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# Multidimensional material response simulations of a full-scale tiled ablative heatshield

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### 1. Introduction

The Mars Science Laboratory (MSL) spacecraft, launched on November 2011, successfully landed the Mars Curiosity rover in the Aeolis Palus region of the Gale Crater on August 2012. The MSL entry vehicle was equipped with a 4.5 m diameter Thermal Protection System (TPS) that effectively protected the spacecraft and its payload during entry into Mars' atmosphere. The MSL TPS used the Phenolic Impregnated Carbon Ablator, or PICA, as heatshield material [1]. PICA is a low density ( $\approx$ 274 kg/m<sup>3</sup>) carbon/resin composite, manufactured via impregnation of a rigid carbon fiber preform (FiberForm) with a phenolic resin (Durite<sup>®</sup> SC-1008), followed by a proprietary high temperature curing and

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### ABSTRACT

The Mars Science Laboratory (MSL) was protected during Mars atmospheric entry by a 4.5 meter diameter heatshield, which was constructed by assembling 113 thermal tiles made of NASA's flagship porous ablative material, Phenolic Impregnated Carbon Ablator (PICA). Analysis and certification of the tiles thickness were based on a one-dimensional model of the PICA response to the entry aerothermal environment. This work provides a detailed three-dimensional heat and mass transfer analysis of the full-scale MSL tiled heatshield. One-dimensional and three-dimensional material response models are compared at different locations of the heatshield. The three-dimensional analysis is made possible by the use of the Porous material Analysis Toolbox based on OpenFOAM (PATO) to simulate the material response. PATO solves the conservation equations of solid mass, gas mass, gas momentum and total energy, using a volume-averaged formulation that includes production of gases from the decomposition of polymeric matrix. Boundary conditions at the heatshield forebody surface were interpolated in time and space from the aerothermal environment computed with the Data Parallel Line Relaxation (DPLR) code at discrete points of the MSL trajectory. A mesh consisting of two million cells was created in Pointwise, and the material response was performed using 840 processors on NASA's Pleiades supercomputer. The present work constitutes the first demonstration of a three-dimensional material response simulation of a full-scale ablative heatshield with tiled interfaces. It is found that three-dimensional effects are pronounced at the heatshield outer flank, where maximum heating and heat loads occur for laminar flows.

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vacuum drying process [2]. The material was successfully used on the Stardust Sample Return Capsule (SRC), assembled in a 0.8 m diameter monolithic aeroshell [3]. Due to manufacturing constraints, it was unfeasible to construct a 4.5 m diameter heatshield out of a single piece of PICA. Instead, the MSL heatshield was developed as an assembly of 113 PICA tiles containing 23 unique shapes. There were also gaps between the TPS tiles to allow for thermal expansion and contraction. These gaps were filled using a silicone elastomer bonding agent. The MSL heatshield was instrumented with temperature and pressure sensors; therefore, the MSL is an established validation case for ablator response models. The MEDLI (MSL Entry, Descent, and Landing Instrument) suite recorded, among others, time-resolved in-depth temperature data using thermocouple sensors assembled in the MEDLI Integrated Sensor Plugs (MISP). Several studies in the literature have used MISP data as a benchmark for state-of-the-art ablation codes [4–6]. Modeling of heat and mass transfer in porous materials during at-

