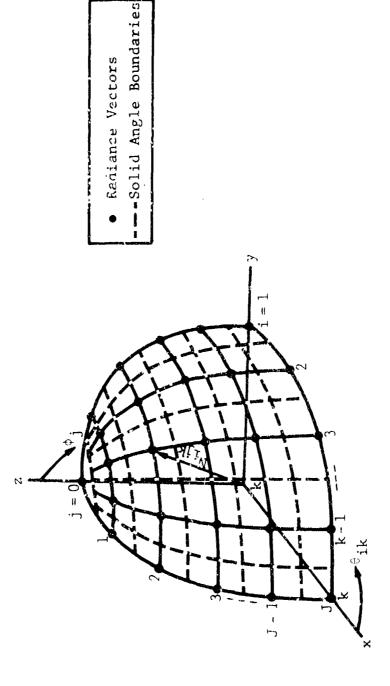


Figure 2-8. Radiance Vectors and Solid Angles

For an exterior element, the solid angle is one-half (edge) or one-fourth (corner) the value given by the above formula; i.e.,

$$\Omega_{1jk} = -\left(\frac{\theta_{2k} - \theta_{1k}}{2}\right) \left[\cos\left(\frac{\phi_{j} + \phi_{j-1}}{2}\right) - \cos\left(\frac{\phi_{j} + \phi_{j+1}}{2}\right)\right]$$

$$i = 1 : j \forall J \text{ (edge)}$$

$$\Omega_{kjk} = -\left(\frac{\theta_{kk} - \theta_{k-1,k}}{2}\right) \left[\cos\left(\frac{\phi_{j} + \phi_{j-1}}{2}\right) - \cos\left(\frac{\phi_{j} + \phi_{j+1}}{2}\right)\right]$$

$$i = k \; ; \; j \neq J \; (edge)$$

$$\Omega_{iJk} = -\left(\frac{\theta_{i+1,k} - \theta_{i-1k,k}}{2}\right) \left[\cos\left(\frac{\phi_{J} + \phi_{J-1}}{2}\right) - \cos\phi_{J}\right]$$

$$j = J \; ; \; i \neq 1, k \; (edge)$$

$$\Omega_{1,Jk} = -\left(\frac{\theta_{2k} - \theta_{1k}}{2}\right) \left[\cos\left(\frac{\phi_{J} + \phi_{J-1}}{2}\right) - \cos\phi_{J}\right]$$

$$i = 1 \; ; \; j = J \; (corner)$$

$$\Omega_{kJk} = -\left(\frac{\theta_{kk} - \theta_{k-1,k}}{2}\right) \left[\cos\left(\frac{\phi_{J} + \phi_{J}}{2}\right) - \cos\phi_{J}\right]$$

$$i = k ; j = J \text{ (corner)}$$

For the axial component, the solid angle is given by

$$\Omega_{O} = \pi \left(1 - \cos \frac{\phi_{1}}{2} \right) .$$

The phase function is evaluated in terms of the angle β formed between the given radiance vector and the line-of-sight.

This angle is readily evaluated from the vector dot product of the corresponding unit vectors $\hat{r}_{i'j'k}$ and \hat{r}_{ijk} where

 $\hat{r}_{ijk} = \hat{x} \sin \phi_{j} \cos \theta_{ik} + \hat{y} \sin \phi_{j} \sin \theta_{ik} + \hat{z} \cos \phi_{j}$ Thus,

$$\hat{\mathbf{r}}_{\mathbf{i}',\mathbf{j}',\mathbf{k}} \cdot \hat{\mathbf{r}}_{\mathbf{i}\mathbf{j}\mathbf{k}} = \sin \phi_{\mathbf{j}'} \sin \phi_{\mathbf{j}} \left[\cos \left(\theta_{\mathbf{i}',\mathbf{k}} - \theta_{\mathbf{i}\mathbf{k}} \right) \right] \\ + \cos \phi_{\mathbf{j}'} \cos \phi_{\mathbf{j}'} \\ = \cos \beta_{\mathbf{i}',\mathbf{j}',\mathbf{k}\mathbf{i}\mathbf{j}}$$

The phase function is defined to include scattering not only from the direction $\theta_{i'k}$, $\phi_{j'}$, but also from the symmetric directions π - $\theta_{i'k}$ and π - $\phi_{j'}$, such that

$$\begin{split} \mathbf{P}_{\mathtt{li'kjj'}} & & = & \mathbb{P}\Big[\beta\big(\theta_{\mathtt{i'k}}, \ \phi_{\mathtt{j'}}, \ \theta_{\mathtt{ik}}, \ \phi_{\mathtt{j}}\big)\Big] \\ & + & \mathbb{P}\Big[\beta\big(\pi - \theta_{\mathtt{i'k}}, \ \phi_{\mathtt{j'}}, \ \theta_{\mathtt{ik}}, \ \phi_{\mathtt{j}}\big)\Big] \\ & + & \mathbb{P}\Big[\beta\big(\theta_{\mathtt{i'k}}, \ \pi - \phi_{\mathtt{j'}}, \ \theta_{\mathtt{ik}}, \ \phi_{\mathtt{j}}\big)\Big] \\ & + & \mathbb{P}\Big[\beta\big(\pi - \theta_{\mathtt{i'k}}, \ \pi - \phi_{\mathtt{j'}}, \ \theta_{\mathtt{ik}}, \ \phi_{\mathtt{j}}\big)\Big] \end{split}$$

For the axial component, the phase function is given by

$$P_{jO}^{\prime} = 2[P(\phi_{j}) + P(\pi - \phi_{j})]$$

In this manner, the computed radiation field for π -steradians is made to account for scattering from 4π -steradians.

The preceding equations may be solved by either iterative or inversion methods. In the iterative approach, the equations are suitable for programming without further manipulation. At

each iteration, the radiances $N_{\rm ok}$, $N_{\rm ijk}$, and $N_{\rm kji}$ are computed based on assumed values for $S_{\rm ok}$, $S_{\rm ijk}$, and $S_{\rm kji}$ from the preceding iteration, subject to the boundary condition $N_{\rm kji}=0$ for all i,j (no external sources); the assumed values for the first iteration are taken equal to zero (i.e., single scattering). In the inversion approach, the equations are manipulated to eliminate the source functions S in favor of the radiances N which are subsequently determined by matrix inversion.

The manipulations required for the inversion method are summarized below. Elimination of $S_{\rm ok}$ in favor of $N_{\rm ok}$ yields

$$N_{ok} \left(1 - \omega_{k} P_{oo} \Omega_{o} \right) = \left(1 - \omega_{k} \right) N_{k}^{o} + \omega_{k} \sum_{i=1}^{k} \sum_{j=1}^{J} P_{jo} \left(N_{ijk} + N_{kji} \right) \Omega_{ijk}$$

Substitution of the N_{ck} into $S_{i,jk}$ yields

$$\begin{split} \mathbf{S}_{\mathbf{i}\mathbf{j}\mathbf{k}} &= \sum_{\mathbf{i}'=1}^{\mathbf{k}} \sum_{\mathbf{j}'=1}^{\mathbf{J}} \left\{ \left[\mathbf{P}_{\mathbf{i}\mathbf{i}'\mathbf{k};\mathbf{j}'} + \mathbf{x}_{\mathbf{j}\mathbf{k}} \mathbf{\omega}_{\mathbf{k}} \mathbf{P}_{\mathbf{j}o} \right] \mathbf{N}_{\mathbf{i}'\mathbf{j}'\mathbf{k}} \right. \\ &+ \left[\mathbf{P}_{\mathbf{i}\mathbf{k};\mathbf{i}',\mathbf{j};\mathbf{j}'} + \mathbf{x}_{\mathbf{j};\mathbf{k}} \mathbf{\omega}_{\mathbf{k}} \mathbf{P}_{\mathbf{j}o} \right] \mathbf{N}_{\mathbf{k};\mathbf{j}';\mathbf{i}'} \right\} \mathbf{\Omega}_{\mathbf{i}',\mathbf{j}'\mathbf{k}} \\ &+ \mathbf{x}_{\mathbf{j};\mathbf{k}} \left(1 - \mathbf{\omega}_{\mathbf{k}} \right) \mathbf{N}_{\mathbf{k}}^{o} \end{split}$$

where

$$y_{jk} = \frac{P_{jo}\Omega_{o}}{1 - \omega_{k}P_{oo}\Omega_{o}}$$

Elimination of S_{ijk} and S_{kji} in favor of N_{ijk} and N_{kji} yields a KJ(K-1) dimensional linear equation set, with KJ(K-1)/2 equations for each quadrant.

Positive Quadrant:

$$\sum_{i'=1}^{k} \sum_{j'=1}^{J} \left[\left(\delta_{i'j'ij} - \psi_{ijk-1} a_{ii'kjj'} \right) N_{i'j'k} \right.$$

$$- \psi_{ijk-1} a_{iki'jj'} N_{kj'i'} \right] - \sum_{i'=1}^{k-1} \sum_{j''=1}^{J} \left[\left(\delta_{i'j'ij} T_{ijk-1} + \psi_{ijk-1} a_{iki'jj'} N_{k-1j'i'} \right) \right]$$

$$+ \psi_{ijk-1} a_{ii'k-1jj'} N_{i'j'k-1} + \psi_{ijk-1} a_{iki'jj'} N_{k-1j'i'} \right]$$

$$+ \psi_{ijk-1} x_{jk} \left(1 - \omega_{k} \right) N_{k}^{\circ} + \left(1 - T_{ijk-1} + \psi_{ijk-1} x_{jk-1} \right) \left(1 - \omega_{k-1} \right) N_{k-1}^{\circ}$$

$$- \omega_{k-1} N_{k-1}^{\circ}$$

$$- \left(\frac{i=1,k}{j=1,J} \right)$$

$$- \left(\frac{i=1,k}{j=1,J} \right)$$

$$- \left(\frac{i=1,k}{j=1,J} \right)$$

$$- \left(\frac{i=1,k}{j=1,J} \right)$$

Negative Quadrant:

$$\sum_{i'=1}^{k} \sum_{j'=1}^{J} \left[\left(\delta_{i'j'ij} - \psi_{ijk} a_{ki'ij'j} \right) N_{kj'i'} - \psi_{ijk} a_{i'kij'j} N_{i'j'k} \right] - \sum_{i'=1}^{k+1} \sum_{j'=1}^{J} \left[\left(\delta_{i'j'ij} T_{ijk} + \psi_{ijk} a_{k+1i'ij'j} \right) N_{k+1j'i'} + \psi_{ijk} a_{i'k+1ij'j} N_{i'j'k+1} \right]$$

$$= \psi_{ijk} x_{jk+1} \left(1 - \omega_{k+1} \right) N_{k+1}^{o} + \left(1 - T_{ijk} + \psi_{ijk} x_{jk} \right) \left(1 - \omega_{k} \right) N_{k}^{o}$$

$$\left\{ \begin{cases} i = 1, k \\ j = 1, J \\ k = 1, J \end{cases} \right\}$$

$$2-49$$

where

$$\psi_{ijk} \equiv \frac{\omega_{k}}{2} \left(1 - T_{ijk} \right)$$

$$a_{ii'kjj'} \equiv \left(P_{ii'kjj'} + x_{jk} \omega_{k} P_{jo} \right) \Omega_{i'j'k}$$

The above equations are subject to KJ boundary conditions

$$N_{Kji} = 0$$
 {i = 1,K; j = 1,J}

The above relations suffice to determine the entire K²J dimensional set of radiances for the cylinder. Repeated application at successive axial stations yields the complete radiation field for the plume.

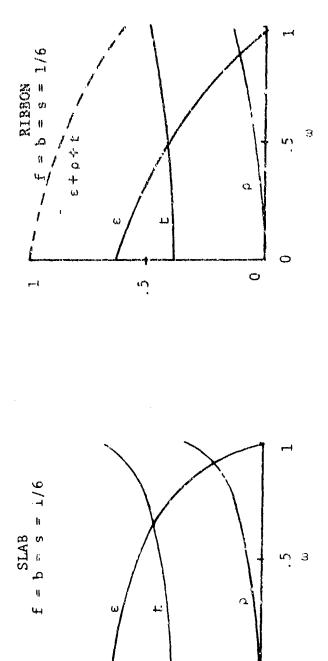
2.6 Computed Results

Extensive parametric calculations were undertaken in order to validate and compare the six-flux and N-flux scattering approaches. The calculations included comparisons with the predictions of other codes: the NASA finite-difference code for plane-parallel geometries and the GAC Monte-Carlo code for arbitrary geometries. The calculations were performed for specified values of the monochromatic scattering parameters (optical depth, scattering albedo, phase function) rather than for a specified spectral region or particle size/density distribution. This procedure enabled the scattering codes to be exercised over the full range of optical environments without the attendant complexities of a complete system

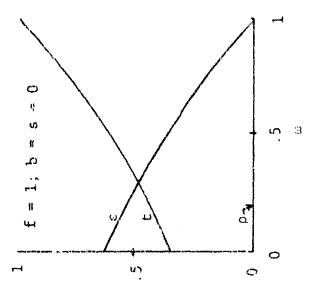
simulation. The evaluation of the optical parameters in terms of the physical properties of a specific gas/particle medium is considered elsewhere in this report.

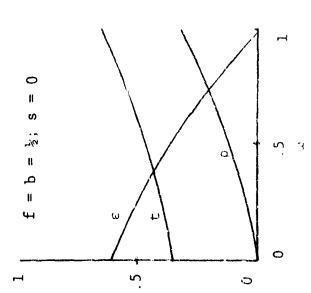
2.6.1 Planar media. The relative simplicity and general availability of solutions for planar media make such solutions ideal for the preliminary test of physical models and parametric effects. This section summarizes results obtained for uniform one- and two-dimensional planar media using an explicit six-flux model (Method I; Figures 2-9 through 2-12) and a standard finite-element program (NASA code; Figures 2-13 through 2-15) modified to include thermal emission.

Figure 2-9 shows the transmissivity, reflectivity, and emissivity of a uniform planar medium of unit optical depth at normal aspect for various phase functions (isotropic, fore and aft, forward only). In addition, isotropic scattering solutions are compared for one- and two-dimensional planar media (slab vs. ribbon) in order to bound the possible range of values for cylindrical media. In all cases, scattering is observed to reduce the emitted component and to increase the transmitted-plus-reflected component of the radiation signature. Comparisons among the slab solution for different phase functions show the effects of backscattering (b>0) and side-scattering (s>0) on the reflected and emitted components. Comparisons between the



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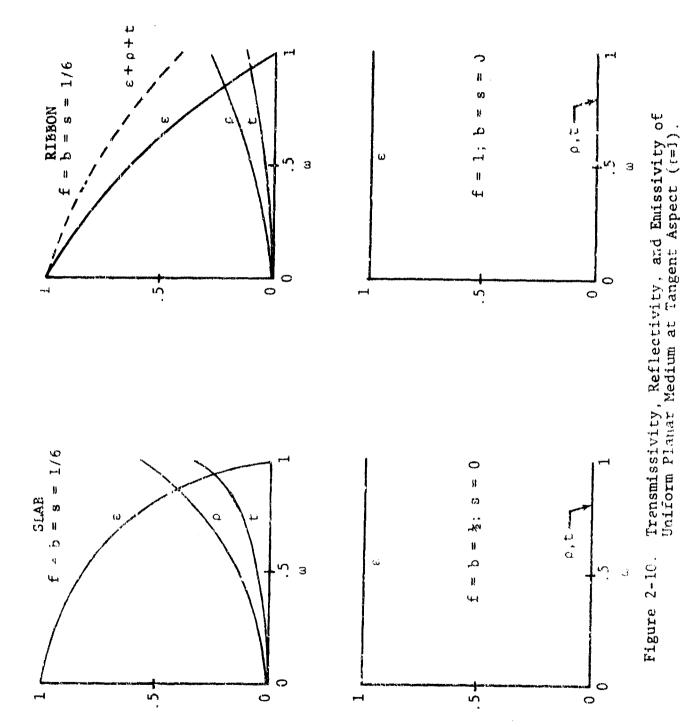
Transmissivity, Reflectivity, and Emissivity of Uniform Planar Medium at Normal Aspect $(\tau=1)$. Figure 2-0.

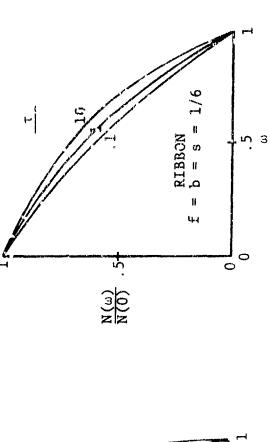
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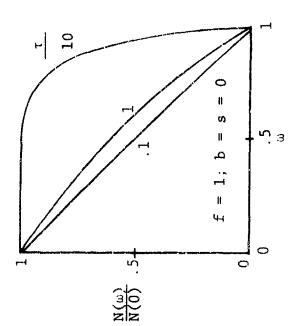
slab and ribbon solutions for isotropic scattering (f=b=s) reveal essential differences between conservative ($\varepsilon + \tau + \rho = 1$) and non-conservative ($\varepsilon + \tau + \rho < 1$) radiation fields which would not be predicted by a two-flux model (s=0).

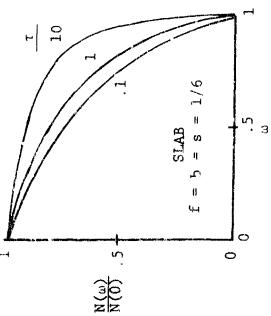
Figure 2-10 shows the transmissivity, reflectivity, and emissivity of a uniform slab at tangent aspect for the same range of parameters as in Figure 2-9. In this case, the incident radiation is assumed to be perpendicular to the surface while the observed radiation is parallel. Thus, the transmitted radiation component undergoes a change in direction from the incident side to the emergent side which accounts for $\tau + 0$ as $\omega + 0$. The emitted component parallel to the slab is black (infinite optical depth) only in the limit $\omega + 0$; for finite ω , the emissivity is reduced due to losses through the slab boundaries. Note that this coupling between the normal and tangential radiance components does not occur for the two-flux models.

Figure 2-11 shows the radiance of a uniform emitting planar medium at normal aspect over a range of optical depths in the absence of external radiation sources. The values are normalized by the non-scattering solution for $\omega=0$ and are therefore equivalent to normalized emissivities. The range of scattering parameters and geometries is the same as in Figures 2-9 and 2-10. In all cases, increased scattering (ω) is observed to reduce the observed radiance for constant τ due to equal but









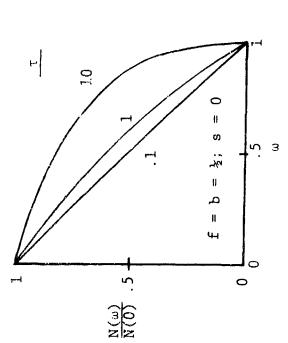


Figure 2-11. Normal Radiance versus Albedo for Uniform Planar Medium

opposite variations in the optical depths for absorption and scattering. The effect of side-scattering is observed to increase the radiance from the slab relative to the ribbon and from the six-flux relative to the two-flux (f=b in both cases) but to decrease the radiance at large optical depths relative to the forward-scattering case (f=1).

Although not shown, the normalized axial radiance component $[N_Z(\omega)/N_Z(0)]$ exhibited close similarity to the normalized radial component $[N_{\mathbf{r}}(\omega)/N_{\mathbf{r}}(0)]$ for isotropic scattering but maintained constant value $[N_Z(\omega)/N_Z(0)=1]$ for two-stream scattering. The normalizing values $N_{\mathbf{r}}(0)$ and $N_Z(0)$ are simply the given black body function (N^0) times the apparent emissivity $(1-e^{-T_{\mathbf{r}}}$ or unity, respectively). Thus, the six-stream approximation accounts for the redistribution of radiance (from axial to radial) due to scattering whereas the two-stream approximation does not.

