

Space Shuttle Debris Impact Tool Assessment Using the Modern Design of Experiments

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Complex computer codes are used to estimate thermal and structural reentry loads on the Shuttle Orbiter induced by ice and foam debris impact during ascent. Such debris can create cavities in the Shuttle Thermal Protection System. The sizes and shapes of these cavities are approximated to accommodate a code limitation that requires simple “shoebox” geometries to describe the cavities—rectangular areas and planar walls that are at constant angles with respect to vertical. These approximations induce uncertainty in the code results. The Modern Design of Experiments (MDOE) has recently been applied to develop a series of resource-minimal computational experiments designed to generate low-order polynomial graduating functions to approximate the more complex underlying codes. These polynomial functions were then used to propagate cavity geometry errors to estimate the uncertainty they induce in the reentry load calculations performed by the underlying code. This paper describes a methodological study focused on evaluating the application of MDOE to future operational codes in a rapid and low-cost way to assess the effects of cavity geometry uncertainty.

Nomenclature

ANOVA	=	ANalysis Of Variance
BLT	=	Boundary Layer Transition
BPZ	=	Body Point Zone, a sub-area on the windward side of the Orbiter
CPU	=	Central Processing Unit
CV	=	Coefficient of Variation. Ratio of standard error to mean response
ET	=	External Tank
FS	=	Factor of Safety

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Inference space	=	A coordinate system in which each axis corresponds to an independent variable
ITA	=	Independent Technical Authority
JSC	=	Johnson Space Center
LaRC	=	Langley Research Center
LOF	=	Lack of Fit
MDOE	=	Modern Design of Experiments
MMT	=	Mission Management Team
MS	=	Margin of Safety
NESC	=	NASA Engineering and Safety Center
OBSS	=	Orbiter Boom Sensor System
OFAT	=	One Factor At a Time
OOPD	=	Out Of Plane Deflection
RCC	=	Reinforced Carbon-Carbon
RSM	=	Response Surface Modeling/Methods
RTF	=	Return to Flight
RTV	=	Room Temperature Vulcanized
SIP	=	Strain Isolation Pad
Site	=	A point in an inference space representing a unique combination of independent variables
TAEM	=	Terminal Area Energy Management
TMM	=	Thermal Math Model
TPS	=	Thermal Protection System
USA	=	United Space Alliance

I. Introduction

AS part of the Space Shuttle Return-to-Flight (RTF) efforts, the ability to perform assessment of damage to the Shuttle Thermal Protection System (TPS), tiles and reinforced carbon-carbon (RCC), became vital to clearing the vehicle for flight readiness and atmospheric entry. In order to accomplish this, a considerable amount of testing and analysis was performed to determine whether a given TPS damage was considered to be survivable. A combination of new and existing math model tools was developed to determine the impact tolerance and damage tolerance of the TPS due to impacts from debris, primarily ice and foam from the external tank (ET). These math model tools include damage prediction, aeroheating analysis, thermal analysis and stress analysis. Some tools are physics-based and others are empirically based. Each tool was created for a specific use and timeframe, such as, probability risk assessment, and real-time on-orbit assessment. In addition, the tools are used together in an integrated strategy for assessing impact damage to tile and RCC.

A peer-review of the tools to assess the engineering readiness for supporting RTF was initiated. The objective of the peer review was to determine if the tools and the end-to-end strategy were suitable to support RTF. As part of the effort to fully understand the limitations and sensitivities of the analytical tools, a task jointly funded by the NASA Engineering and Safety Center (NESC) and the Independent Technical Authority (ITA) was initiated as a collaboration of the NASA Johnson Space Center, the NASA Langley Research Center, and Jacobs Technology team members. The scope of this assessment was limited to tile damage.

Utilizing these tools requires knowledge of a variety of input parameters. As with any analysis, an understanding of the sensitivity of the model response to the variation of the input parameters is essential. Additionally, since the end-to-end analysis involves a variety of analysis steps utilizing multiple models and crossing disciplines, an assessment of end-to-end uncertainty in the analysis was deemed necessary.