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### Nomenclature

$\mathcal{A}_{i,j}$	Arrhenius law pre-exponential factor $K^{-n_{i,j}} s^{-1}$	Xi, j	Advancement of pyrolysis reaction $j$ within phase $i$
$\mathcal{E}_{i,j}$	Arrhenius law activation energy $\dots$ J mol <sup>-1</sup>	$\epsilon$	Porosity
$\mathcal{R}^{+}$	Perfect gas constant Jmol <sup>-1</sup> K <sup>-1</sup>	$\epsilon_{i,0}$	Initial porosity of phase <i>i</i>
$\dot{m}_{ca}$	Char ablation rate $\dots$ kg m <sup>-2</sup> s <sup>-1</sup>	λ	Scaling factor for $C'_H$
$\dot{m}_{pg}$	Pyrolysis gas production rate $\dots$ kg m <sup>-2</sup> s <sup>-1</sup>	$\mu$	Viscosity kg m <sup>-1</sup> s <sup>-1</sup>
St <sub>H</sub>	Stanton number for heat transfer	$v_{i,k}$	Number of atoms of element <i>k</i> in molecule of specie <i>i</i>
$St_M$	Stanton number for mass transfer	П	Total pyrolysis gas production rate $kg m^{-3} s^{-1}$
<u>K</u>	Permeability tensor m <sup>2</sup>	$\pi_k$	Pyrolysis gas production rate of specie k kg m <sup>-3</sup> s <sup>-1</sup>
k	Effective thermal conductivity tensor $W m^{-1} K^{-1}$	ρ	Mass density kg m <sup>-3</sup>
n	Front surface normal	$\rho_{i,0}$	Initial mass density of phase $i$ kg m <sup>-3</sup>
νσ	Gas velocity m s <sup>-1</sup>	σ	Stefan-Boltzmann constant,
V <sub>ca</sub>	Char ablation velocity ms <sup>-1</sup>		$5.670367 \times 10^{-8}$ W m <sup>-2</sup> K <sup>-4</sup>
$A_k$	Element <i>k</i>	β	Klinkenberg correction tensor m <sup>2</sup> s <sup>-1</sup>
B'	Dimensionless mass blowing rate	E	Fmissivity
$C'_{II}$	Corrected heat transfer coefficient $kg m^{-2} s^{-1}$	۲ ۲	Mass stoichiometric coefficient
Сн	Heat transfer coefficient = $\rho_e u_e St_H \dots kg m^{-2} s^{-1}$	, 	
См	Mass transfer coefficient = $\rho_e u_e St_M \dots kg m^{-2} s^{-1}$	Subscripts	
Cn	Specific heat capacity	0	Initial
e	Specific energy Ikg <sup>-1</sup>	$\infty$	Infinity
F <sub>i</sub> i	Fraction of subphase <i>i</i> in phase <i>i</i>	а	Ablation
h	Specific absolute enthalpy	adv	Advection
Ki	Equilibrium constant	С	Char
m <sub>i, i</sub>	Arrhenius law advancement pyrolysis reaction factor	cond	Conduction
Ns	Number of gaseous species	conv	Convection
Np	Number of solid phases	diff	Diffusion
Ňe	Number of gaseous elements	е	Boundary layer edge
n <sub>i.i</sub>	Arrhenius law temperature factor	flux	Corrected convection
р	Pressure Pa	g	Gas phase
Pi	Number of subphases in solid phase <i>i</i>	р	Pyrolysis reaction
q	Heat flux $\dots$ W m <sup>-2</sup>	pla	Plasma
Si	Specie i	rad	Radiation
Т	Temperature K	S	Solid phase
$x_i$	Mole fraction of specie <i>i</i>	t	Total (solid and gas phases)
$x_k$	Mole fraction of element <i>k</i>	ν	Virgin
Уi	Mass fraction of specie <i>i</i>	w	Wall
$z_k$	Mass fraction of element k	Suhnerse	crints
Greek			
GIECK		in	Inside the material
α	Absorptivity	out	Outside the material

mospheric entry of spacecrafts is a complex and computationally expensive problem. Traditionally, NASA TPS design has been done using one-dimensional ablation and thermal response solvers [7,8]. Research by Chen and Milos [9,10] investigated multi-dimensional effects on the thermal response of a monolithic Apollo-shaped heatshield, using the 3dFIAT code developed at the NASA Ames Research Center. The convective aerothermal environment over the exterior TPS surface were derived from the Configuration Based Aerodynamics (CBAERO) analysis of Lunar return trajectories [11]. The analysis indicated that, for a high angle of attack entry, the peak heat flux and heat load are located at the windside heatshield outer flank. At this location, the planar approximation was shown to underpredict the peak bondline temperature. In this study, de-tailed three-dimensional heat and mass transfer analyses are un-dertaken on the tiled MSL heatshield to assess the validity and limitations of the traditional one-dimensional design assumption. The Porous material Analysis Toolbox based on OpenFOAM (PATO) software program is used to simulate the TPS response [12]. PATO is released as open source software by NASA.<sup>1</sup> The aerothermal 

environment at the heatshield surface, at discrete points along the MSL entry trajectory, is obtained from hypersonic Computational Fluid Dynamics (CFD) simulations performed using the Data Parallel Line Relaxation (DPLR)<sup>2</sup> Navier–Stokes software program [13]. A procedure for temporal and spatial interpolation was used to loosely couple the aerothermal environment to the material response. Simulations were performed for both monolithic and tiled heatshield configurations. Tiles are normally not included in 3D material response analysis because of the high computational costs associated with such simulations. The massively parallel simulation support inherited in PATO from its OpenFOAM architecture now makes such studies possible.

The paper is organized as follows. In section 2, we describe the MSL monolithic and tiled heatshield grid used for the simulations. Section 3 presents the governing equations used in the material response model. Section 4 details the spatial and temporal interpolation from DPLR to PATO. In section 5, the overall material response is presented where monolithic and tiles configurations are compared. Detailed heat and mass transfer analyses

#### <sup>1</sup> https://software.nasa.gov/software/ARC-16680-1A.

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