Figure 2-12 shows analogous results to Figure 2-11 except that the optical depth parameter is based on the absorption component rather than the total extinction (absorption + scattering). In this case, the emitted radiation remains constant as ω increases rather than decreasing, as before, in proportion to $(1-\omega)$. As a result, the curves lie above those in Figure 2-11 by an amount which increases with increasing ω . In the case of forward scattering, the radiance remains finite in the limit $\omega \to 1$.

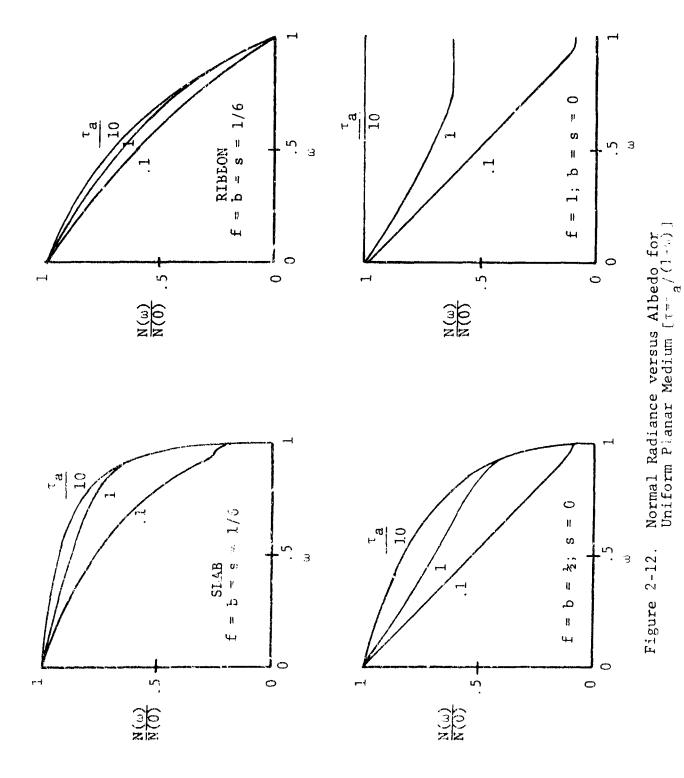
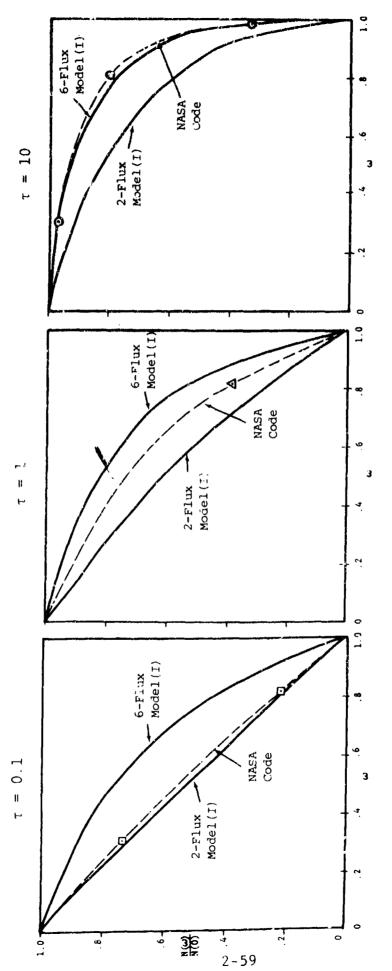
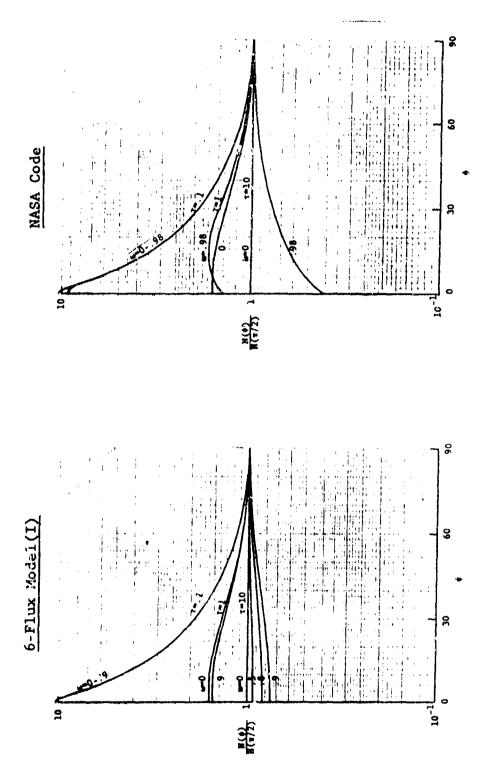


Figure 2-13 shows a comparison between the predictions of the two-flux and six-flux models and the NASA finite-element code for the normalized radiance at normal aspect from an isotropically-scattering uniform slab of specified optical depth. For small optical depth $(\tau=.1)$, the NASA code is in close agreement with the two-flux model $(f=b=\frac{1}{2})$. For large optical depth (T010), the NASA code is in close agreement with the six-flux model (f=b=s=1/6). For intermediate optical depth (based on a single NASA code calculation), the NASA code prediction is intermediate between the other two. Taking the NASA code as the standard, this comparison provides physical insight into the sources of error which can result from coarse angular quadrature. In short, the two- and six-flux quadratures, respectively, underpredict and over-predict the effective emissivity of the medium. However, from the comparative results of Figures 2-9 to 2-13 for slabs and ribbons, the error in the six-flux approximation at small optical depths should be appreciably reduced for cylindrical geometries (i.e. plumes).

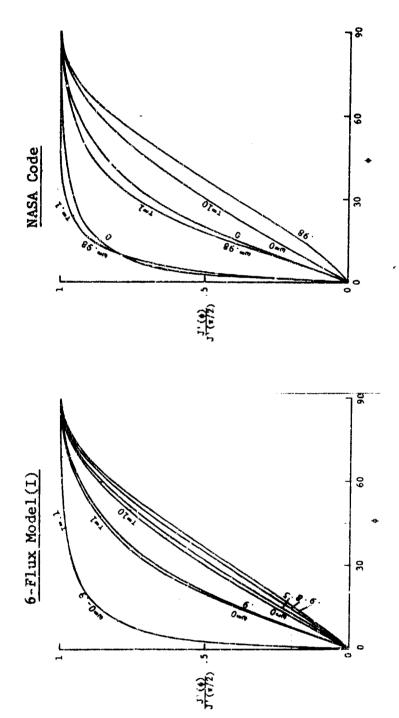
Figures 2-14 and 2-15 show further comparisons between predictions of the six-flux model and the NASA code for an isotropically scattering uniform slab. These figures show the variation with aspect angle of the radiance and radiant intensity (radiance times projected area) normalized to their respective values at normal aspect. The predictions of the six-flux model (Method I) represent interpolations between the values actually computed for normal and tangent aspect and, therefore,



Normal Radiance versus Albedo for Uniform Planar Medium (Isotropic Scattering). Figure 2-13.



Radiance versus Aspect for Uniform Planar Medium (Isctropic Scattering) Figure 2-14.



Radiant Intensity versus Aspect for Uniform Planar Medium (Isotropic Scattering) Figure 2-15.

have limited quantitative accuracy between those limits. However, the predictions and comparisons are useful to illustrate In particular, both models predict the same qualitative variation of radiance with aspect (Figure 2-14) which is strongly dependent on optical depth. For small optical depths and low scattering albedos, the radiance increases toward tangent aspect due to the increased length of the optical path. However, for large optical depths and high scattering albedos, the radiance decreases toward tangent aspect due to side-scattering losses not predicted by two-flux models. The aspect dependence of the radiant intensity (Figure 2-15) includes the variation in projected area (proportional to sin \$\phi\$) and, consequently, shows a monotonic decrease from normal to tangent aspect. In both Figures, the six-flux model predicts lower levels of radiation for higher scattering albedos, with greater sensitivity for larger optical depths. The opposite behavior of the NASA predictions for low optical depth does not lead to larger absolute differences.

2.6.2 Cylindrical media. Compared to the vast literature on multiple-scattering solutions for planar media, relatively little work has been published for cylindrical media. Accounting for this essential geometric effect was a central feature of the SIRRM model development. Hence, the logical first test of a new solution for cylindrical media is a comparison with previous solutions for planar media.

Figure 2-16 shows the normal radiance of a uniform infinite cylinder (centered line-of-sight) and of a uniform slab of equal optical depth. The predictions are based on the six-flux model (Method I or II) and on the GAC Monte Carlo code. Except in the non-scattering limit (ω =0), the radiances are observed to be lower for the cylinder than for the slab due to side-scattering losses (not predicted by two-flux models). The six-flux and Monte Carlo results are shown to be in essential agreement for both isotropic and forward scattering phase functions. The forward scattering result is only slightly greater than the isotropic scattering result due to the assumption of uniform properties; larger sensitivities would be expected for the non-uniform conditions of actual plumes.

Figure 2-17 shows the normal radiance for isotropic scattering of a uniform infinite cylinder (centered line-of-sight) and of a uniform slab and ribbon of equal optical depth. The predictions are based on the 6-flux model (Method I or II) with comparative calculations from the N-flux model and the Monte Carlo code. The variation of radiance with optical depth (upper left plot) is qualitatively consistent with previous predictions for planar media (Figure 2-11). Specific comparisons for equal optical depths (remaining plots) show a uniform progression of the cylinder radiances from the ribbon values at small optical depths to the slab values at large optical depths. This behavior is the result of side-scattering losses from the cylinder which increase with decreasing optical depth. These losses

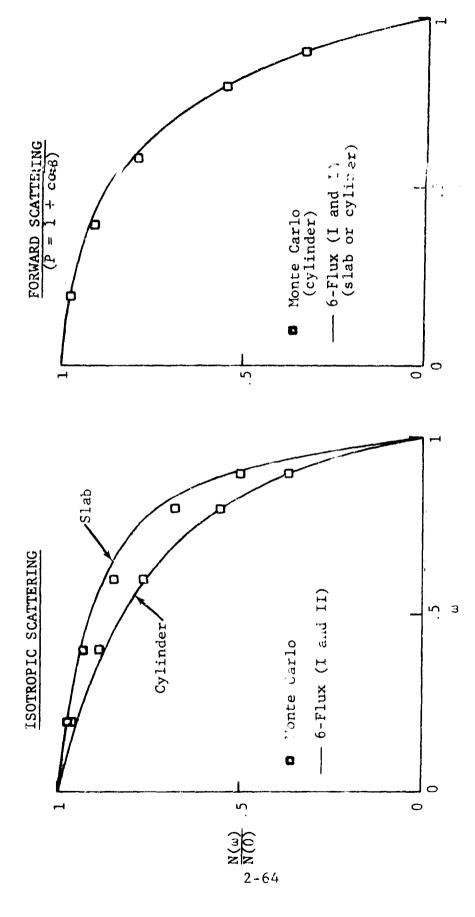
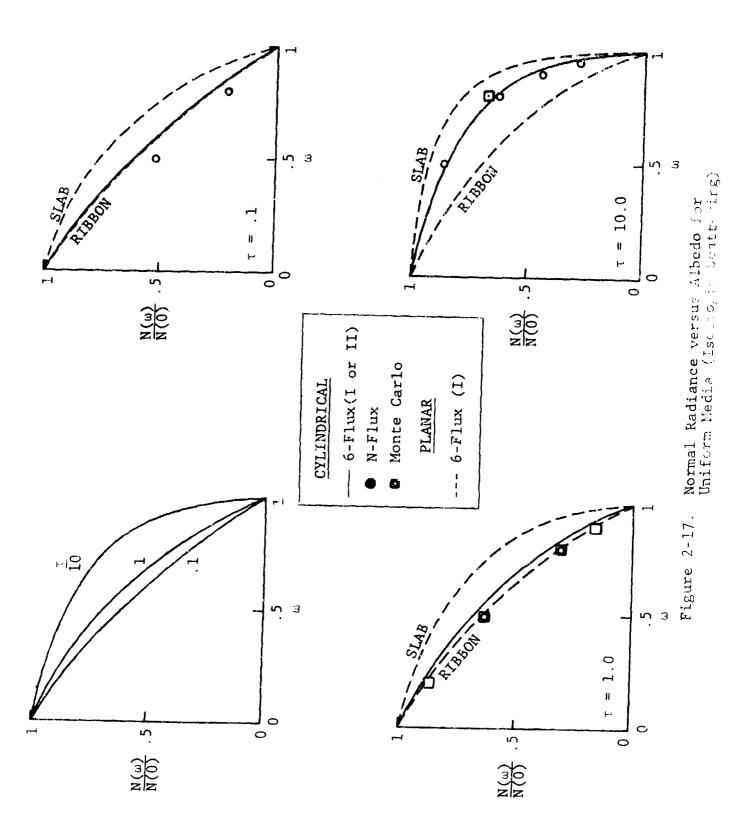


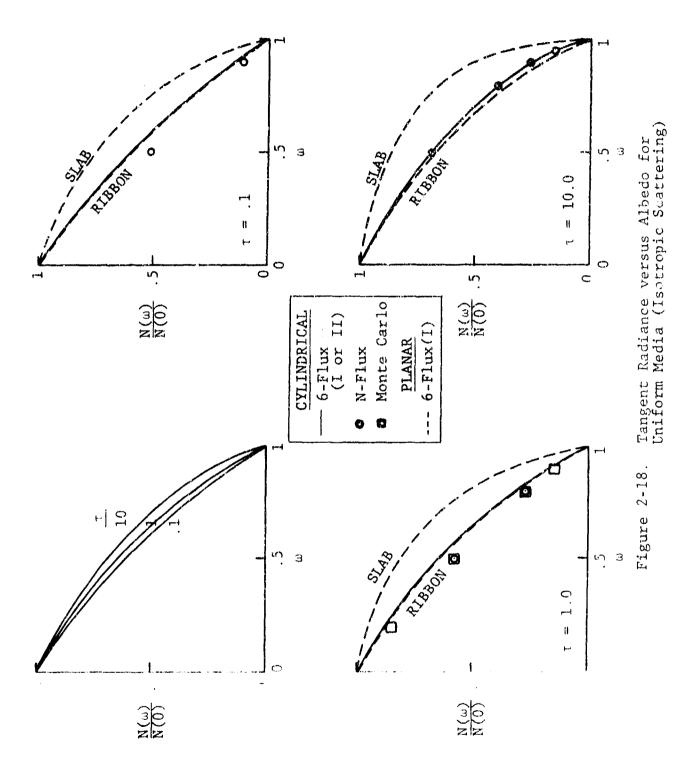
Figure 2-16. Normal Radiance versus Albedo for Uniform Media (1.4.0)

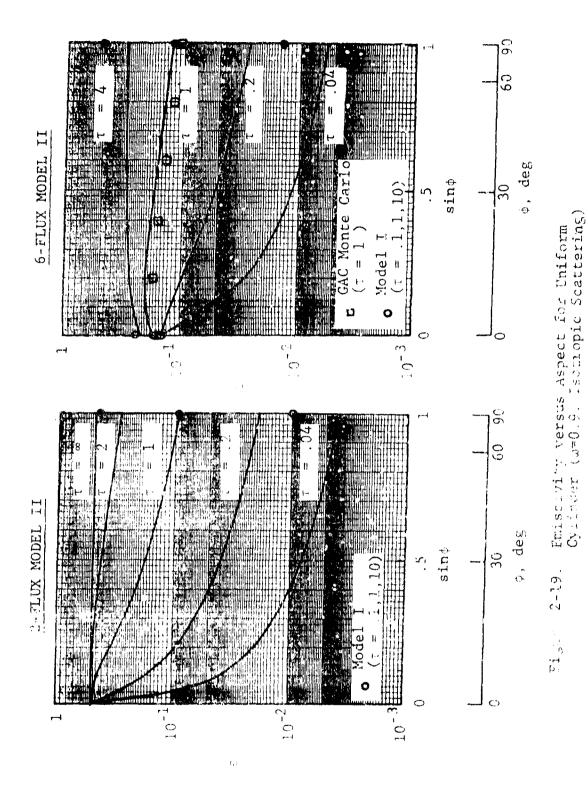


result in improved predictive accuracy for the six-flux model at small optical depths. Comparison of the six-flux model predictions with the more exact predictions of the N-flux model and Monte Carlo codes reveals a residual over-prediction for small optical depths but substantial improvement relative to similar predictions for a slab (compare Figure 2-13). Comparison of the N-flux and Monte Carlo predictions for unit optical depth revals excellent agreement. The comparison is close but less exact at an optical depth of ten.

Figure 2-18 shows the tangent radiances (i.e., aspect = 0 or 180 degrees) for the same conditions as the normal radiances in Figure 2-17. As before, the curves are shown normalized by their values for a non-scattering medium, i.e., the local Planck function $\left[N_Z(\omega=0)=N^0\right]$. The curves reveal the same qualitative trends as the normal radiances in Figure 2-17 except for a reduced sensitivity to optical depth due to cylindrical curvature. As a result, the tangent radiances remain close to the ribbon values for all optical depths. This effect is confirmed by the more exact predictions of the N-flux model and Monte Carlo code.

Figure 2-19 shows the aspect dependence of the emissivity of a uniform infinite cylinder (centered line-of-sight) as predicted for a given albedo by two-flux and six-flux model approximations (Method II). The curves are qualitatively consistent (but quantitatively different) for small optical depths near normal aspect. However, for large optical depths near tangent aspect, the two approximations predict opposite trends due to





the effect of side-scattering losses near grazing incidence. This behavior is consistent with previous predictions of radiance vs. aspect for a planar slab (Figure 2-14) and may have significant negative impact on the detectability of plumes from nose aspect. Comparison of the predictions of Method I and II for normal and tangent aspect reveals good agreement for both phase function approximations at comparable optical depths. Comparison of the predictions of Method II and the Monte Carlo code for isotropic scattering at unit optical depth reveals agreement within 20 percent for all aspect angles.

Figure 2-20 shows a comparison between the six-flux (Method I) and N-flux models for the normalized radiance from an isotropically scattering uniform infinite cylinder (centered line-of-sight) over the full range of albedo and optical depth. These curves are equivalent to normalized emissivities from Figure 2-19. As in Figure 2-14 (slab), the present six-flux predictions represent interpolations between the values actually computed for normal and tangent aspect and, therefore, have limited quantitative accuracy between those limits. In particular, the Method I interpolation formulae are observed to under-predict relative to the N-flux model but to yield valid qualitative trends. Both models yield comparable predictions at normal and tangent aspects. Comparison of N-flux and Monte Carlo predictions are in excellent agreement and appear to confirm the 6-flux Method II predictions in Figure 2 - 19.

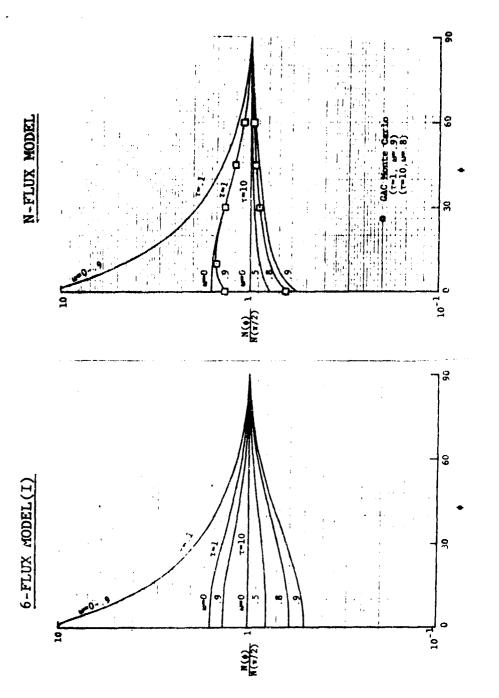


Figure 2-20. Centerline Radiance versus Aspect for Uniform Cylinder (Isotropic Scattering)

Figure 2-21 shows the N-flux model prediction for the variation of normal radiance with distance from the cylinder axis for the same conditions as Figure 2-20. Integration under these curves would yield the normalized station radiation at side aspect. The effect of scattering is to reduce the radiance from the off-axis regions, with greater sensitivity for larger optical depths. The assumption of the six-flux model I that the radiance profile is independent of albedo is, therefore, seen to be in error for large optical depth. This assumption is not made by six-flux model II.