The suitability of a Modern Design of Experiments (MDOE) approach was investigated for assessing the analytical models' sensitivity and end-to-end uncertainty. These analyses involve a large number of input variables. To fully investigate all possible combinations of the parameters deterministically or by Monte Carlo sampling was deemed infeasible. The MDOE approach showed promise in making the assessments more efficiently.

II. Analytical Model Overview

Three-dimensional TPS damage thermal math models were developed to aid in the assessment of damage sustained by the Orbiter tile during ascent. On-orbit inspection of the vehicle TPS by the Orbiter Boom Sensor System (OBSS) provides a point cloud geometry of the damage that is subsequently used to develop simplified representations of tile damage for use in the thermal model.

The thermal response of the damaged region during atmospheric entry is assessed to predict tile and structural temperatures in and around the damaged area. These temperatures are then superimposed with the entry mechanical loads on a structural model to determine the structural integrity of the tile and vehicle structure. The results of this analysis will be used to aid the determination by the Mission Management Team (MMT) of whether an observed damage is acceptable to “fly as is,” or a repair will be attempted before reentry, or that safe reentry is not an option in the damaged state and the Space Shuttle crew must remain on the International Space Station until a rescue mission can be launched. Because of the dire consequences of an ill-informed decision, understanding the uncertainty in this analysis is vital to the MMT.

Three analytical tools make up the core of the damage assessment modeling capability developed by the United Space Alliance and the Boeing Company team for the NASA Orbiter Project Office:

- 1) The 3-D Thermal Math Model (3-D TMM) tool predicts the temperature response of the Orbiter lower surface acreage TPS materials and structure at damaged tile locations, using simplified tile damage geometries, structure definition, tile thickness, and material thermal properties. The tool predicts time histories of tile and structure temperatures, and structural thermal gradients.
- 2) The RTV Bondline Tool calculates stresses for the Orbiter lower surface 6 in. × 6 in. acreage tiles at the tile-to-strain isolation pad (SIP) interface using tile damage dimensions, tile bondline temperature data from the 3-D TMM tool and out-of plane deflections (OOPD) from the stress assessor tool. The tool defines factors of safety at the tile/strain isolator pad (SIP) interface of the damaged tile for tension in the RTV bond or tile, compression and shear of the tile, and shear in the SIP.
- 3) The Stress Assessor Tool converts damage site local temperature increases into thermal loads. The tool combines these increased thermal loads with existing mechanical loads and analyzes the new set of combined loads using existing documented flight certification analysis programs to define Orbiter structural margins of safety and Orbiter surface out-of-plane deflections in the area near a tile damage site or tile repair site.

The tool parameter inputs describing the idealized tile damage cavity (an inverted rectangular pyramid geometry) are the length, width, and depth of the damage as measured at the top of the tile surface, and the angles of the entrance, side, and exit faces. Length, width, depth, entrance angle, and side angle were the five independent cavity geometry variables selected to define the polynomial response models for this study, with the decision to hold exit angle fixed at 75°. This decision was predicated on preliminary results that revealed relatively little impact of exit angle changes on computed reentry loads, at least over the range of 70° to 80°, which prior experience suggests is typical for exit angles.

In addition to the five independent geometry variables, two Shuttle state variables were identified. The first, boundary layer transition time, serves as a surrogate for surface roughness and was defined as a categorical variable with two fixed levels, 900 seconds and 1200 seconds, selected to bracket the anticipated range of this variable. The second, baseline heating, was likewise established as a two-level categorical variable, representing low and moderate baseline heating conditions.

Tool output parameters of interest are tile bondline temperature, structural temperature, tile factor of safety, and structural margin of safety. In addition, tile out of plane deflections for two regimes of the reentry trajectory were also examined. The first regime was from Mach 25 to Mach 4, and the second regime was the Terminal Area Energy Management (TAEM) regime, from Mach 4 to touchdown.

III. Overview of Modern Design of Experiments

The Modern Design of Experiments (MDOE) is a formal method of empirical investigation first introduced at Langley Research Center in 1997 to enhance quality and productivity in large-scale, high-cost experiments, initially wind tunnel experiments.¹ It consists of integrated design, execution, and analysis elements intended to extend conventional industrial design of experiments technology (generally used for process and product improvement) to cover the special circumstances of aerospace research. Such circumstances include a very low tolerance for unexplained variance in experimental data, coupled with generally complex relationships between system response variables of interest and the independent variables upon which they depend. The MDOE method is a general approach to experimentation, and while initially motivated by productivity and quality issues in wind tunnel testing, MDOE applications have rapidly expanded to include a wide range of aerospace disciplines, including computational experiments.

The mechanism by which MDOE improves both quality and productivity is to alter the focus of experimentation from a traditional emphasis on data acquisition to an explicit concentration on the extension of knowledge. While data acquisition rate and resulting data volume have been traditional measures of productivity in large-scale

aerospace testing, MDOE practitioners recognize that direct operating costs and cycle time are proportional to the volume of data acquired and therefore seek to organize experimental investigations around the smallest volume of data that is still ample to achieve technical objectives.² Likewise, quality—traditionally assessed in terms of variability in the data—is regarded differently in an MDOE experiment.^{3,4} Low variability in the raw data is not sufficient to assert high quality in an MDOE experiment. From the MDOE perspective, quality is assessed in the context of minimized inference error risk. There must be a demonstrably high probability that the conclusions of the study are correct, which places a strong emphasis on uncertainty assessment. This focus on quantifying uncertainty, coupled with an emphasis on minimizing data volume requirements, made the MDOE method a likely candidate for evaluating the debris impact assessment models developed in support of RTF activities.

IV. Preliminary MDOE Screening Experiment Designs

Four sources of uncertainty in estimating the thermal and structural reentry loads were identified. These are

- 1) **Model verification errors:** Errors due to improper modeling of the underlying behavior of the system. These errors occur when the relationship between dependent system response variables and independent input variables is imperfectly understood. The computer code may accurately represent the mechanism that is programmed, but verification errors will occur if that mechanism is incorrect. Verification tests for modeling errors rather than coding errors.
- 2) **Model validation errors:** Errors due to “bugs” in the software that cause improper results of calculations even when the relationships that are specified for coding represent the true underlying behavior of the system under study. These are coding errors, not modeling errors.
- 3) **Errors due to independent variable uncertainty:** Even if the code is validated and verified, its output can only be as good as its input. The TPS cavity geometry variables are imperfectly known. This is due to ordinary experimental error in measuring them, and also because they are modeled with idealized rectangular “shoebox” geometries that are only approximations to the true geometry. Uncertainty in the independent variable specifications translates into uncertainty in the model response predictions.
- 4) **Lack of Fit (LOF) errors in estimating response surface models (RSM):** In order to facilitate independent variable error propagation, it is convenient to represent the complex reentry load models with simpler polynomial response functions. This simplification introduces a source of error known as lack of fit error.

While MDOE methods could be used to develop low-cost, high-quality model validation and verification experiments, validation and verification considerations⁵ were beyond the scope of the current activity. The computer codes developed by the United Space Alliance and the Boeing Company team for the NASA Orbiter Project Office were assumed to have been sufficiently validated and verified that reentry load estimates made with those codes adequately represented the true loads that would be experienced by the Orbiter on reentry. That is, validation and verification errors were assumed to be negligible. The current effort focused on the propagation of input variable uncertainty, including considerations of LOF errors associated with polynomial models that were developed to facilitate the error propagation calculations.

MDOE computational experiments were designed for each of 33 body point zones on the Orbiter’s windward side. These zones are shown in Fig. 1. Initially, a series of two-level full factorial screening experiments were designed, executed, and analyzed to rank order independent variables by their correlation with response variables of interest, in order to identify variables with sufficiently small influence on the responses to justify dropping them from further consideration. Eliminating irrelevant factors from detailed consideration is an important quality and productivity step. It is important for productivity because the cases that must be run to fit a given order of polynomial grows rapidly with the number of independent variables in the model, with attendant growth in computational time and costs. A full d^{th} -order polynomial model in k independent variables features p parameters, where

$$p = \frac{(d + k)!}{d!k!} \quad (1)$$

The test matrix must include at least one data point for each parameter in the model, and generally some additional degrees of freedom are desirable beyond this minimum to allow an assessment of uncertainty in addition to fitting the data. For a full third-order polynomial function in 10 variables as originally contemplated for this study $d = 3$, $k = 10$, and by Eq. (1), $p = 286$. Since two of the variables were two-level categorical factors for which quadratic and cubic terms are not necessary, there were actually only 269 points required to fit a third-order model

in the original 10 variables. With 20 residual degrees of freedom specified to assess uncertainty, and 15 additional confirmation points, the original planned scale of the experiment was $269 + 20 + 15 = 304$ points.