Figure 2-22 shows a comparison between the six-flux (Method I) and N-flux model predictions for station radiation as a function of aspect angle for the same conditions as Figure 2-20. Both models predict the same qualitative effect of increasing albedo, namely, decreased radiation relative to side aspect with greater sensitivity for large optical depth. This behavior is consistent with previous 6-flux model predictions for a planar slab (compare Figure 2-15) except that the sensitivity to albedo is greater in the present case. However, the six-flux model (Method I) interpolation formulae appear to over-predict the sensitivity as predicted by the N-flux model and as confirmed by additional predictions with the GAC Monte Carlo code. The latter two predictions are shown to be in excellent agreement.

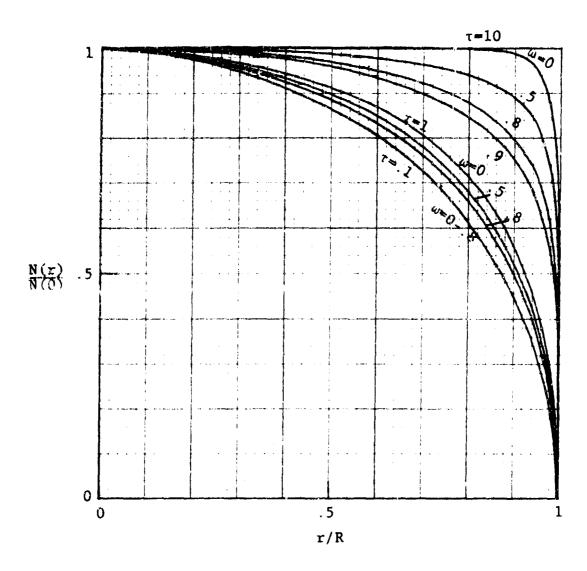
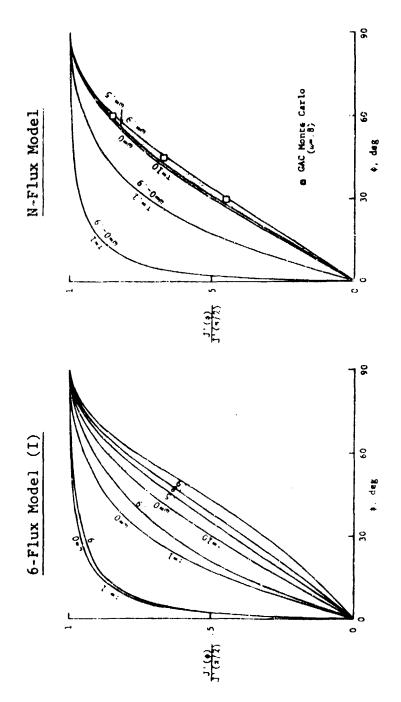
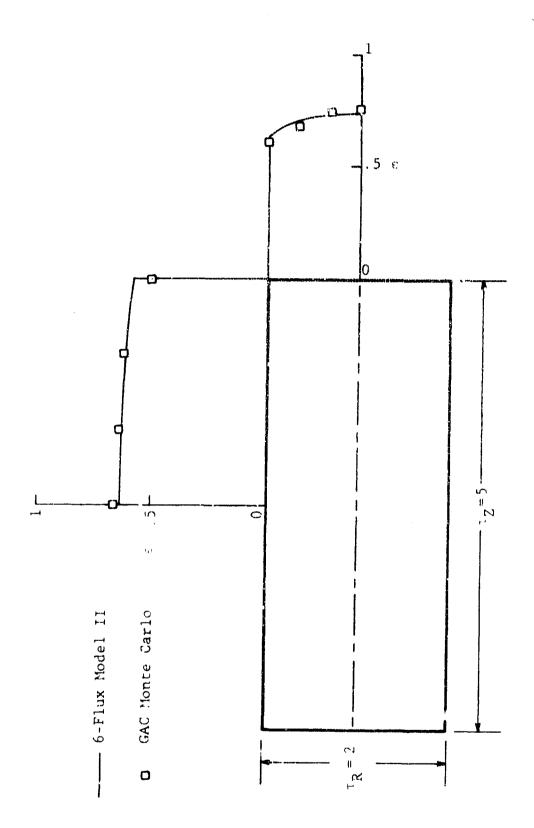


Figure 2-21. Normal Radiance versus Distance from Centerline (Isotropic Scattering)



The preceding results for infinite cylindrical media represent valid solutions for rocket plumes provided that the variation in axial radiance is negligible over a distance of one optical depth (i.e., $dN_2/d\tau_2 << N_2$). This criterion is satisfied throughout most of the plume mixing region but may not be satisfied in local regions of the plume inviscid core when strong shocks are present. In order to assess the impact of shock-induced gradients on the coupled gas/particle radiation field, the six-flux model (Method II) was extended to finite cylindrical media with arbitrary end-wall boundary conditions. Numerical experiments were recently initiated to assess the impact of aallong external radiation source parallel to the axis on the radiation scattered to side aspect. A key objective of this assessment is the determination of the axial region of influence beyond which the infinte cylinder model becomes a valid approx mor ico

Figure 2-23 shows the result of a preliminary test calculation with the extended six-flux model. In this case, the maintain is assumed uniform with no external radiation seem and that the computed radiation field is symmetric about the cylinder mid-point. The radial and axial emissivities are observed to decrease from the mid-points toward the corners as a result of three-dimensional radiation-coupling effects. For the given optical depths, the mid-point radial emissivity is nearly equal to the emissivity of an infinite cylinder; i.e., the mid-point



Axial and Radial Emissivities for Finite Cylinder (Isotropic Scattering; $\omega=.5$) Figure 2-23.

values are nearly independent of end-wall boundary conditions.

Thus, the axial region of influence is contained within the finite cylinder considered. Comparative Monte Carlo predictions appear to validate the six-flux model result.

As a last test case, a condition of practical significance was chosen. It involves an occulted emitting core and a non-emitting scattering envelope, simulating a nose aspect signature (see Figure 2-24). This is a stringent test case because the nose aspect signature is due entirely to scattering and because the core is a localized source leading to a highly anisotropic radiation field. A comparison between the forward scattered radiance predicted by the 6-flux model (Method II) and the Monte Carlo method showed agreement within 20%.

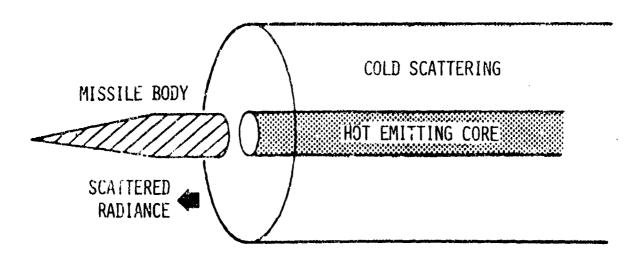


Figure 2-24. Core/Envelope Structure for Nose Aspect Test Case.

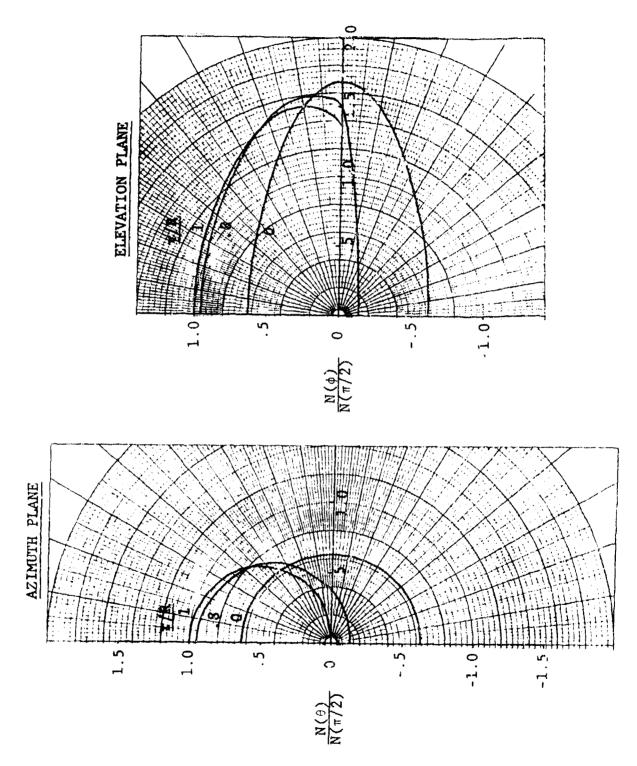
2.7 Summary of Models and Results

The results discussed in Section 2.6 include calculations for planar and cylindrical media using 6-flux, finite-difference, and Monte Carlo solution techniques. The calculations include the full range of optical parameters [albedo (0 to 1), optical depth (.1 to 10), phase function (isotropic to delta function)] and observables (transmissivity, reflectivity, emissivity; radiance and station radiation versus position and aspect). The results demonstrate the importance of three-dimensional multiple-scattering effects and the relative ability of different physical models to provide an accurate simulation.

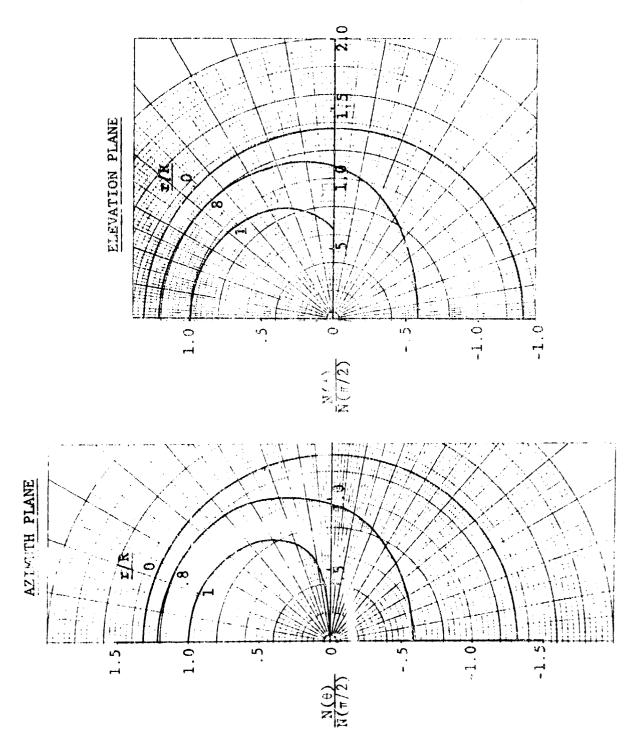
Three new models were developed and verified in this phase of the effort. The first was an explicit six-flux model which is exact (for six-flux phase functions) for planar media and approximate for cylindrical media. The second was an implicit six-flux model which is exact (for six-flux phase functions) for cylindrical media. The third was an implicit multiple-flux ("N-flux") model which is exact (for arbitrary phase functions) for cylindrical media. The six-flux models provide the minimum angular resolution necessary to account for three-dimensional scattering effects. As such, they constitute valid engineering approximations which are highly efficient for production calculations. The N-flux model provides complete flexibility in the selection of angular resolution (elevation and azimuth). As a result, alternate levels of resolution can be tested and compared with respect to accuracy and efficiency.

The six-flux models are distinguished principally by the degree of rigor in treating side-scattering effects. That is, Method I approximates these effects via simple closure relations whereas Method II computes these effects via exact coupling equations. Both methods solve initially for the normal and tangential radiances and subsequently for the radiance at arbitrary aspect (if required). The fundamental difference is associated with the initial solution step.

To the authors' knowledge, the N-flux model represents the first application of general finite-element techniques to emitting cylindrical media. Previous application of this technique was to non-emitting planar media (Barkstrom, 1976). information content of the calculation is considerable. For each radial station, the model predicts the complete threedimensional radiance distribution in spherical coordinates. Typical results for two uniform isotropically scattering cylinders of different optical depth are illustrated in Figure 2-25. The figure shows polar plots in azimuth and elevation of the local radiance distribution at a series of radial stations from the dester of each cylinder to the edge. In each case, the radian is are normalized by the edge value for a centered line-of-sight at side aspect $(\Theta = \phi = \pi/2)$. In the azimuth plane, the radiance distribution is observed to be isotropic at the center and to evolve toward a forward lobe at the edge. In the elevation plane, the radiance distribution is roughly eliptic at the center.



Radiance Distributions in Azimuth and Elevation (N-Flux Model; Isotropic Scattering; $\tau=1.\%$; $\omega=0.8$) Figure 2-25a.



. Adiance Distributions in Azimuth and Elevation (-Flux Model: Tectropic Scartering: τ =10.; w=0.8) Figure 2-25b.

(approaching isotropic for large optical depths) and evolves toward a composite forward/side lobe at the edge. The radiances are observed to increase or decrease from the center to the edge for small or large optical depths, respectively, with the maximum rate of change occuring near the surface. This result may be compared to a non-scattering medium in which the radiance always increases monotonically outward.

The N-flux model provides a powerful tool for the analysis of plume scattering problems of arbitrary complexity. The accurac, of the prediction is limited only by the computational resources of the user. Because of the significant cost implications of repetitive calculations for multiple spectral points, an important aspect of the continuing effort is the optimization of the mathematical procedures and the grid selection (spatial and angular). Mathematical optimization includes the selection of iterative or inversion procedures for small or large scattering optical depth $(\omega\tau)$ and the use of a different (finer) radial grid for evaluation of the emission/extinction terms. Grid optimization involves the correlation of accuracy and efficiency with both the number of elements (radius and aspect) and the positions of their boundaries.

A rough preliminary indication of the degree of resolution required for a given accuracy is shown in Figure 2-26. This figure is representative of the limited computational experience

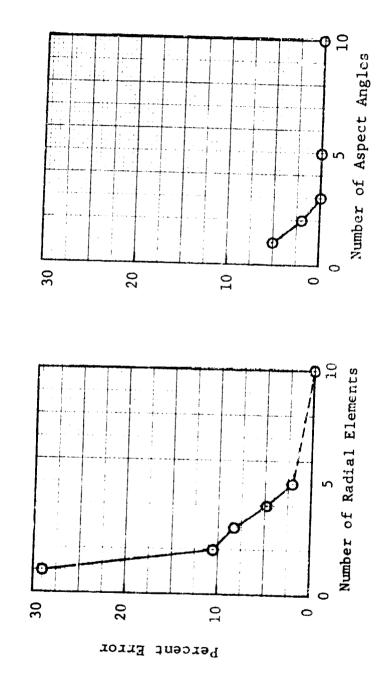


Figure 2-26. Accuracy versus Resolution for N-Flux Model (τ =10; ω =.8)

acquired to date but does not necessarily possess broader quantitative applicability. The results shown are for an optically thick uniform cylinder with isotropic scattering (τ =10, ω =.8). No particular effort was made to optimize the grid selection for a given number of points. The accuracy is observed to be much more sensitive to the number of radial elements than to the number of aspect angles.

By far the largest portion of the computing time is in the calculation of the source functions. As a result, CPU time is roughly proportional to the square of the number of aspect angles $(\phi_j\phi_j)$ and to the cube of the number of radial segments (r_k,θ_i,θ_i) . Thus, doubling both the number of aspect angles and the number of radial segments results in a computing time increase of about a factor of 30. Optimization of the selection of these numbers is, therefore, of considerable importance for production calculations.

Figure 2-27 summarizes the parametric and comparative calculations performed to date as reported in Sections 2.6 and 2.7. This summary illustrates the interrelationships among the various techniques and the scope of physical problems considered. The emphasis in this phase of the study was on model development and verification. Additional work is required in the areas of model optimization and application. In particular, final quantitative assessment of the relative accuracy and efficiency of the different calculation methods in different physical situations must necessarily await further model refinements and

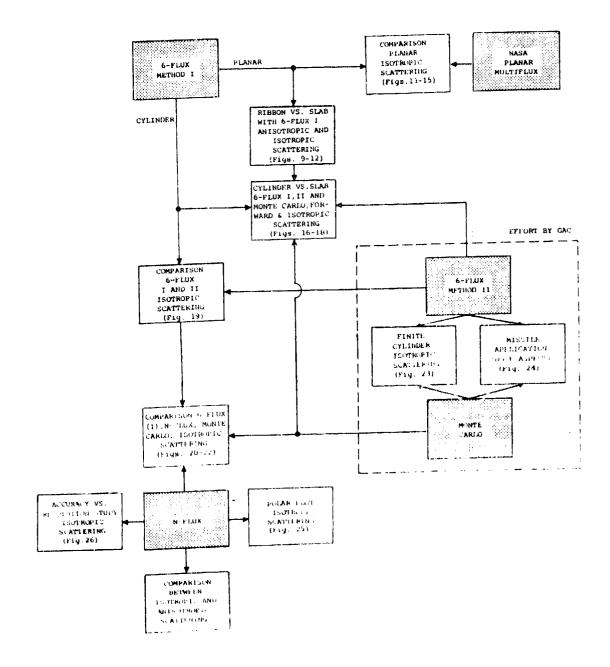


Figure 2-27. Overview of Parametric and Comparative Calculations Presented in Subsections 2.6 and 2.7.

additional parametric predictions. These efforts are currently underway. However, based on the parametric and comparative calculations performed to date, certain general conclusions can be drawn:

- 1) The six-flux angular quadrature (Method I or II) yields quantitatively accurate predictions for normal and tangent aspect in the cases of intermediate-to-large optical depth or anisotropic scattering. The standards for comparison are the NASA finite-element code for slabs and the N-Flux and Monte-Carlo codes for cylinders.
- The six-flux angular quadrature (Method I or II) tends to overpredict the normal and tangent radiation for isotropic scattering at small optical depths. The error is greatest for slabs and least for ribbons due to the corrective effect of side-scattering losses. For cylindrical plume geometries, the error appears to be acceptable (less than 20 percent).
- 3) The aspect dependence predicted by the six-flux methods is based on prior solution for normal and tangent aspect. For intermediate aspect angles, simple interpolation (Method I) gave the correct qualitative behavior, but overestimated the sensitivity to albedo by comparison to the N-flux model. Line-of-sight integration (Method II) gave the correct quantitative behavior by comparison to limited Monte Carlo calculations for intermediate optical depth.
- 4) The N-Flux and Monte Carlo calculations demonstrated uniformly excellent agreement for all cases investigated.

BAND MODEL APPROACHES

3.1 Introduction

The equation of radiative transfer, which is rigorous on a monochromatic basis, is only approximate when applied over a finite spectral interval. The spectral averaging procedures required for the calculation of radiant transfer in a multiple scattering medium must be considered.

The integral form of the equation of transfer for a finite spectral interval $\Delta\lambda$ may be written

$$\vec{N}_{\lambda}(\vec{s}) = \int_{s_{0}}^{s} \left[N_{\lambda}^{o}(s') \left(\frac{dT_{\lambda}(s',s)}{ds'} - \omega_{\lambda} \frac{dT_{\lambda}(s',s)}{ds'} \right) + \frac{1}{4\pi} \int_{4\pi}^{s} P_{\lambda}(\vec{s}'', \vec{s}') \omega_{\lambda} N_{\lambda}(\vec{s}'') \frac{dT_{\lambda}(s',s)}{ds'} du(\vec{s}'') \right] \cdot ds'$$

where a spectrally averaged quantity $\bar{\mathbf{x}}$ is defined as

$$\overline{\mathbf{x}}_{\lambda} \equiv \frac{1}{\Delta \lambda} \int_{\Delta \lambda} \mathbf{x}_{\lambda} d\lambda$$

and where the relatively small spectral variations of N_{λ}^{o} and P_{λ} within $\Delta\lambda$ are neglected. The integrand contains three spectrally-averaged terms involving the transmittance derivative and its

product with the albedo and radiance. Of these terms, only the transmittance derivative is amenable to simple averaging through superposition of the individual species averages. Evaluation of the product terms requires additional approximations which are investigated in this section. The goal of this investigation is to develop as comprehensive a band model theory of combined scattering and gaseous absorption as possible and to determine the functional dependence of the mean transmissivity and emissivity on the gaseous band model and particle scattering parameters as a guide to their most reasonable combination in the more general three-dimensional numerical solutions of the radiative transfer equation.

In developing a combined absorption-scattering band model formalism, the approach has been used which is analogous to that of the classical gaseous band-model theory, namely, from a monochromatic solution of the radiative transfer equation, to perform a spectral integration over a defined gaseous absorption band, making approximations as required to obtain a closed form solution (and studying the limitations on the accuracy of the solution imposed by the required approximations).

3.2. Band Model Formulation For One-Dimensional Scattering

A starting point has been selected to be the "one-dimensional" (i.e., two-flux) scattering formalism of Vanderbilt and Slack (1976), which has the advantage of having been developed to the state in which various closed-form solutions exist which are explicit in certain limiting cases.

It is important to note that the two-flux and six-flux approaches to the problem of radiation from a plane-parallel slab yield similar functional forms, differing only in the definitions of the effective backward and forward scattering cross sections. Thus, the so-called "one-dimensional" scattering solution is capable of reproducing the effects of three-dimensional scattering in this particular case.

It is noted also that the condition of completely anisotropic scattering which is sometimes evoked as justification for application of the one-dimensional formalism is a sufficient condition, but by no means a necessary condition for its applicability.

Because of its relative mathematical simplicity (in addition to the fact that the previously mentioned sufficient condition for its validity, i.e., the assumption of complete anisotropy, is very closely approached in many real scattering problems), this formalism will always be of great value in a large number of practical situations.

We start with one approximate solution of the two-flux scattering equation, which is exact in a limiting case and which provides an iterative solution under more general conditions (Vanderbilt and Slack, 1976):

$$I^{-}(0) = \chi_{0}^{\frac{1}{2}} \frac{(\chi_{m}^{\frac{1}{2}}+1)e^{W_{m}}\int_{0}^{W_{m}} \frac{2B}{\chi_{2}^{\frac{1}{2}}} e^{-W'}dW' - (\chi_{m}^{\frac{1}{2}}-1)e^{-W_{m}}\int_{0}^{W_{m}} \frac{2B}{\chi_{2}^{\frac{1}{2}}} e^{W'}dW'}{(\chi_{m}^{\frac{1}{2}}+1)(\chi_{0}^{\frac{1}{2}}+1)e^{W_{m}} - (\chi_{m}^{\frac{1}{2}}-1)(\chi_{0}^{\frac{1}{2}}-1)e^{-W_{m}}} (3-1)}$$

where

$$W = \int_0^s N\sigma_a \chi^{\frac{1}{2}} ds' \qquad (3-2)$$

and

$$\chi^{\frac{1}{2}} = \left(1 + \frac{2\sigma_{sb}}{\sigma_{a}}\right)^{\frac{1}{2}} = \left(1 + \frac{2\beta\omega}{1-\omega}\right)^{\frac{1}{2}} \tag{3-3}$$

In these equations, N is the particle density, σ the absorption a cross section and σ_{sb} the backscattering cross section, and β is the ratio of backscattering to total scattering. For simplicity, the absorption and scattering cross sections are referenced to a single particle density; however, for mixtures of particles or particles plus gases, the expressions are readily generalized.

For the case of constant χ , an exact solution is provided for an otherwise inhomogeneous path, i.e.,

$$I^{-}(0) = \frac{\frac{2}{\chi^{\frac{1}{2}}+1} \left[\int_{0}^{W_{m}} Be^{-W'} dw' - \frac{\chi^{\frac{1}{2}}-1}{\chi^{\frac{1}{2}}+1} e^{-2W_{m}} \int_{0}^{W_{m}} Be^{W'} dW' \right]}{1 - \left(\frac{\chi^{\frac{1}{2}}-1}{\chi^{\frac{1}{2}}+1} \right) e^{-2W_{m}}}$$
(3-4)

where
$$W_{\rm m} = \chi^{\frac{1}{2}} \tau_{\rm g}$$
 (3-5)

The absorption optical depth is denoted by τ_a , which may represent the sum of contributions from a number of gaseous and particulate species.

For the isothermal case, direct integration provides (on division by the Planck blackbody function) the known expression for emissivity:

$$\varepsilon = \frac{1^{-}(0)}{B} = \frac{\frac{2}{\chi^{\frac{1}{2}}+1} \left[\left(1 - e^{-\chi^{\frac{1}{2}}\tau} a \right) - \frac{\chi^{\frac{1}{2}}-1}{\chi^{\frac{1}{2}}+1} \left(e^{-\chi^{\frac{1}{2}}\tau} a - e^{-2\chi^{\frac{1}{2}}\tau} a \right) \right]}{1 - \left(\frac{\chi^{\frac{1}{2}}-1}{\chi^{\frac{1}{2}}+1} \right)^{2} e^{-2\chi^{\frac{1}{2}}\tau} a}$$
(3-6)

On factoring and rearranging into a form more suitable for our purposes, we obtain

$$\varepsilon = \frac{1 - e^{-\chi^{\frac{1}{2}}\tau} a}{1 + \frac{1}{2} (\chi^{\frac{1}{2}} - 1) (1 + e^{-\chi^{\frac{1}{2}}\tau} a)}$$
(3-7)

Briefly considering the limit of small $\beta\omega$ ($\chi \% 1$), we obtain a somewhat more transparent form

$$\varepsilon \approx \frac{1-e^{-\tau}s^{-\tau}a}{1+\frac{1}{2}\frac{\beta\omega}{1-\omega}\left(1+e^{-\tau}s^{-\tau}a\right)}$$
(3-8)

Because of the relatively small variation of the denominator from unity, the mean value over a spectral interval can be written as

$$\overline{\varepsilon} = \frac{1}{\Delta \omega} \int_{\Delta \omega} \varepsilon d\omega \approx \frac{1 - e^{-\tau} s e^{-f(\overline{\tau}_a)}}{1 + \frac{1}{2} (\frac{\overline{\beta} \omega}{1 - \omega}) \left(1 + e^{-\tau} s e^{-f(\overline{\tau}_a)}\right)}$$
(3-9)

where

$$f(\overline{\tau}_a) = -\ln \left\{ \frac{1}{\Delta \omega} \int_{\Delta \omega} e^{-\tau_a} d\omega \right\}$$
 (3-10)

represents the band-model curve-of-growth for the gas plus particle absorption component (that is, a known function). Equation (3-9) explicitly shows the modifications resulting from the simultaneous presence of scattering.

For χ significantly greater than unity, it still appears reasonable to consider the preponderant effect of spectral variation to occur in the numerator (which varies from unity to zero). The percentage error introduced by this assumption will be investigated in the numerical studies in Subtask II. We write

$$\overline{\epsilon} \approx \frac{1-e^{-f(\overline{\tau}_a, \tau_s)}}{1+\frac{1}{2}\left[(\overline{\chi})^{\frac{1}{2}}-1\right]\left(1+e^{-f(\overline{\tau}_a, \tau_s)}\right)}$$
(3-11)

where

$$f(\overline{\tau}_{a}, \tau_{s}) = -\ln \left\{ \frac{1}{\Delta \omega} \int_{\Delta \omega} e^{-(\tau_{a}^{2} + 2\tau_{a}\tau_{s})^{\frac{1}{2}}} d\omega \right\}$$
 (3-12)

We designate the numerator of Eq. (3-11) as 1-t' (since it

corresponds to the emissivity = 1-transmissivity in the

3.2.1 Elsasser model. We examine the behavior of the function \overline{t} ' when the absorption results from one gaseous component represented by the Elsasser model (for which case the effects of the gaseous spectral structure will be the greatest):

$$\tau_{a} = \overline{\tau}_{a} \frac{\sinh \beta}{\cosh \beta - \cos x}$$
 (3-13)

where

$$\overline{\tau}_{a} = \overline{k}u = Su/d$$
 (3-14)

(mean absorption optical depth)

$$\beta = 2\pi \gamma/d \tag{3-15}$$

(fine-structure parameter)

and

$$x = 2\pi(\omega - \omega_0)/d. \qquad (3-16)$$

Thus

$$\overline{t}' = \frac{1}{\pi} \int_{0}^{\pi} \exp \left[-\left(\overline{\tau}_{a}^{2} \frac{\tanh^{2} \beta}{(1 - \cos x/\cosh \beta)^{2}} + 2\overline{\tau}_{a}^{\tau} s \frac{\tanh \beta}{(1 - \cos x/\cosh \beta)}\right)^{\frac{1}{2}} \right] dx$$
(3-17)

The properties of this integral will be examined in detail for the limiting cases of small β and large β .

For large values of the parameter β (i.e., for significant overlapping of the spectral lines), the function \overline{t}' becomes (on expansion of the terms in 1-cosx/cosh β),

$$\overline{t}' = \exp\left[-(\overline{\tau}_a^2 \tanh^2 \beta + 2\overline{\tau}_a \tau_s \tanh \beta)^{\frac{1}{2}}\right]$$

$$\times I_o\left[(\overline{\tau}_a^2 \tanh^2 \beta + 2\overline{\tau}_a \tau_s \tanh \beta)^{\frac{1}{2}} \left(\frac{\overline{\tau}_a \tanh \beta + \tau_s}{\overline{\tau}_a \tanh \beta + 2\tau_s}\right)/\cosh\beta\right]$$
(3-18)

where I_0 is the modified Bessel function of the first kind. Since $\exp(-x) \approx 1-x$ for small x, and $I_0(y) \approx 1+\frac{1}{4}y^2$ for small y (and y << x), the I_0 term provides a higher-order correction to the exponential term. Thus a useful approximation is

$$\overline{t}' = \exp\left[-(\overline{\tau}_a^2 \tanh^2 \beta + 2\overline{\tau}_a \tau_s \tanh^2)\right]$$
 (3-19)

Note that for no scattering ($\tau_s \equiv 0$), we have

$$\overline{t} = 1 - \overline{\varepsilon} = \overline{t}' = \exp(-\overline{\tau}_a \tanh \beta),$$
 (3-20)

the classical Elsasser model result.

For the more interesting case of β small (little over-lapping of spectral lines), we have (using cosh $\beta \approx 1$)

$$(\overline{\tau}_{\mathbf{a}}^{2} + 2\overline{\tau}_{\mathbf{a}}\tau_{s})^{\frac{1}{2}} = \overline{\tau}_{\mathbf{a}} \left(\frac{\sinh^{2}\beta}{(1-\cos x)^{2}} + \frac{2\tau_{s}}{\overline{\tau}_{\mathbf{a}}} \frac{\sinh\beta}{(1-\cos x)} \right)^{\frac{1}{2}}$$

$$= \left(\overline{\tau}_{\mathbf{a}}^{2} \frac{\sinh^{2}\beta}{4 \sin^{4}x/2} + \overline{\tau}_{\mathbf{a}}\tau_{s} \frac{\sinh\beta}{\sin^{2}x/2} \right)^{\frac{1}{2}}$$

$$= \left(\frac{b}{\sin^{4}y} + \frac{a}{\sin^{2}y} \right)^{\frac{1}{2}}$$

$$(3-21)$$

where

$$a = \overline{\tau}_a \tau_s \sinh \beta \tag{3-22}$$

$$b = \frac{1}{4} \tau_a^2 \sinh^2 \beta$$
 (3-23)

and

$$y = \frac{1}{2} x$$

By making the substitution

$$(a+b)^{\frac{1}{2}}z = \left(\frac{a}{\sin^2 y} + \frac{b}{\sin^4 y}\right)^{\frac{1}{2}}$$

to transform to a more standard form (e.g., that of a Laplace transform), the function

$$\overline{t}' = \frac{2}{\pi} \int_{0}^{\pi/2} \exp\left[-\left(\frac{a}{\sin^2 y} + \frac{b}{\sin^4 y}\right)^2\right] dy$$
 (3-24)

becomes

$$\overline{t}' = -\frac{2}{\pi} \int_{1}^{\infty} \exp\left[-(a+b)^{\frac{1}{2}}z\right] \left\{ \frac{z}{\left[z^{2} - \frac{a+(a^{2}+4b(a+b)z^{2})^{\frac{1}{2}}}{2(a+b)}\right]^{\frac{1}{2}}} \right\} \\
\left\{ \left[2(a+b)\right]^{\frac{1}{2}} b \left[a+(a^{2}+4b(a+b)z^{2})^{\frac{1}{2}}\right]^{-\frac{1}{2}} (a^{2}+4b(a+b)z^{2})^{-\frac{1}{2}} \right\} \\
- \frac{1}{\left[2(a+b)\right]^{\frac{1}{2}}z^{2}} \left[a+(a^{2}+4b(a+b)z^{2})^{\frac{1}{2}}\right]^{\frac{1}{2}} dz .$$
(3-25)

This form is somewhat unwieldy, and does not appear to be immediately integrable. For the limiting case a=0 (no scattering), the expression above reduces to

$$\bar{t}' = \frac{1}{\pi} \int_{1}^{\infty} \exp(-\sqrt{b}z) \frac{1}{z(z-1)^{\frac{1}{2}}} dz$$
 (3-26)

$$= \operatorname{erfc}(b^{\frac{1}{4}}) = \operatorname{erfc}\left|\left(\overline{\tau}_{a}\beta/2\right)^{\frac{1}{2}}\right| , \qquad (3-27)$$

the usual Elsasser model result for small β .

In the other limiting case of b = 0 (more strictly speaking, $b/a \rightarrow 0$) the integral for \bar{t} ' reduces to

$$\bar{t}' = \frac{2}{\pi} \int_{1}^{\infty} \exp(-a^{\frac{1}{2}}z) \frac{1}{z(z^2-1)^{\frac{1}{2}}} dz$$
 (3-28)

This expression may be converted to a known Laplace transform, yielding

$$\bar{t}' = 1 - \frac{2}{\pi} \int_{0}^{a^{\frac{1}{2}}} K_{o}(x) dx$$
 (3-29)

where

$$a^{\frac{1}{2}} = (\overline{\tau}_a \tau_s \sinh \beta)^{\frac{1}{2}},$$

as before, and $K_O(x)$ is the modified Bessel function of the second kind. Such functions (of integral order) appear to be the least tractable of the Bessel functions. However, since $K_O(x)$ is only weakly (logarithmically) divergent at the origin, series expressions may be used and integrated over the finite interval indicated.

Because of the mathematical intractability of this Bessel function, a numerical study was made of \overline{t}' . From the power series definition of K_0 , and a term-by-term integration of Equation (3-29) we obtain

$$\int_{0}^{z} K_{0}(x) dx = \sum_{k=0}^{\infty} \left[-b_{k} (in(\frac{1}{2}z) + \gamma) + c_{k} \right] (\frac{2}{2k+1}) (\frac{1}{2}z)^{2k+1} , \quad (3-30)$$

where $\gamma = .5772156649...$, (Euler's constant)

$$b_{k} = \frac{1}{(k!)^{2}},$$

$$a_{k} = \frac{1}{(k!)^{2}} \sum_{i=1}^{k} \left(\frac{1}{i}\right),$$

$$c_{k} = a_{k} + \frac{1}{2k+1} b_{k}.$$

$$(3-31)$$

By use of recursion relations between the coefficients, the series is readily programmable, with high accuracy, even on a pocket calculator, for values of z as large as 12. However, for z greater than about 8, higher accuracy may be achieved by an asymptotic series. By expanding the exponential in a power series and integrating, the following expressions are obtained

$$\bar{t}' = \sqrt{\frac{2}{\pi}} \exp(-a^{\frac{1}{2}}) \sum_{j=0}^{\infty} (-1)^{j} d_{j} a^{-\frac{1}{2}(j+\frac{1}{2})},$$
 (3-32)

where

$$d_{j} = \sum_{i=0}^{j} \left(\frac{1 \cdot 3 \cdot 5 \cdot \dots (2i-1)}{i! \ 2^{2i}} \right) \left(1 \cdot \frac{1}{2} \cdot \frac{3}{2} \cdot \dots \ \frac{2j-1}{2} \right)$$
 (3-33)

This series is "semi-convergent" and must be truncated when the magnitude of the terms no longer decreases. For $a^{\frac{1}{2}}$ greater than about 8, it is capable of providing 3-place accuracy for \overline{t} '

deteriorates; for $a^2 \approx 1$, it is effectively worthless.

The function $t' = 1 - \frac{2}{\pi} \int_0^{a^{\frac{1}{2}}} K_0(x) dx$ is plotted in Figure 3-1 in Cartesian coordinates and the curve-of-growth $-\ln t'$ is plotted versus $a^{\frac{1}{2}}$ (log-log scale) over a much wider range of values of $a^{\frac{1}{2}}$ in Figure 3-2.

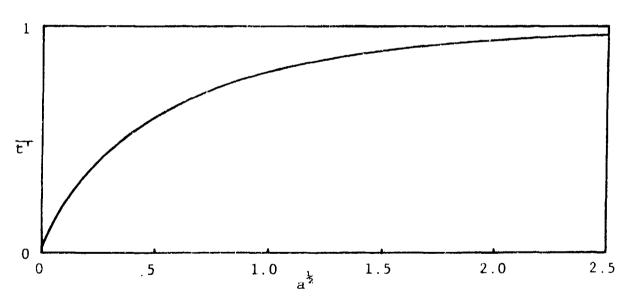


Figure 3-1. Function $\bar{t}' = 1 - \frac{2}{\pi} \int_{0}^{a^{\frac{1}{2}}} K_{o}(x) dx \text{ vs. } a^{\frac{1}{2}}$.

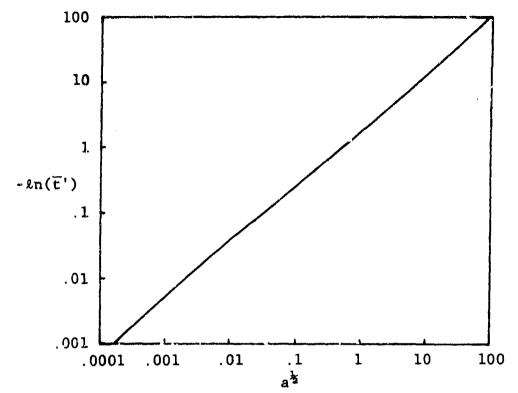


Figure 3-2. Curve-of-Growth $-\ln(\overline{t}')$ versus $a^{\frac{1}{2}}$ for $\overline{t}' = 1 - \frac{2}{\pi} \int_{0}^{a^{\frac{1}{2}}} K_{o}(x) dx$.