The preliminary screening experiments justified the elimination of three variables initially identified as candidates for more detailed investigation. The cavity exit angle was eliminated because the screening experiment revealed that over the relatively narrow interval that experience dictated this variable ranges in practice (70° to 80°), the regression coefficient was not large compared to the uncertainty in estimating it. It was therefore not possible to resolve a difference between this coefficient and zero with the level of confidence (99%) adopted as criterion in this experiment. A constant exit angle of 75° was therefore modeled, with no substantive loss in predictive power.

Lateral and longitudinal coordinates were eliminated after anomalous behavior in the screening experiment lead to further investigations that showed it was not possible to specify arbitrary cavity location coordinates due to the discrete distribution of Shuttle body points. In an experiment design featuring X and Y cavity centroid coordinates, the underlying code computed responses for the nearest Shuttle body point and assigned them to the coordinate specified in the test matrix, notwithstanding the fact that the specified coordinates were not the coordinates for which the response was computed. Since this would produce a false relationship between location coordinates and response predictions, the detailed investigations were conducted for a specified, representative body point within each zone. The option was reserved to model multiple body points within a zone if within-zone spatial response variation was later deemed to be of special interest.

A full third-order polynomial function in seven variables instead of 10 as originally contemplated for this study requires a minimum of 120 points by Eq. (1). Since again two of the variables were two-level categorical factors for which quadratic and cubic terms are not necessary, only 104 points are required to fit a third-order model in the retained seven variables. Twenty residual degrees of freedom were added to assess uncertainty, so 124 points were specified to fit the response surface models. There were 15 additional confirmation points for a total of 139 points that were specified for each body point zone, down from the 304 points required for 10 variables. The savings of 165 case runs is realized for each of the 33 body point zones, a total of 5,445 cases that did not have to be run. Each case required an average of approximately 45 minutes to run, with eight such cases typically run in parallel. The savings of 5,445 cases therefore translates into more than 500 hours of CPU time that were saved. The actual wall-clock savings were considerably greater, since the 45 minutes per run does not include the considerable time required to prepare inputs for each run and other activities that consume time in addition to the CPU time of the case runs.

Eliminating candidate variables in a preliminary screening experiment can reduce the computer resources required in a computational experiment, as shown. It can also improve quality, by reducing the uncertainty in predictions made by response surface models developed from such experiments.

The prediction variance of a polynomial regression model varies from one combination of independent variables to another, but across all n points used to fit the model, the average prediction variance is as follows⁶

$$\frac{\sum_{i=1}^n \text{Var}(\hat{y}_i)}{n} = \frac{p\sigma^2}{n} \quad (2)$$

where p is the number of terms in the model and σ represents the standard error in the response variable. This result holds for any model that is linear in the coefficients, including polynomial models of any order. Note that while σ cannot be empirically determined from replicated data in a computational experiment as it can be in a physical experiment, the response estimates have uncertainty nonetheless, and σ is a measure of that uncertainty, however it is determined in practice.

Equation (2) indicates that the average prediction variance increases linearly with the number of terms in the model, so there is potential to reduce the average prediction variance simply by reducing the number of model terms. This is because each term carries with it some contribution to the total uncertainty in the model prediction. As Eq. (1) shows, reducing the number of variables to be fitted with a polynomial response function results in a rapid reduction in the number of terms required for the model, and thus a rapid reduction in prediction uncertainty for models developed by fitting a given number of data points. Alternatively, reducing the number of independent variables can lead to the same prediction uncertainty with substantially fewer data points used to fit the response model.

Based on the screening experiment results, an experiment was designed to acquire data to fit a third-order response model in seven independent variables. These variables are noted in Table 1.

Table 1. Independent variables in response surface modeling experiments.