3.2.2 Random band models: The mean trasnmissivity for a given band model may be written alternatively

$$\overline{t} \left[= \frac{1}{\Delta \omega} \int_{\Delta \omega} t(k(\omega)) d\omega \right] = \int_{0}^{\infty} f(k) t(k) dk, \qquad (3-34)$$

in which f(k) is the probability density function of the absorption coefficient k. Because the properties of random band models without scattering are relatively simply derivable from those of single lines, explicit expressions for such probability densities have not been required for study of the

gaseous band models. However, because of the exponential form of the monochromatic transmissivity, the probability density function in the expression

$$\overline{t}(u) = \int_{0}^{\infty} f(k)e^{-ku} dk$$
 (3-35)

can be seen to be the inverse Laplace transform of the mean transmissivity function. A study of the inverse Laplace transforms of known transmissivity functions has been made by Domoto (1974), who has found that for the exponential-tailed S^{-1} model a relatively simple functional form exists, namely,

$$f(k) = \left[\frac{\overline{k} \ a}{\pi \ k^3}\right]^{\frac{1}{2}} \exp \left[a\left(2 - \frac{k}{\overline{k}} - \frac{\overline{k}}{k}\right)\right] , \qquad (3-36)$$

corresponding to the transmissivity function

$$\overline{t}(u) = \exp\left\{-2a\left[\left(1 + \frac{\overline{k}u}{a}\right)^{\frac{1}{2}} - 1\right]\right\} . \tag{3-37}$$

We consider here the case of no particulate absorption, for which the effects of the gaseous spectral structure will be the greatest.

The function $\overline{\mathbf{t}}'$ previously defined as

$$\overline{t}' = \frac{1}{\Delta \omega} \int_{\Delta \omega} \exp \left[- \left(\tau_a^2 + 2\tau_a \tau_s \right)^{\frac{1}{2}} \right] d\omega$$
 (3-38)

(where $\tau_a = ku$) can be rewritten

$$\overline{t}' = \int_0^\infty \left[\frac{\overline{\tau}_a}{\pi \tau_a} \frac{a}{3} \right]^{\frac{1}{2}} \exp \left[a \left(2 - \frac{\overline{\tau}_a}{\overline{\tau}_a} - \frac{\overline{\tau}_a}{\tau_a} \right) - \left(\tau_a^2 + 2\tau_a \tau_s \right)^{\frac{1}{2}} \right] d\tau_a \quad (3-39)$$

Direct integration is rendered difficult by the presence of the radical in the exponential term. Before studying the general behavior of \overline{t} ', some limiting cases (small τ_s , large τ_s) are examined.

a) $\underline{Small} \tau_s$

For small values of $\tau_{_{\bf S}}$ $(\tau_{_{\bf S}}<<\overline{\tau}_{_{\bf d}}),$ expansion of the radical yields

$$(\tau_a^2 + 2\tau_a^2 \tau_s)^{\frac{1}{2}} \approx \tau_a^2 + \tau_s^2 - \frac{1}{2} \frac{\tau_s^2}{\tau_a} \dots,$$
 (3-40)

in which terms in τ_s^3/τ_a^2 etc. have been suppressed. Substitution produces

$$\bar{t}' = \int_{0}^{\infty} \left(\frac{\bar{\tau}_{a} a}{\pi \tau_{a}^{3}}\right)^{\frac{1}{2}} \exp\left[2a - \tau_{s} - \left(\frac{a}{\bar{\tau}_{a}} + 1\right)\tau_{a} - \left(a\bar{\tau}_{a} - \frac{1}{2}\tau_{s}^{2}\right)\tau_{a}^{-1}\right] d\tau_{a}, (3-41)$$

or, defining

$$a^* = a \left(1 - \frac{\tau_s}{2a\overline{\tau}_a}\right)^{\frac{1}{2}}$$
, (3-42)

and

$$\overline{\tau}_{a}^{*} = \overline{\tau}_{a} \left(1 - \frac{\tau_{s}^{2}}{2a\overline{\tau}_{a}} \right)^{\frac{1}{2}} , \qquad (3-43)$$

$$\overline{t}' = e^{-\tau} s \left(1 - \frac{\tau_s^2}{2a\tau_a} \right)^{-\frac{1}{2}} \int_0^\infty \left(\frac{\overline{\tau}_a * a^*}{\pi \tau_a 3} \right)^{\frac{1}{2}} exp \left[2a - \left(\frac{a^*}{\overline{\tau}_a *} + 1 \right) \tau_a - \frac{a^* \overline{\tau}_a *}{\tau_a} \right] d\tau_a^{(3-44)}$$

Since the integral is now in the standard form with new constants, we have

$$\bar{t}' = e^{-\tau_s} \left(1 - \frac{\tau_s^2}{2a\tau_a} \right)^{-\frac{1}{2}} \exp \left\{ -2a^* \left[\left(1 + \frac{\tau_a^*}{a^*} \right)^{\frac{1}{2}} - 1 \right] \right\}$$
 (3-45)

or

$$\bar{t}' = e^{-\tau s} \left(1 - \frac{\tau s^2}{2a\bar{\tau}_a} \right)^{-\frac{1}{2}} \exp \left\{ -2a \left(1 - \frac{\tau s^2}{2a\bar{\tau}_a} \right)^{\frac{1}{2}} \left[\left(1 + \frac{\bar{\tau}_a}{a} \right)^{\frac{1}{2}} - 1 \right] \right\}$$
 (3-46)

To the same degree of approximation we may write

$$\overline{t}' = \exp\left\{-2a\left(1 - \frac{\tau_s^2}{2a\overline{\tau}_a}\right)^{\frac{1}{2}} \left[\left(1 + \frac{\overline{\tau}_a}{a}\right)^{\frac{1}{2}} - 1\right] - \tau_s + \frac{\tau_s^2}{4a\overline{\tau}_a}\right\}$$
 (3-47)

b) Large τ_s

In the limit of large τ_s the magnitude of the argument of the exponential term in the integral becomes large, and the minimum progressively less shallow, and occurring at smaller values of τ_a . If $\tau_s >> \tau_a$, so that

$$(\tau_a^2 + 2\tau_a\tau_s)^{\frac{1}{2}} \approx (2\tau_a\tau_s)^{\frac{1}{2}}$$
, (3-48)

the argument of the exponential term becomes

$$\mathbf{E}(\tau_{\mathbf{a}}) = -\mathbf{a} \frac{\overline{\tau}_{\mathbf{a}}}{\tau_{\mathbf{a}}} \left(1 - \frac{\tau_{\mathbf{a}}}{\overline{\tau}_{\mathbf{a}}}\right)^{2} - \left(2\tau_{\mathbf{a}}\tau_{\mathbf{s}}\right)^{\frac{1}{2}} . \tag{3-49}$$

The magnitude of this argument has a minimum at approximately

$$\tau_a^{\min} \approx (\tau_s/2)^{-\frac{1}{3}} (a\overline{\tau}_a)^{\frac{2}{3}}$$
 , (3-50)

so that for large τ_s , $\tau_a^{min} << \overline{\tau}_a$.

In the region of this minimum, the argument may be expanded:

$$E(\tau_{\mathbf{a}}) \approx -3(\frac{1}{2}a\overline{\tau}_{\mathbf{a}}\tau_{\mathbf{s}})^{1/3}$$

$$-\frac{3}{4}(\frac{\tau_{\mathbf{s}}}{2a\overline{\tau}_{\mathbf{a}}})(\tau_{\mathbf{a}}-\tau_{\mathbf{a}}^{\min})^{2}+\dots \qquad (3-51)$$

the expression for $\overline{\overline{v}}$ ' then becomes

$$\overline{t}' = \int_{0}^{\infty} \left(\frac{\overline{\tau}_{a} a}{\pi \tau_{a}^{3}} \right)^{\frac{1}{2}} \exp \left[-3 \left(\frac{1}{2} a \overline{\tau}_{a} \tau_{s} \right)^{\frac{1}{3}} - \frac{3}{4} \left(\frac{\tau_{s}}{2 a \overline{\tau}_{a}} \right) \left(\tau_{a} - \tau_{a}^{\min} \right)^{2} \right] d\tau_{a} . \tag{3-52}$$

By recourse to the mean value theorem, noting the narrowness of the maximum of the exponential in the neighborhood of τ_a^{min} , the multiplicative factor of $\tau_a^{-3/2}$ may be taken outside the integral and evaluated at τ_a^{min} :

$$\overline{t}' \approx \left(\frac{\overline{\tau_a} a}{\pi \left(\tau_a^{\min}\right)^3}\right)^{\frac{1}{2}} \exp\left[-3\left(\frac{1}{2}a\overline{\tau_a}\tau_s\right)^{\frac{1}{3}}\right] \int_{0}^{\infty} \exp\left[-\frac{3}{4}\left(\frac{\tau_s}{2a\overline{\tau_a}}\right)\left(\tau_a - \tau_a^{\min}\right)^2\right] d\tau_a. \quad (3-53)$$

The remaining integration may be performed by letting the lower bound $\rightarrow -\infty$:

$$\overline{t}' \approx \sqrt{\frac{4}{3}} \exp\left[-3\left(\frac{1}{2}a\overline{\tau}_a\tau_s\right)^{1/3}\right]. \tag{3-54}$$

This functional dependence provides an interesting analogy to the square-root region for pure gaseous absorption in which

$$\overline{t} \rightarrow \exp\left[-2(a_{\tau_a})^{\frac{1}{2}}\right]$$

c) Arbitrary T

Numerical studies have been carried out on the integral representation of the function \overline{t} ', given by Eq. (3-39). Some results of these calculations are shown in Figures 3-3 to 3-5, in which \overline{t} ' is plotted against $\overline{\tau}_a$, the mean absorption optical depth, for various values of τ_s , the scattering optical depth and a , the fine structure coefficient. The linear scale for $\overline{\tau}$ ' does not show the precise asymptotic behavior at $\overline{\tau}$ ' very close to 0 or 1, but does indicate a smooth, well-behaved dependence on τ_s , encouraging confidence in the approach of seeking an analytical or semi-empirical functional representation.

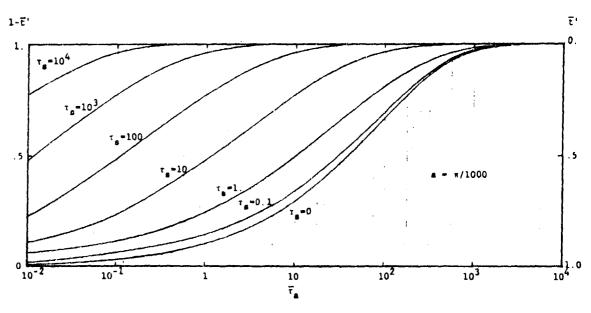


Figure 3-3. Plot of \overline{t}' vs. $\overline{\tau}_a$ for $a = \pi/1000$ and τ_s ranging from 0 to 10⁴.

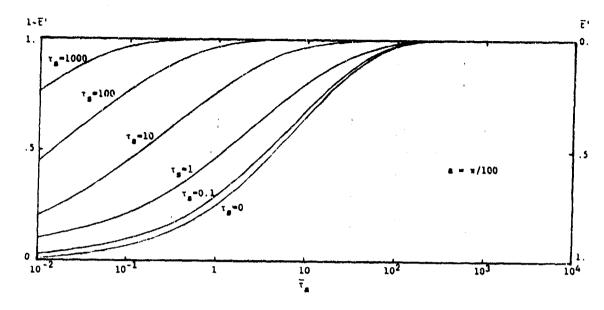


Figure 3-4. Plot of \bar{t} ' vs. $\bar{\tau}_a$ for $a = \pi/100$ and τ_s ranging from 0 to 10^3 .

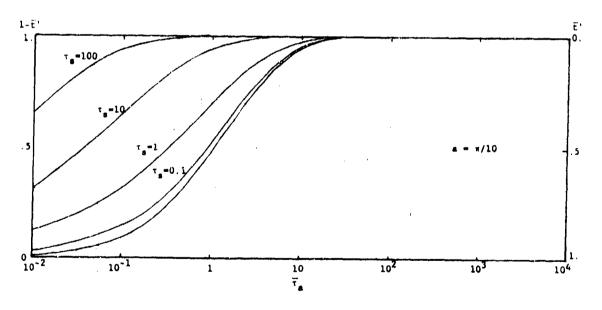


Figure 3-5. Plot of \overline{t} ' vs. $\overline{\tau}_a$ for $a = \pi/10$ and τ_s ranging from 0 to 10^2 .

The asymptotic region of small \overline{t} ' was examined in detail. The conclusion was reached that the particular asymptotic form developed above [see Eq. (3-54)] is of major practical significance. A plot of the ratio \overline{t} '/exp[-3($\frac{1}{2}a\overline{\tau}_a\tau_s$) $\frac{1}{3}$] is presented in Figure 3-6. This ratio can be seen to approach the predicted theoretical value of $(4/3)^{\frac{1}{2}} = 1.1547$ closely as a+0 even for small values of

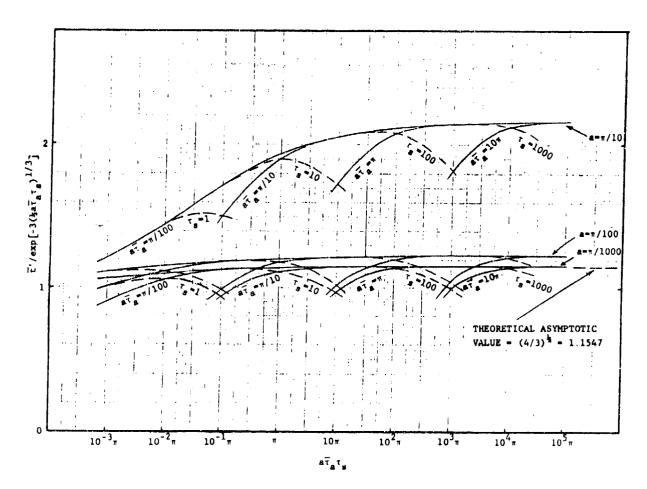


Figure 3-6. Plot of $\overline{t}'/\exp[-3(\frac{1}{2}a\overline{\tau}_a\tau_s)^{1/3}]$ vs. $a\overline{\tau}_a\tau_s$ showing asymptotic behavior for "large" τ_s (ranging from 1 to 10^3), for various values of a and $a\overline{\tau}_a$. The range of values of $a\overline{\tau}_a\tau_s$ corresponds to a range of \overline{t}' from about 0.7 at the left to essentially zero at the right.

 $a\overline{\tau}_a \tau_s$, when τ_s is sufficiently large. Empirically, for a $\lesssim 0.05$ and $\tau_s \gtrsim 10 a \overline{\tau}_a$, the asymptotic approach is quite close (within about 10%, even for $a\overline{\tau}_a \tau_s$ as small as 10^{-2} , corresponding to transmissivities as large as 0.6). For larger values of a, an asymptotic value is approached, whose magnitude depends on the value of a. The particular functional dependence could be determined theoretically or empirically (e.g., an asymptotic value of $(4/3)^{\frac{1}{2}} \exp(2a)$ provides an excellent fit over the range of values of a considered here).

3.2.3 Empirical formulations for scattering band models. A study was made of possible empirical forms in the scattering band model representation based on the previous theoretical studies and numerical calculations. No exact explicit expressions were derived for the general case, but various limiting cases were studied, and any empirical form considered should have the appropriate theoretical asymptotic behavior. The numerical calculations then serve to select the best representation for intermediate values of the parameters.

Specifically, the function previously defined at \overline{t} ' (analogous to the mean transmissivity function \overline{t} for the no-scattering limit) is desired to have the following limits:

- A) No scattering $(\tau_s=0)$ t' $\rightarrow \overline{t}=\exp\left[-f(\overline{\tau}_a,a)\right]$ where f is the appropriate gaseous band model curve of growth, in which a is the fine-structure parameters and $\overline{\tau}_a$ the mean absorption optical depth (= \overline{ku}_x where \overline{k} is the gaseous absorption coefficient and u_x the amount of the absorbing gas, in reciprocal units). The function f may include particulate absorption, i.e., be replaced by $f(\overline{\tau}_a, a) + \overline{\tau}_o$, where f is the pure gas curve of growth and $\overline{\tau}_o$ is a spectrally constant absorption optical depth (gaseous and/or particulate).
- B) Weak scattering $(\tau_s \leqslant \overline{\tau}_a)$ $\overline{t}' \approx \exp \left[-f(\overline{\tau}_a, a) \tau_s\right]$

in which f may include particulate absorption as in (A).

C) Constant absorption coefficient (gaseous and/or particulate).

 $(a \to \infty; f(\overline{\tau}_a, a) \to \overline{\tau}_a \text{ for any band model})$ $\overline{t}' \to \exp\left[-(\overline{\tau}_a^2 + 2\overline{\tau}_a\tau_s)^{\frac{1}{2}}\right]$ (this is a functional requirement of either the two-flux or six-flux treatment for constant absorption coefficient).

D) Strong-line, large scattering limit (no particulate absorption)

$$(\overline{\tau}_a/a \text{ large}, \tau_s \text{ large})$$
 $\overline{t}' \rightarrow c_1 \exp \left[-c_2 (\overline{\tau}_a a \tau_s)^{1/3} \right]$

for a random band model, with $c_1 \sim 1$ and $c_2 \sim 2.4$. This important limiting behavior was derived in the previous study and is analogous to the strong-line limit for gaseous absorption $\overline{t} \rightarrow \exp^{-2(\sqrt{t_a} - a)^{\frac{t_a}{2}}}$.

An empirical form satisfying (A), (B), and (C) is

$$\overline{t}' = \exp \left[-f(\overline{\tau}_a, a) \left\{ 1 + \frac{2 \tau_s}{f(\overline{\tau}_a, a)} \right\}_{2}^{\frac{1}{2}} \right]$$
 (3-55)

This is the form utilized intuitively in the Vanderbilt-Slack program. However, in limiting case (D) it approaches $\overline{t}' \rightarrow \exp(-2\overline{\tau_a}^{\frac{1}{4}} a^{\frac{1}{4}} \tau_s^{\frac{1}{2}}).$

Alternatively, a form satisfying (A), (B), and (D) is

$$\overline{t}' = \exp \left[-f(\overline{\tau}_a, a) \right] \left\{ 1 + \frac{3 \tau_s}{f(\overline{\tau}_a, a)} \right\}$$
 (3-56)

Although this form provides a better fit to the numerical values calculated for intermediate ranges of the parameters, it does not have the proper form to match limit (C); instead $\overline{t}' \rightarrow \exp \left[-(\overline{\tau_a}^3 + 3\overline{\tau_a}^2\tau_s)^{1/3}\right].$ For this reason, a more elaborate form containing all four limiting cases above is sought.

A convenient form having these required limits is

$$\overline{t}' = \exp\left\{-f(\overline{\tau}_a, a)\left[1 + \frac{2\tau_s}{f(\overline{\tau}_a, a)}\right]^{\frac{1}{2}} \left[\frac{\left|(\overline{\tau}_a + c)a\right|^n}{\left|(\overline{\tau}_a + c)a\right|^n + \left|b\tau_s^2(1 - \frac{f(\overline{\tau}_a, a)}{\tau_a})\right|^n}\right]^{\frac{1}{12n}}\right\} (3-57)$$

where b, c and n are empirical constants. Numerical evaluation of this equation has shown significantly improved agreement relative to the previous forms. The curve of growth $f(\overline{\tau}_a,a)$ may include a particulate absorption optical depth.

Figure 3-7, -9, and -11 show comparisons of the exact values of \overline{t} ' previously calculated for $\pi/10^3$ < a < $\pi/10$ (solid lines) with the approximation given by Equation (3-55) (dashed line). This approximation is seen to be especially poor for large τ_o and small a (i.e., when condition (D) should be met).

Figure 3-8, -10 and -12 show comparisons of the same values of \bar{t}' with the empirical form given by Equation (3-57) for the case c=1, n=1, b=1/8. The improved match overall is evident. No effort was made at optimization of the parameters b, c, and n. It is anticiapted that an optimized version of this form (or a very similar one) will be used in the basic formulation of the SIRRM band model representation.