Variable	Units	Low	High
Cavity Length	in.	2	10
Cavity Width	in.	1.4	4
Cavity Depth	in.	0.1	0.6
Entry Angle	deg.	10	60
Sidewall Angle	deg.	10	87
Baseline Heating		Low	Nominal
BLT Time	sec.	900	1200

V. MDOE Experiment Designs for Response Surface Modeling

MDOE response surface experiments were designed for each of 33 body point zones in Fig. 1. A case matrix was developed for each of the 33 experiments, consisting of 124 unique combinations of the seven independent variables in Table 1. The USA/Boeing computer codes were used to generate responses previously noted for each of these variable combinations, plus 15 more that were held in reserve as confirmation points to test the models.

To develop the test matrix for an MDOE experiment, the total number of points must be determined as described above, and then specific points must be selected. While not generally taken into account in the design of test matrices for aerospace research, the *selection* of points (not just the *number* of points) is an important determinant of the quality of experimental and computational research results. Figure 2 illustrates the concept for a simple first-order function of one variable.

For an experiment as simple as the one illustrated in Fig. 2, the order of the model, d , is 1 and the number of variables, k , is 1, so that Eq. (1) reminds us of the familiar fact that two points are required to fit a straight line. Figure 2 illustrates how the uncertainty in the coefficient estimates for this model—the slope of the straight line and the intercept—depend on the selection of two points used to fit the straight line. In this case the points that are farther apart give a more satisfactory result than points that are closer together. This is because each point has some uncertainty associated with it, and this uncertainty allows a range of slope and y intercepts to be consistent with the data. This range is more restricted for one data set than another in Fig. 2.

The term “inference space” is used to denote a k -dimensional space in which each axis corresponds to one of the independent variables. Each site in the inference space represents a unique combination of independent variable levels. Figure 2 illustrates the effect of site selection in a simple one-dimensional inference space. The phenomenon illustrated in this figure for a simple first-order function of one variable is completely general. It applies to RSM experiment designs no matter how complex the response function or how many variables are in the model. Much of the quality enhancement of MDOE RSM experiment designs is achieved by optimizing site selections in different ways to optimize results by various criteria.

For the RSM experiments developed in this study, a D-optimal design strategy was implemented, in which the points are chosen to minimize the variance associated with the estimates of model coefficients. Standard references describe the D-optimal algorithm in detail,^{7,8} but modifications to the method are discussed here that were needed to accommodate certain constraints among the independent variables in this problem.

Consider first the design matrix, \mathbf{X} . The design matrix is an extension of the test matrix. It has rows for each point in the design just as a conventional test matrix does, but it also has columns for every term in the response model, including the intercept term. The design matrix for this experiment therefore has 104 columns and 124 rows.

The covariance matrix, \mathbf{C} , is a $(p \times p)$ square matrix computed by pre-multiplying the design matrix by its transpose, inverting the product, and multiplying each element of the resulting matrix by the unexplained variance of the residuals, σ^2 :

$$\mathbf{C} = (\mathbf{X}'\mathbf{X})^{-1} \sigma^2 \quad (3)$$

It can be shown that the diagonal elements of the covariance matrix represent the variance in estimates of the regression coefficients. That is, the variance in the i^{th} regression coefficient, β_i is simply:

$$\text{Var}(\beta_i) = \mathbf{C}_{ii} \quad (4)$$

To minimize the uncertainty in the regression coefficients, we must minimize the determinant of the covariance matrix. This minimizes the confidence *ellipsoids* for each regression coefficient, the multidimensional equivalent of minimizing confidence *intervals* about a measured data point. The D-optimal design algorithm selects the subset of points to do this from a candidate points list.

A. D-Optimal Candidate Point List

A candidate point list consists of a large number of potential points (typically thousands for an experiment of this complexity), that satisfy certain criteria. Specialized software routines identify the candidate points automatically. One such algorithm⁹ starts with a point selected at random from within the inference space. Subsequent points are picked at random and retained if they increase the rank of the design matrix corresponding to a proposed regression model that is specified at the beginning of this process. That is, points are retained if they do not feature linear relationships among the columns of the design matrix constructed from points selected so far. This process is continued until a full rank design matrix is obtained—one with as many unique rows (p) as columns and with no linear dependence among any of the columns. This ensures that at least one full rank design matrix is available from the candidate point set, which is necessary to determine the model coefficients by linear regression. Additional points necessary to meet a total point count specified at the start of the process (in order to satisfy precision requirements, for example) are then simply selected at random to provide broad coverage within the inference space.