All the results shown here have been obtained for the exponential-tailed 1/S random band model. The expression for the absorption coefficient probability distribution function for the Goody model is so mathematically intractable as to make it extremely unattractive for similar theoretical or numerical studies. However, the conclusions reached here regarding proper asymptotic and approximate forms are certainly also applicable to the Goody model (whose curve of growth for the non-scattering case is only slightly different from that of the exponential-tailed 1/S model).

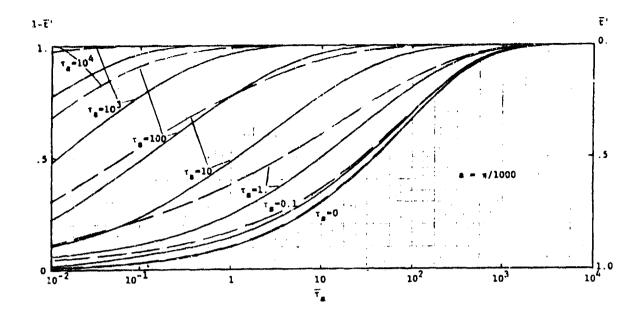


Figure 3-7. Plot of \overline{t} ' vs. $\overline{\tau}_a$ for $a=\pi/1000$ and τ_s ranging from 0 to 10° compared with approximation (1).

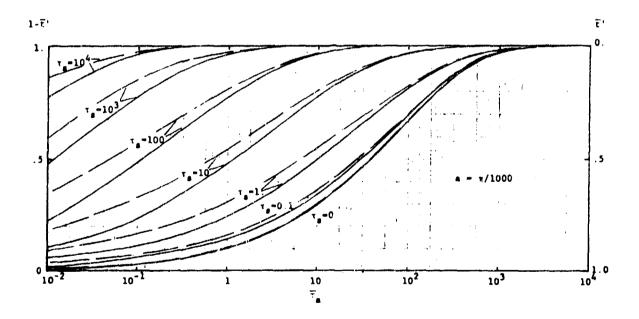


Figure 3-8. Plot of \overline{t} ' vs. $\overline{\tau}_a$ for $a = \pi/1000$ and τ_s ranging from 0 to 10° compared with approximation (3).

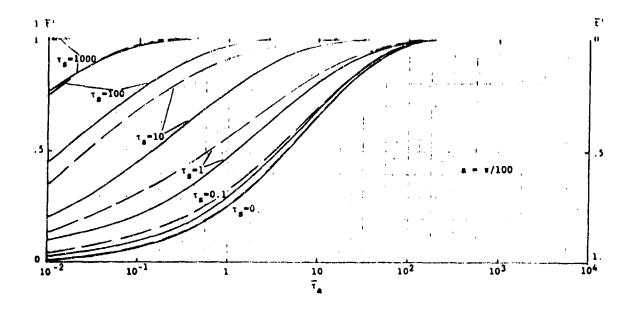


Figure 3-9. Plot of \overline{t} ' vs. $\overline{\tau}_a$ for $a = \pi/100$ and τ_s ranging from 0 to 10^3 compared with approximation (1).

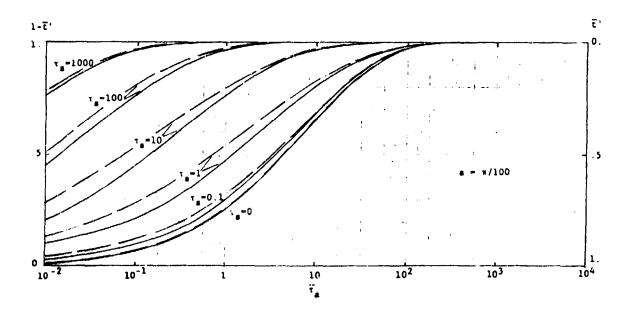


Figure 3-10. Plot of \overline{t} ' vs. $\overline{\tau}_a$ for $a=\pi/100$ and τ_s ranging from 0 to 10^3 compared with approximation (3).

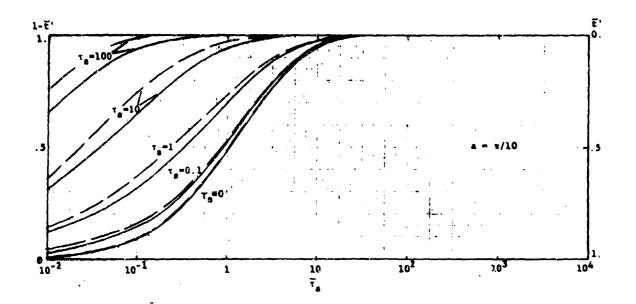


Figure 3-11. Plot of \overline{t}' vs. $\overline{\tau}_a$ for $a = \pi/10$ and τ_s ranging from 0 to 10^2 compared with approximation (1).

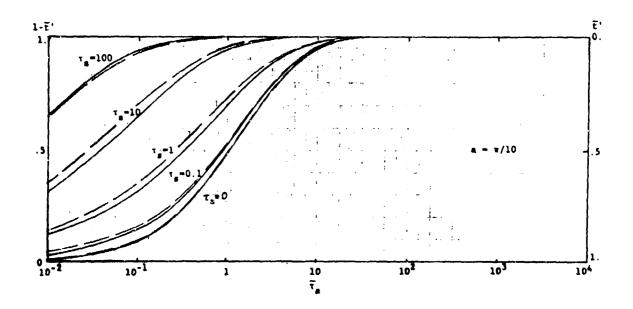


Figure 3-12. Plot of \bar{t}' vs. $\bar{\tau}_a$ for a = $\pi/10$ and τ_s ranging from 0 to 10^2 compated with approximation (3).

The preceding development is for single band models (with or without particle or other gaseous absorption). In practice, it is anticipated that methods of combining band model parameters will be utilized which are analogous to those in current use in gaseous radiative transfer codes (e.g., the Curtis-Godson approximation).

OPTICAL PARAMETERS OF PARTICLES

4.1. Introduction

The treatment of the gas/particles radiative transfer requires as inputs the volume extinction coefficients and angular scattering phase functions. These quantities can be derived from the primary optical parameters by use of the Mie theory, which describes the scattering and absorption properties of a homogeneous sphere whose optical properties are given by the complex index of refraction. Although the Mie theory is based upon the idealization of homogeneous spherical particles, it can be considered as a very reasonable first approximation since real plume particles were often found to be solid spheres (Rochelle, 1967). In addition, higher order account of particle non-uniformity and non-sphericity is prohibitively complex (Kerker, 1969) and not warranted by the present state of knowledge. Exact solutions of the wave equations are obtained by the matching of boundary conditions at the surface of the sphere. The solutions are expressed as infinite series involving Legendre polynomials and Ricatti-Bessel functions. Evaluation of the formal series solution proves computationally difficult for some ranges of the parameters and considerable effort has been expended in the past on the development of efficient and accurate computational methods of evaluating these solutions. The most recent treatment is by Dave (1972) and it is his approach that is being used in the present program. A brief discussion together with the extension from single particles to particle distributions is given in Section 4.2.

A preliminary listing of the particulate species of interest is given in Table 4-1 together with an indication of the availability of the index of refraction (n_1) and the absorption index (n_2) measured in a given temperature interval. Only four out of nine particulate species are sufficiently known, and these will have to serve as representatives until more information about the remaining species becomes available.

TABLE 4-1. LIST OF EXHAUST PARTICULATE SPECIES AND THE AVAILABILITY OF n₁ and n₂.

PARTICLE		η ₁ (Τ,λ)	η ₂ (Τ,λ)	SOURCE
Al ₂ 0 ₃	(s)	300-1800K	300-1800	Aerospace
ç	(s)	300K	300-3000	Aerospace/PRA Estimate
Mg0	(s)	300-2225K	300-2225	Aerospace
ZrO ₂	(s)	300-600K	300-600	Aerospace
ZrC	(s)			
BN	(s)	300K(Na-D)		Kordes (1960)
$B_{2}O_{3}$	(1)	300K(Na-D)		Handbook Chem. Phys. for (s)
Be0	(s)	300K(Na-D)		Handbook Chem. Phys.
Be ₃ N ₂				

Representative results of volume coefficients and angular scattering phase functions for Al_2O_3 and carbon particle distributions are given in Section 4.3. The results indicate that the large variations in coefficients and phase functions typically observed as a function of wavelength and scattering angle for single particles are greatly reduced when realistic size distributions are introduced. In addition, the influence of the

A preliminary listing of the particulate species of interest is given in Table 4-1 together with an indication of the availability of the index of refraction (n_1) and the absorption index (n_2) measured in a given temperature interval. Only four out of nine particulate species are sufficiently known, and these will have to serve as representatives until more information about the remaining species becomes available.

TABLE 4-1. LIST OF EXHAUST PARTICULATE SPECIES AND THE AVAILABILITY OF \mathbf{n}_1 and \mathbf{n}_2 .

PARTICLE		$\eta_1(T,\lambda)$	η ₂ (Τ,λ)	SOURCE	
A203	(s)	300-1800K	300-1800	Aerospace	
ç į	(s)	300K	300-3000	Aerospace/PRA Estimate	
Mg0	(s)	300-2225K	300-2225	Aerospace	
ZrO2	(s)	300-600K	300-600	Aerospace	
ZrC	(s)				
BN	(s)	300K(Na-D)		Kordes (1960)	
B ₂ 0 ₃	(%)	300K(Na-D)		Handbook Chem. Phys. for (s)	
Be0	(s)	300K(Na-D)		Handbook Chem. Phys.	
Be ₃ N ₂					

Representative results of volume coefficients and angular scattering phase functions for $\mathrm{Al}_2\mathrm{O}_3$ and carbon particle distributions are given in Section 4.3. The results indicate that the large variations in coefficients and phase functions typically observed as a function of wavelength and scattering angle for single particles are greatly reduced when realistic size distributions are introduced. In addition, the influence of the

OPTICAL PARAMETERS OF PARTICLES

4.1. Introduction

The treatment of the gas/particles radiative transfer requires as inputs the volume extinction coefficients and angular scattering phase functions. These quantities can be derived from the primary optical parameters by use of the Mie theory, which describes the scattering and absorption properties of a homogeneous sphere whose optical properties are given by the complex index of refraction. Although the Mie theory is based upon the idealization of homogeneous spherical particles, it can be considered as a very reasonable first approximation since real plume particles were often found to be solid spheres (Rochelle, 1967). addition, higher order account of particle non-uniformity and non-sphericity is prohibitively complex (Kerker, 1969) and not warranted by the present state of knowledge. Exact solutions of the wave equations are obtained by the matching of boundary conditions at the surface of the sphere. The solutions are expressed as infinite series involving Legendre polynomials and Ricatti-Bessel functions. Evaluation of the formal series solution proves computationally difficult for some ranges of the parameters and considerable effort has been expended in the past on the development of efficient and accurate computational methods of evaluating these solutions. The most recent treatment is by Dave (1972) and it is his approach that is being used in the present program. A brief discussion together with the extension from single particles to particle distributions is given in Section 4.2.

temperature on the coefficients and phase functions for solid carbon is minimal.

In the last section, we recommend a scheme for the Data File that combines flexibility and minimum requirements in the operation of the SIRRM code.

4.2 Formalism of the Mie Scattering

4.2.1 Single particle. In the Mie theory, the primary quantities:

$$S_1(a,n,\Theta)$$

$$S_2(a,n,\Theta)$$

$$Q_{sca}(a,n)$$

are calculated (Van de Hulst, 1957) where S_1 and S_2 are amplitude functions related to the planes of polarization, $Q_{\rm ext}$ and $Q_{\rm sca}$ are the extinction and scattering efficiencies, a is the size parameter (a=2 π r/ λ , with radius r of scattering sphere and wavelength λ), n is the refractive index (n=n(λ) = n₁(λ)-in₂(λ)) and θ is the angle between the directions of incoming and scattered radiation. After the primary quantities are calculated, the intensities of the scattered radiation and their degrees of polarization can be determined. The calculation is simplified by repeated use of recursion relationships and efficient computing methods have been developed (Deirmendjian, 1969; Dave 1969, 1970A, 1970B). In the following, the approach taken by Dave (1972) is briefly described.

The intensity scattered by a single spherical particle per unit solid angle for unity incident intensity is given by

$$I(a,n,\theta) = \frac{1}{4\pi} (\pi r^2) Q_s(a,n) M(a,n,\theta)$$
 (4-1)

where $M(a,n,\theta)$ is the normalized scattering phase function, which is given by

$$M(a,n,\theta) = \sum_{k=1}^{n} \Lambda_k(a,n) P_{k-1}(\cos \theta) \qquad (4-2)$$

where $P_{k-1}(\cos \theta)$ are the Legendre polynomials. The coefficients of the Legendre series are given by

$$\Lambda_{k}(a,n) = \frac{2}{a^{2}Q_{s}(a,n)} \left[L_{k}^{(1)}(a,n) + L_{k}^{(2)}(a,n) \right]$$

where the functions $L_k^{(1),(2)}(a,n)$ are computed from the values of the complex Mie amplitude $a_n(a,n)$ and $b_n(a,n)$. The extinction and scattering efficiencies are given by

$$Q_{\text{ext}}(a,n) = \frac{2}{a^2} \sum_{n=1}^{\infty} (2n+1) \operatorname{Re}\{a_n + b_n\}$$

$$Q_{\text{sca}}(a,n) = \frac{2}{a^2} \sum_{n=1}^{\infty} (2n+1) \{|a_n|^2 + |b_n|^2\}$$
(4-3)

The amplitudes a_n and b_n are expressible in terms of the spherical Bessel functions of the first and second kind. Using the method of logarithmic derivative function $A_n(a,n)$ of the Bessel functions, the numerical computations are facilitated. However, difficulties

exist in computing $A_n(a,n)$ under certain conditions (Deirmendjian, 1969). Kattawar and Plass (1967) found that these difficulties can be avoided by using a downward recurrence scheme for $A_n(a,n)$ which is basically stable. Dave (1972) has adopted this scheme in his approach.

4.2.2 Particle size distribution. Since rocket exhaust plumes do not, in general, consist of particles with only one fixed size, it is necessary to investigate the characteristics of size distributions. These characteristics are described by the volume extinction coefficient $\beta_{\rm ext}$, the volume scattering coefficient $\beta_{\rm sca}$, the volume absorption coefficient $\beta_{\rm abs}$ and the normalized scattering phase function $M(\theta)$. These quantities are dependent on $n(\lambda)$ and various parameters describing different size distributions. They can be expressed as

$$M(\Theta) = \sum_{k=1}^{K} \Lambda_k(r_j) P_{k-1}(\cos \Theta)$$
 (4-4)

and

$$\beta_{z}(r_{j}) = \pi \int_{r_{min}}^{r_{j}} Q_{z}(a,n)r^{2} N(r)dr \qquad (4-5)$$

where z stands for ext, sca or abs. The normalized Legendre coefficients become

$$\Lambda_{k}(r_{j}) = \frac{\lambda^{2}}{\pi \beta_{sca}(r_{j})} \int_{r_{min}}^{r_{j}} L_{k}(a,n) N(r) dr \qquad (4-6)$$

The phase function is normalized according to

$$\int_{4\pi} M(\theta) \sin\theta d\theta d\phi = 4\pi$$

The distribution N(r) is given by

$$N(r) = Nf(r) (4-7)$$

where N is total number of particles per unit volume and $\int_{0}^{\infty} f(r)dr = 1$. Dave (1972) has made provisions to use any one of three different size distribution functions. We are interested in the discontinuous power law type

$$f(r) = C \qquad \text{for } r_{\min} \leqslant r \leqslant r_{m}$$

$$f(r) = C(r_{m}/r)^{\nu+1} \text{ for } r_{m} \leqslant r \leqslant r_{\max}$$

$$(4-8)$$

and the modified gamma function

$$f(r) = a r^{\alpha} e^{-br^{\gamma}} for r_{min} \leq r \leq r_{max}$$
 (4-9)

It is found that the modified gamma function is representative of size distributions encountered in exhaust plumes with Al_2O_3 particles (Dawbarn, 1978). However, many different size distributions have been encountered which sometimes depend upon the method of collection (see Figure 4-1). The coefficients a, α , b and γ can be adjusted to represent any of the measured or calculated size distributions.

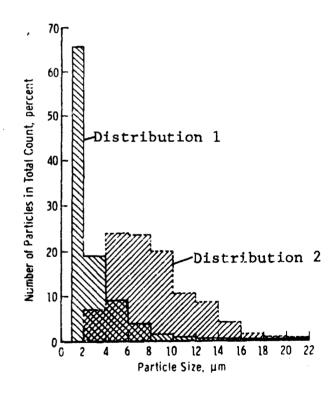


Figure 4-1. Two Al₂O₃ particle size distributions of Titan IIIc using two different collection methods (Dawbarn, 1978)

The parameters for the modified gamma function are not independent of each other and they have to be adjusted in a certain manner to obtain a given size distribution that has its maximum at radius $\mathbf{r}_{\mathbf{m}}$. This interdependence was investigated. The four parameters may be reduced to two by utilizing the conditions

$$\frac{d f(r)}{dr} = 0 \text{ for } r = r_{m}$$
 (4-10)

and

$$\int_{c}^{\infty} f(r) = 1$$

These conditions lead to

$$\alpha/b = \gamma r_{m}^{\gamma} \qquad (4-11)$$

$$1 = a \gamma^{-1} b^{-\xi} \Gamma(\xi)$$
 (4-12)

where $\xi=\frac{\alpha+1}{\gamma}$ and $\Gamma(\xi)$ is the gamma function. At $r=r_m$, f(r) must assume a value between zero and unity (as determined by experiment). Introducing Equations (4-11) and (4-12) into (4-10) and solving for r_m results in

$$f(r) = \gamma^{1-\xi} \frac{\alpha^{\xi}}{\Gamma(\xi)} \frac{e^{-\alpha/\gamma}}{r_{m}}$$
 (4-13)

Assuming upper and lower values for $f(r_m)$, Equation (4-13) has been solved for $f(r_m) \sim 1$ and $f(r_m) \sim .2$ using different values of α and γ . The solutions are shown graphically in Figure 4-2. A family of curves is made up for a given α , the upper and

lower limits of which are given by $f(r_m) \sim 1$ and $f(r_m) \sim 2$. Thus, pairs of α and γ can be chosen to describe any size distribution for r_m between 0.01 and 10 μm . The choice is dictated by the known (measured) standard deviation of the distribution function. The higher values of α and γ produce smaller standard deviations.

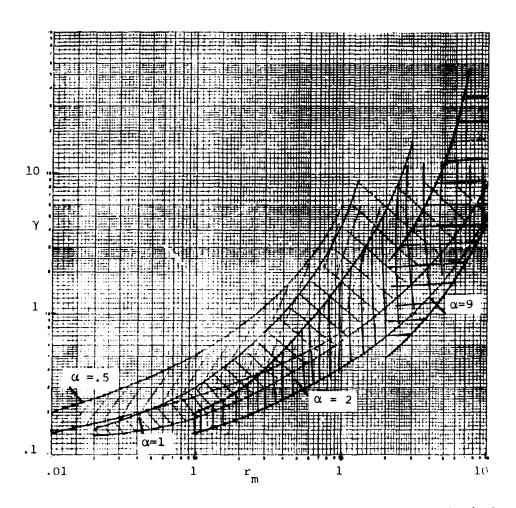


Figure 4-2. Families of Curves for which $0.2^{\circ}\,f(r_m)\!\simeq\!1$ at a given r_m as determined by pairs of γ and α .

Following a similar argument as was presented above for the modified gamma function, the adjustable parameters in the discontinuous power law (Eq. 4-8) are determined through

$$\int_{0}^{\infty} f(\mathbf{r}) = C \int_{0}^{\mathbf{r}_{m}} d\mathbf{r} + C r_{m}^{\nu+1} \int_{\mathbf{r}_{m}}^{\infty} r^{-(\nu+1)} d\mathbf{r} = 1$$

$$v = \frac{r_{m}}{1/C - r_{m}}$$

Based on a review of existing experimental data (Rochelle, 1967), we have chosen $r_m = 0.03~\mu m$ for solid carbon particles and C = 1. With these assumptions, then $v \approx 0$. The size distribution is shown in Figure 4-3. It resembles that used by Stull and Plass (1960) which was based on data by Tesner (1959). Included in the same figure is also the size distribution based on Equation (4-9).

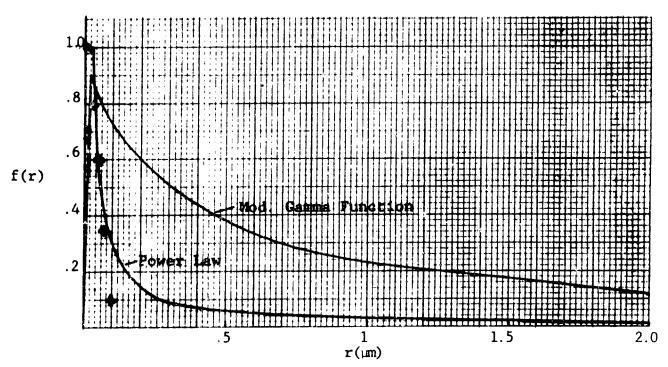


Figure 4-3. Size Distribution of Solid Carbon Particles in Rocket Exhaust, based on the Two Distribution Functions Described in the Text. Distribution used by Stull and Plass (1960) is indicated by (+).

4.3 Optical Parameters

A set of representative values of n_1 and n_2 were collected by the Aerospace Corporation for Al_2O_3 , C, MgO, and ZrO_2 (Whitson, 1975). Additional data for these particles were presented by Dowling and Randall (1977) in part based on experiments conducted during the "Particle Optical Properties Measurement (POPM) Program". Our effort was concentrated in comparing these data and study the influence of higher temperatures.

Carbon

As an example of the wide differences of measured refractive index n_1 , and absorption n_2 , Dowling and Randall have compared data on polycrystalline and pyrolytic graphite, pressed carbon soot, and a microcrystalline polished slab of graphite. The results for n_2 are shown in Figure 4-4. We have added the NASA Handbook

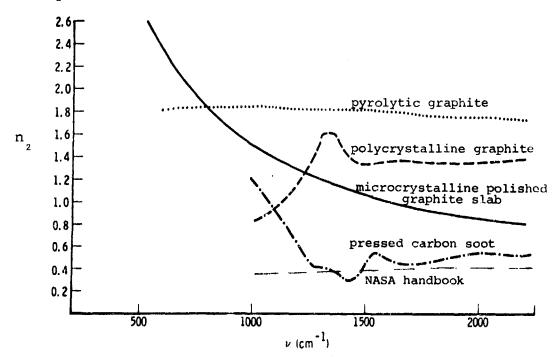


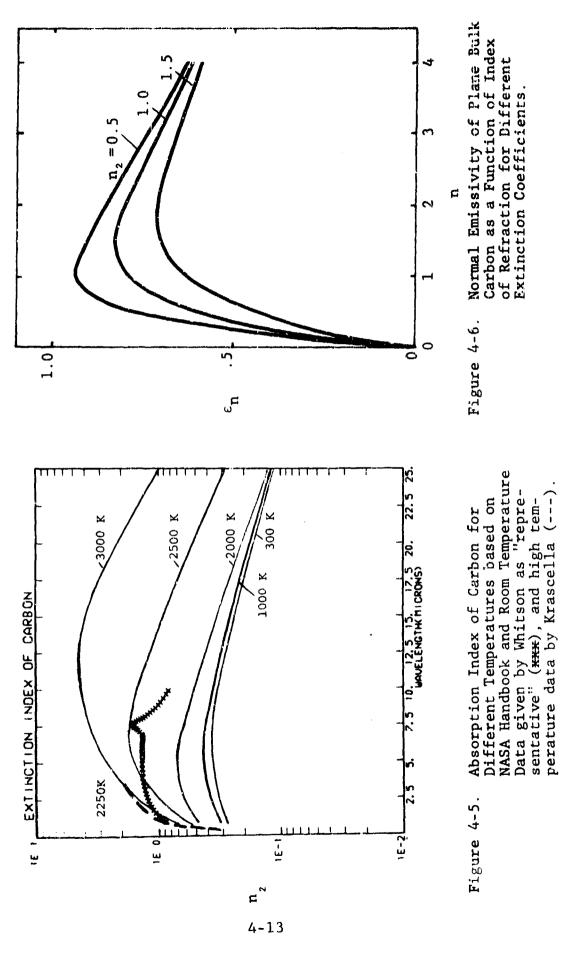
Figure 4-4. The Absorption Index of Carbon Materials in the Wavenumber Region 400 cm⁻¹ to 2200 cm⁻¹.

data for room temperature. The "representative" values chosen by Whitson are those for the polycrystalline graphite.

In addition to room temperature data, the NASA Handbook lists high temperature data for the absorption coefficient of small carbon particles. These data were based in part on theoretical work and in part on measurements. We have extrapolated those data to 3000 K and to λ = 25 μm , using the relationship

$$k = k_c \frac{\lambda}{2\pi} \rho_0$$

to convert the absorption coefficients k_c (cm²/gm) into absorption indices, where ρ_0 = 2g/cm³ was chosen as a representative value for the density of carbon. The results are plotted in Figure 4-5. We recommend to use these values for n_2 until more experimental data become available. For large particles compared with λ , the computed particle emissivity approaches the emissivity of a plane bulk sample. Dowling and Randall have calculated ϵ_n = .71 for λ < 25 µm, choosing n_1 = 3 and n_2 = 1. Using the same formula, we have calculated ϵ_n parametrically for $0 < n_1 < 4$ and $0.5 < n_2 < 1.5$. The results are shown in Figure 4-6. Measured spectral emissivities as reported by Whitson range from .8 to 1.0 for different samples of graphite (except for pyrolytic, c-face) in the temperature regime from 1100 to 2800 K, pointing to a somewhat lower value of n_1 at higher temperatures.



Normal Emissivity of Plane Bulk Carbon as a Function of Index of Refraction for Different Extinction Coefficients.

Figure 4-6.

Figure 4-5.

The only high temperature data for n listed by Whitson are due to N. K. Krascella (1966), and are plotted in Figure 4-7 together with Whitson's "representative" values at room temperature and our extrapolation to longer wavelengths.

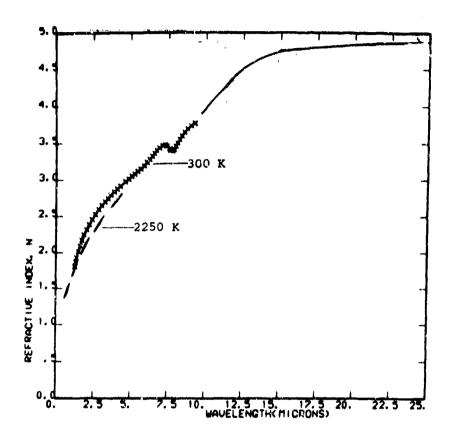


Figure 4-7. n_1 (Carbon) vs. λ for T = 300 and 2250K, based on the Compilation by Whitson and our Extrapolation.

Al₂O₃

Experimental reflectivity measurements in the POPM program at 300 and 678 K were used to determine n_1 and n_2 of AL_2O_3 (Dowling and Randall, 1977). The results are shown in Figure 4-8 and 4-9 and they indicate that n_1 and n_2 are reduced at

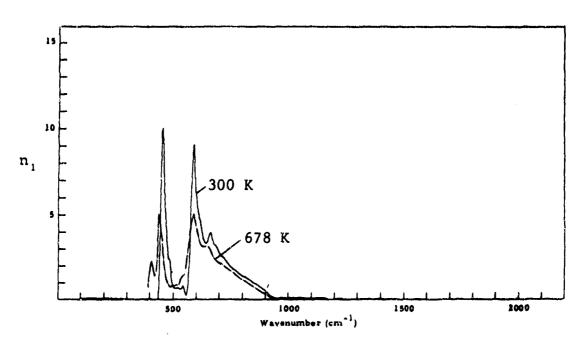


Figure 4-8. The Refractive Index (n_1) of Al_2O_3 in the Region 400 cm⁻¹ to 2200 cm⁻¹ as Determined from a Kramers-Kronig Analysis of Reflectivity Data taken at T=300 and 678 K (Dowling and Randall, 1977).

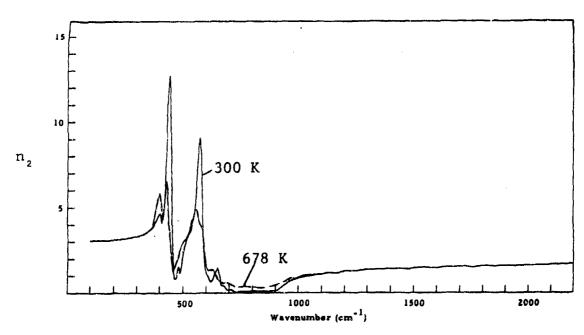


Figure 4-9. The Absorption Index (n_2) of Al_2O_3 in the 400 cm⁻¹ to 2200 cm⁻¹ as Determined from a Kramers-Kronig Analysis of Reflectivity Data taken at T-300 and 678 K (Dowling and Randall, 1977).

higher temperatures in the resonance region. A comparison with the "representative" values by Whitson shows good agreement for the room temperature data. Also the qualitative behavior at higher temperatures in the resonance region has been observed by B. Piriou (1966). The data as presented by Whitson are shown in Figures 4-10 and 4-11.

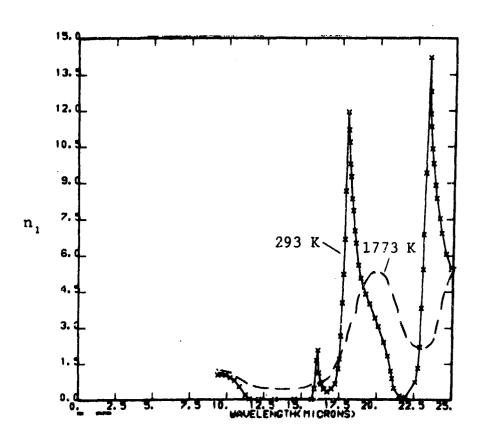


Figure 4-10. $n_1(Al_2O_3)$ as a Function of λ for Two Temperatures as Presented by Whitson (1975).

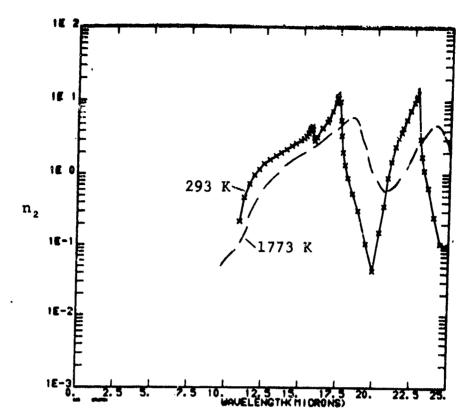


Figure 4-11. $n_2(Al_2O_3)$ as a Function of λ for Two Temperatures as Presented by Whitson (1975).

Mg0

M. E. Whitson (1975) has listed experimental data of n_1 and n_2 taken by several workers. All of these are in general agreement and the "representative" values at room temperature are based on these values. There appears to be little or no dependency on temperature, as the data by B. Piriou (1966) indicate (see Figure 4-12 and 4-13). The experimental study and subsequent data analysis using classical dispersion theory by Dowling and Randall (1977) confirm these data. A comparison of n_1 and n_2 with Whitson's representative room temperature data

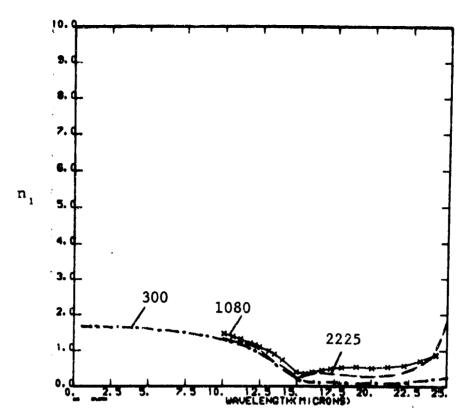


Figure 4-12. n_1 of MgO vs. λ for Three Different Temperatures.

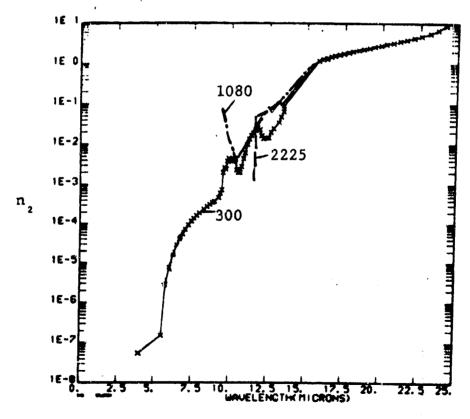


Figure 4-13. $\boldsymbol{n}_{_{2}}$ of MgO vs. λ for Three Different Temperatures.

show good agreement (see Figures 4-14 and 4-15). In addition, Dowling and Randall's data at elevated temperature (573 K) show little or no temperature dependency (see Figures 4-16 and 4-17).

ZrO2

Whitson's report does not contain any data of n_1 and only a limited set of n_2 at room temperature between 5 and 9 μm . Thus, the data obtained by Dowling and Randall filled an important gap. The reflectivity between 400 and 2200 cm⁻¹ was measured at room temperature and 573°K. The spectral structure is similar to that reported by Whitson, albeit lower by about 20%. The results for n_1 and n_2 are shown in Figures 4-18 and 4-19, respectively.

4.4 Numerical Results for Al₂O₃ and Carbon

The volume extinction, scattering and absorption coefficients and the scattering phase function have been calculated for Al_2O_3 and carbon with a code developed by Dave (1972). These two species are the most important particulate constituents in present missile exhausts and possess very different optical and geometrical properties. Thus, the results give insight into the functional dependencies of β and M on size distributions, wavelength and temperature. For Al_2O_3 , the modified gamma function was used for representing the size distribution 2 as given in Figure 4-1, viz.

$$N(r) = N \left[.0551 \text{ y e}^{-.004r^4} \right]$$

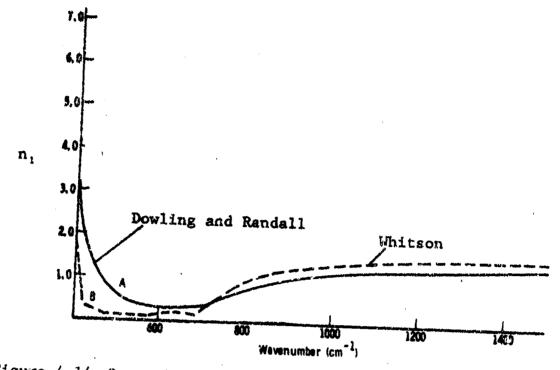


Figure 4-14. Comparison of Index of Refraction (n₁) for MgO.

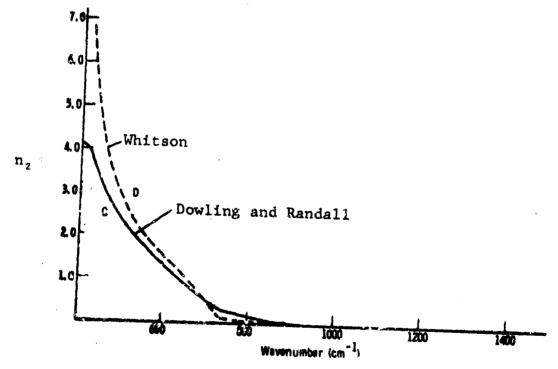


Figure 4-15. Comparison of Absorption Index (n_2) for MgO.

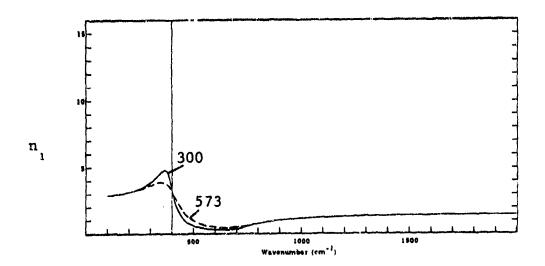


Figure 4-16. The Index of Refraction (n_1) of MgO from 400 cm⁻¹ to 2000 cm⁻¹ as Derived by Classical Dispersion Analysis $(T = 300^{\circ}K)$.

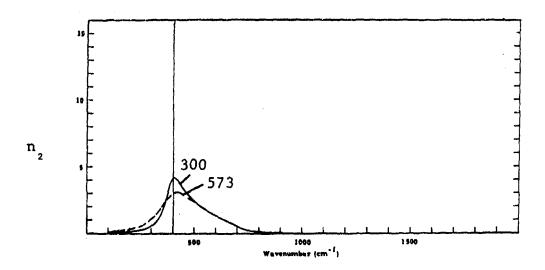


Figure 4-17. The Absorption Index (n_2) of MgO from 400 cm⁻¹ to 2000 cm⁻¹ as Derived by Classical Dispersion Analysis $(T = 300^{\circ}K)$.

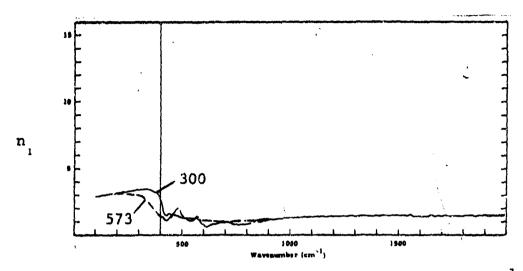


Figure 4-18. The Refractive Index (n₁) of ZrO₂ from 400 cm⁻¹ to 2000 cm⁻² as Determined from a Kramers-Kronig Analysis of the Reflectivity Data.

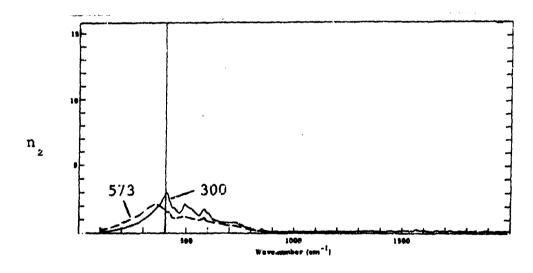


Figure 4-19. The Absorption Index (n₂) of ZrO₂ from 400 cm⁻¹ to 2000 cm⁻¹ as Determined from a Kramers-Kronig Analysis of the Reflectivity Data.

The total number of particles chosen was $10^5 \, \mathrm{cm}^{-3}$ which corresponds to a mole fraction of about 30% when the mean radius is 5 $\mu \mathrm{m}$. The min and max radii chosen were 1 and 9 $\mu \mathrm{m}$, respectively. The input for a_{\min} , a_{\max} , n_1 and n_2 as a function of λ for a temperature of 1773 K are given in Table 4-2.

TABLE 4-2. INPUT PARAMETERS USED FOR AL203

λ(μ m)	a min	a max	n ₁	n ₂
1	6.29	56.55	2	1.5-3
3	2.09	18.85	2	2.0-3
5	1.26	11.31	2	6.0 ⁻³
7	. 90	8.08	1.5	1.8-2
9	. 70	6.28	1.2	8.0-2
11	. 57	5.14	. 5	. 3
13	. 48	4.35	. 3	1.0
15	. 42	3.77	. 3	2.5
17	. 37	3.33	. 5	6.0
19	. 33	2.98	4.0	. 6
21	. 30	2.69	3.5	. 8
23	. 27	2.46	3.0	6.0
25	. 25	2.26	5.0	1.0

The results for the volume extinction, scattering and absorption coefficients and albedo ($\overline{\omega} = \beta_{\rm sca}/\beta_{\rm ext}$) as a function of λ (in steps of 2 µm) are listed in Table 4-3 and the angular scattering functions M(θ) as functions of θ are shown in Figures 4-20 through 4-32. It can be seen that the scattering becomes more and more isotropic as the wavelength increases. This is also shown in Figure 4-33, where the ratio of the side to forward scattering component is plotted vs. λ .

TABLE 4-3. Results for $\beta_{\rm ext}$, $\beta_{\rm sca}$, $\beta_{\rm abs}$ and $\overline{\omega}$ as a Function of λ for Al₂O, Particles in Rocket Plume Exhaust at 1773K.

λ (μ m)	SERE	Beca	⁶ aba	ä
1	2.182	1.961	. 221	. 899
3	2.374	2.237	. 138	. 942
5	2.517	2.259	. 258	. 897
7	2.983	2.550	. 433	. 855
9).641	. 893	. 748	. 544
11	1.#00	1.164	637	. 647
13	2.083	1.877	. 805	. 700
15	3.890	3.331	. 559	. 856
17	2.760	2.637	. 123	. 955
.9	2.848	1.540	1.307	. 541
11	2.941	1.499	1.442	. 510
3	2.760	2.217	. 543	. 803
25	2.808	1.584	1.224	. 564

The dependency of $\beta_{\rm ext}$, $\beta_{\rm sca}$, $\beta_{\rm abs}$ and M(θ) on the temperature is not expected to be strong, since n_1 and n_2 do not show large variations due to the temperature. As will be seen for carbon particles, relatively large variations in n_2 due to temperature produce only moderate variations in the coefficients and negligible ones in the scattering function. Similar results for the efficiency factors of single aluminum particles were also found by Plass (1965).

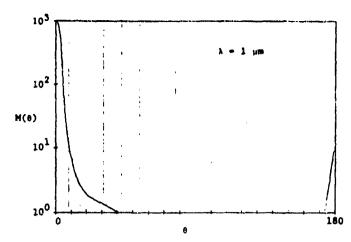


Figure 4-20. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 1 μm .

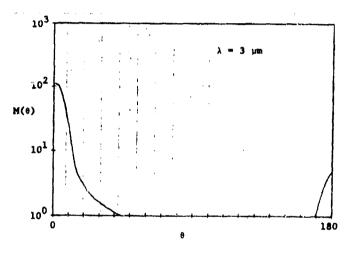


Figure 4-21. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 3 μ m.

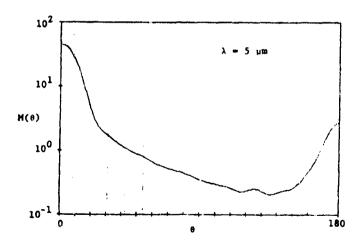


Figure 4-22. Angular Scattering Phase Function $M(\theta)$ for Al_2O_3 at $\lambda = 5 \mu m$.

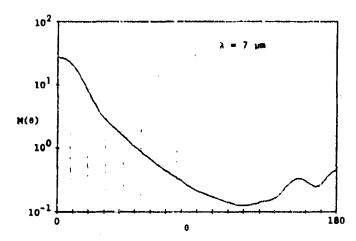


Figure 4-23. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 7 μm .

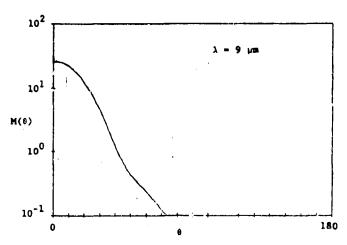


Figure 4-24. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 9 µm.

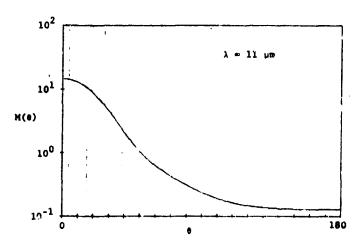


Figure 4-25. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 11 μm .

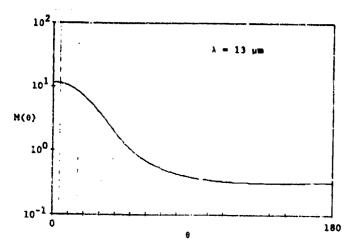


Figure 4-26. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 13 μ m.

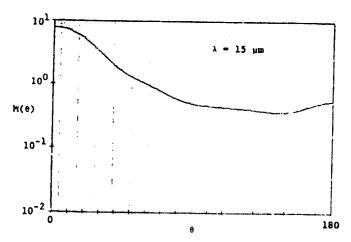


Figure 4-27. Angular Scattering Phase Function M(θ) for Al₂O₃ at λ = 15 μ m.

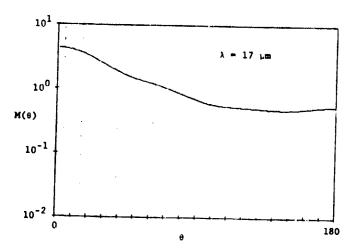


Figure 4-28. Angular Scattering Phase Function $M(\theta)$ for Al_2O_3 at $\lambda=17\,\mu m$.

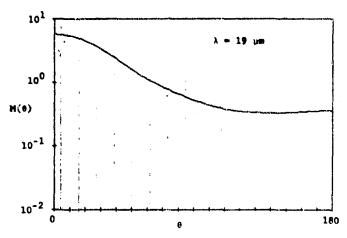


Figure 4-29. Angular Scattering Phase Function M(0) for Al $_2$ O $_3$ at λ = 19 μm .

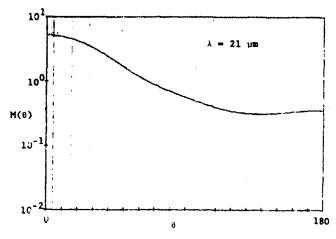


Figure 4-30. Angular Scattering Phase Function M(0) for Al₂O₃ at λ = 21 μm .

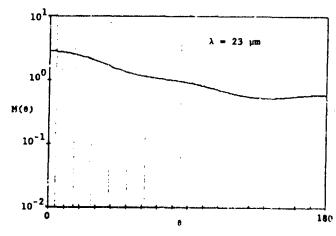


Figure 4-31. Angular Scattering Phase Function M(θ) for At₂O₃ at λ = 23 μ m.

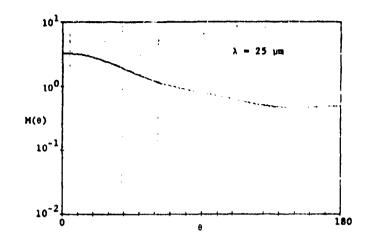


Figure 4-32. Angular Scattering Phase Function $M(\theta)$ for Al_2O_3 at λ = 25 μm .

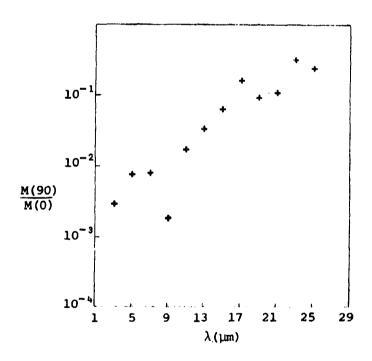


Figure 4-33. Ratio of Side to Forward Scattering Components vs. Wavelength for ${\rm Al}_2{\rm O}_3$

For the second particulate species, the volume coefficient and angular scattering functions have been calculated for carbon. The discontinuous power law was used to represent the size distribution as reported in the literature (Rochelle, 1967), viz.,

$$N(r) = N$$
 for .01 < r < .03 µm

$$N(r) = N(\frac{.03}{r})^2$$
 for .03 < r < 2 µm

The total number of particles was arbitrarily chosen to be $10^{13} \, \text{cm}^{-3}$ which corresponds to about 20% mole fraction ($r_{\text{mean}} \sim .02 \mu \text{m}$). The inputs for a_{min} , a_{max} , n_1 and n_2 are listed in Table 4-4. In order to study the influence

TABLE 4-4. INPUT PARAMETERS FOR CARBON PARTICLES

λ	a _{min}	a max	n ₁	n ₂ (1000K)	n ₂ (2500K)
1	. 013	18.8	1.5	. 3	. 5
3	, 0042	6.25	2.7	. 37	1.1
5	. 0025	3.77	3.0	. 4	1.5
7	.0018	2.69	3.5	. 4	1.9
9	.0014	2.09	3.7	. 4	1.7
11	.0011	1.71	4.1	. 32	1.3
13	. 0010	1.45	4.4	. 29	1.1
.5	. 0009	1.26	4.7	. 23	. 9
١7	. 0007	1.11	4.8	. 21	. 75
.9	. 0006	. 99	4.8	. 18	. 6
1	. 0006	. 90	4.8	. 15	. 5
23	. 0005	. 82	4.8	. 13	. 2
5	. 0005	. 75	4.8	. 12	. 3

of the temperature, n_2 is given for 1000 and 2500K. The results of $\beta_{\rm ext}$, $\beta_{\rm sca}$, $\beta_{\rm abs}$ and $\overline{\omega}$ for 1000 and 2500 K are shown in Tables 4-5 and 4-6. Although the absorption index is higher by a factor of about 3 for the higher temperature throughout the wavelength region from 1 to 25 μm , the parameters $\beta_{\rm ext}$, $\beta_{\rm sca}$, $\beta_{\rm abs}$ and $\overline{\omega}$ are not greatly affected. The influence on the scattering phase function is even less, as can be seen in Figures 4-34 through 4-46.

As for Al_2O_3 , we have used the ratio of the phase function at 90° and 0° to indicate the degree of isotropic scattering of C(s). The ratio 1s plotted in Figure 4-47 and compared with that of Al_2O_3 .

TABLE 4-5. Results for $\beta_{\rm ext}$, $\beta_{\rm sca}$, $\beta_{\rm abs}$ and $\overline{\omega}$ as a Function of λ for C(s) Particles in Rocket Plume Exhaust (T = 1000 K, N = $10^{13}/{\rm cm}^3$.

λ (μπα)	β _{ext}	β _{eca}	⁶ aba	ω
1	1921	934	987	. 486
3	2134	1107	1027	. 519
5	2096	1075	1022	. 513
7	2006	1.039	968	.518
9	1885	962	922	. 510
11	1698	812	886	. 478
13	1578	800	778	. 307
15	1407	717	690	. 510
17	1305	661	643	. 507
19	1308	693	614	. 530
21	1229	746	484	. 607
23	954	588	365	.616
25	819	475	344	. 580

TABLE 4-6. Results for $\beta_{\rm ext}$, $\beta_{\rm sca}$, $\beta_{\rm abs}$ and $\overline{\omega}$ as a Function of λ for C(s) Particles in Rocket Plume Exhaust (T = 2500 K, N = $10^{13}/{\rm cm}^3$).

λ (μm)	\$ _{ext}	Baca	Sabe	<u>.</u>
1	1927	936	992	. 486
3	2137	1106	1031	. 518
5	2095	1082	1013	. 524
7	1.977	1044	932	. 528
9	1844	927	917	. 503
11	1698	812	886	. 478
13	1560	714	846	. 458
15	1415	610	805	. 431
17	1283	520	763	. 405
19	1198	481	717	. 402
21	1090	461	629	. 423
23	913	397	517	. 435
25	777	327	450	. 421

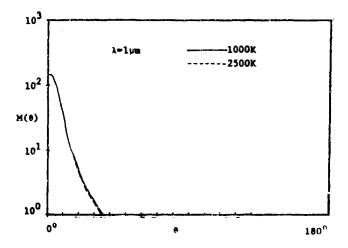


Figure 4-34. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 1 μm .

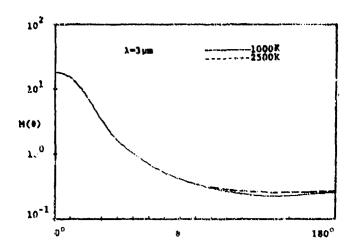


Figure 4-35. Angular Scattering Phase Function $M(\theta)$ for Carbon Particles at λ = 3 μm .

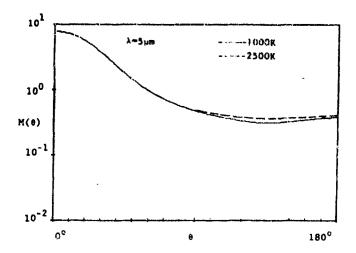


Figure 4-36. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 5 μm .

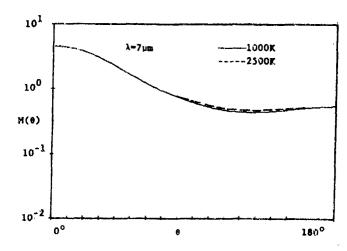


Figure 4-37. Angular Scattering Phase Function $M(\theta)$ for Carbon Particles at λ = 7 μm .

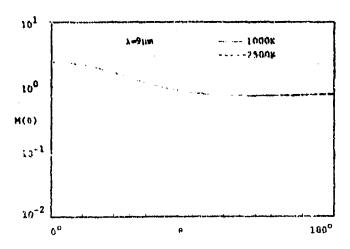


Figure 4-38. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 9 μm .

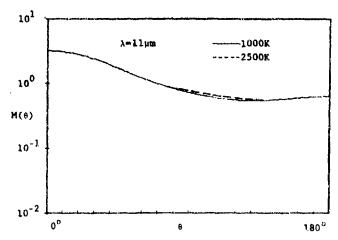


Figure 4-39. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 11 μm .

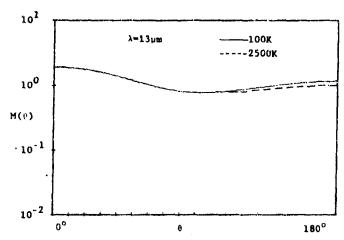


Figure 4-40. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 13 μm .

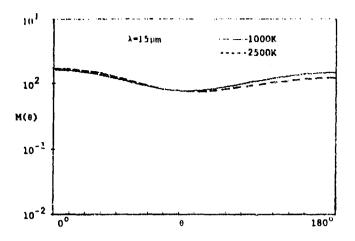


Figure 4-41. Angular Scattering Phase Function M(0) for Carbon Particles at $\lambda = 15 \, \mu m$.

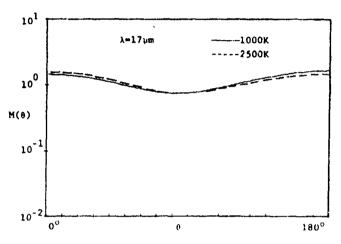


Figure 4-42. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 17 μ m.

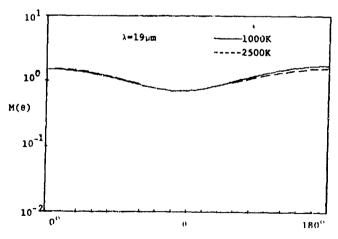


Figure 4-43. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 19 μm .

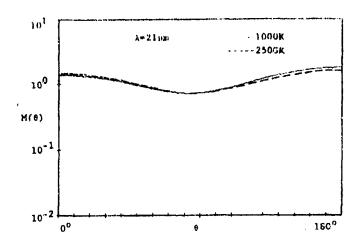


Figure 4-44. Angular Scattering Phase Function M(θ) for Carbon Farticles at λ = 21 μ m.

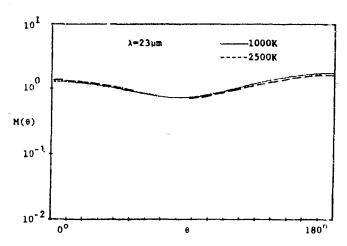


Figure 4-45. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 23 μ m.

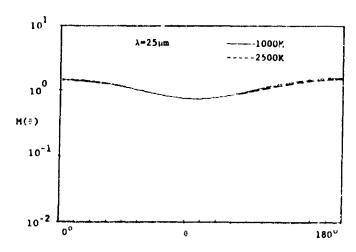


Figure 4-46. Angular Scattering Phase Function M(θ) for Carbon Particles at λ = 25 μ m.

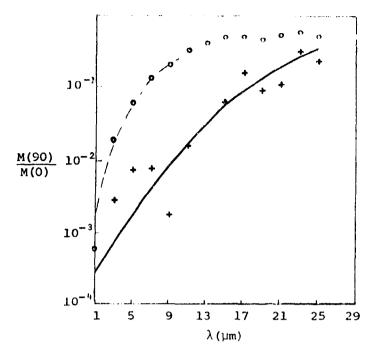


Figure 4-47. The Ratio of the Phase Function at $90^{\rm o}$ and $0^{\rm o}$ Scattering Angle as a Function of λ_a for Al₂0₃ (+ — +) and C(s) (o --- o).

4.5 Recommended Data File and Update Procedure

Based on the studies of the scattering properties of particles performed during Subtask IA it is recommended that SIRRM shall include

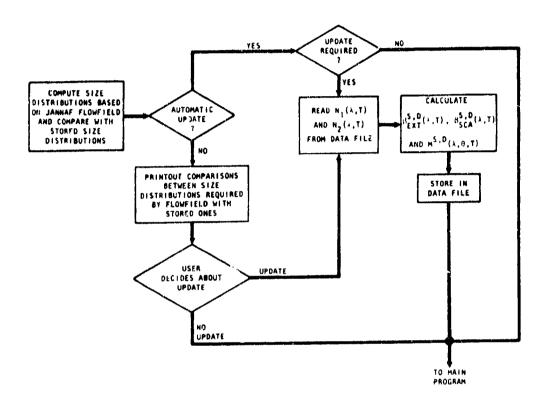
(1) a permanent Data File of $n_{(\lambda,T)}^{s}$, $n_{(\lambda,T)}^{s}$, for each species s and for each size distribution d; (this Data File may be minimized to as few temperatures and size distributions as indicated by experience);

- (2) a subroutine that allows the user to update this Data File or add to it as new optical parameters $(n_1(\lambda,T), n_2(\lambda,T))$ and size distributions) become available;
- (3) a subroutine that compares the size distribution as calculated by the JANNAF flow field with those stored in the Data File and that will indicate to the user whether or not the stored size distribution(s) is (are) sufficient. We consider also the option of an automatic updating.

The advantages of the recommended approach are the following:

- Computer time is minimized by using a stored Data File;
- Storage requirements are minimized by using as few size distributions for each species as possible;
- Flexibility is maintained by allowing the user to update the Data File at his discretion;
- Indication is given to the user before execution whether an update of the Data File in terms of new and/or additional size distributions is necessary.

A tentative logic flow diagram of the recommended approach is shown in Figure 4-48. After the JANNAF flow field is read in and the relevant flow field positions for the line-of-sight calculations are determined, a subroutine calculates the analytical expression for the size distributions for each position. This is necessary since the JANNAF flow field does not give the analytical expression but does provide the number density for several specific particle sizes (private communications, S. Dash, May 1979). Different analytical expressions will be used.



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Figure 4-48. Logic Flow Diagram of Data File Update Procedure

such as the discontinuous power law for particles with very small radii (carbon) and modified gamma function for particles with larger radii ($\Delta \ell_0 0$, etc.)

The computed size distributions will be compared with those already stored in the Data File. If the differences are minor (criteria must yet be established) no update is required. If the differences are major, $\beta_{\rm ext}$, $\beta_{\rm sca}$ and M(0) may have to be calculated and stored in the Data File. If the decision for an update is desired to be automatic (probably a default input), SIRRM will read the appropriate n and n and will calculate $\beta_{\rm ext}$, $\beta_{\rm sca}$ and M(0). If the decision is not to be automatic, the user has the option whether or not new values of $\beta_{\rm ext}$, $\beta_{\rm sca}$ and M(0) should be calculated. As more experience is gained about the influence of different size distributions on the radiation field using the six-flux and N-flux models, the decision about updating the Data File may become straightforward.

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