While design points chosen by this process will generally yield well-formulated regression models, not all of these points necessarily represent physically realizable cavity geometries. This is because entry angles must exceed a certain depth-dependent minimum to ensure realizable cavity floor lengths, and side angles must exceed a certain depth-dependent minimum to ensure realizable cavity floor widths. These constraints form a concave hull within the design space, outside of which all points are infeasible. Therefore, a significant departure from this D-optimal point selection process had to be developed to accommodate these constraints on cavity geometry.

Figure 3 illustrates the cavity geometry leading to the following equations governing these constraints:

$$\xi_L = L - \left[\frac{D}{\tan(\alpha)} + \frac{D}{\tan(\beta)} \right] - L_B \geq 0 \quad (5a)$$

$$\xi_W = W - \left[\frac{2D}{\tan(\phi)} \right] - W_B \geq 0 \quad (5b)$$

where ξ_L and ξ_W are zero when the constraints are just satisfied. If these values are less than zero, the constraints are not met. If they are positive but relatively small, the constraints are satisfied but the point is said to be “near the constraint boundary.” The minimum length and width of the bottom of each cavity, L_B and W_B , were each selected to be 0.5 inch.

Initially, points with negative values of either ξ_L or ξ_W were simply dropped. Unfortunately, this did not completely address the problems with this design approach because there were also too few “near constraint” points. This would mean that smaller cavities were not likely to be modeled as well as larger cavities. It was therefore necessary not only to eliminate “out of constraint” points but also to add “near constraint” points.

To ensure a large number of candidate points near the constraint boundaries as needed for good fits in that part of the inference space, 2000 additional points were selected by a maximal distance algorithm that picks points as far away from each other in the inference space as possible. This has the effect of “filling holes” in the inference space by forcing a certain fraction of the points into the near-constraint region. Points were eliminated for which either $\xi_L < 0$ or $\xi_W < 0$. This added some additional points near the boundary, and also ensured a generally uniform distribution of candidate points throughout the seven-dimensional inference space.

A number of additional points guaranteed to be close to the constraint boundary were then added in the following way: A total of 50,000,000 points randomly distributed within the inference space were generated. All of

these points for which both ξ_L and ξ_w were greater than 0 and less than 0.1 were retained. Only 278 such points were identified. These were added to the candidate point list. These points were selected via an algorithm written by Wayne Adams of Stat-Ease, Inc., to run in the “R” statistical computing language. The Adams algorithm was used in the same way to generate a further 772 points defined as “near” points. Near points feature values of ξ_L and ξ_w greater than 0.1 and less than 0.5. This again assured a good representation within the candidate point list of points that were in the region moderately close to the constraint boundary. The process of adding points within specified ranges of ξ_L and ξ_w while eliminating points with negative values of those variables resulted in a final candidate point list of 2,507 points. These points all satisfied the constraints of Eq. (5), provided good coverage (few “holes”) throughout the seven-dimensional inference space, and concentrated points near the constraint boundary as necessary to achieve as good a result with small cavities as with large cavities. The final set of design points was selected from this list by the D-Optimal selection process we now describe:

B. D-Optimal Point Selection

The 124 points necessary to produce each RSM case matrix for this study were all drawn from the 2,507 candidate points created as described above. (Fifteen additional conformation points were selected at random, subject to the constraints in Eq. (5).) The algorithm for selecting the points to fit the models begins by picking the candidate point with the largest Euclidean norm as the first design point. This is the point that is the farthest from the center of the design space. Figure 2 illustrates why points far from the center are desirable.

Each subsequent design point is selected to minimize the determinant of the covariance matrix in Eq. (3) while increasing the rank of the design matrix. When a full rank design matrix has been constructed, the remaining points in the candidate list are evaluated through a series of exchange steps to see if substituting them for points provisionally in the design will result in an improvement.

The process begins with a one-point exchange. The point in the candidate list that decreases the determinant of the covariance matrix the most is added to the design. The point in this augmented design that increases the determinant of the covariance matrix the most is then dropped from the design and placed back in the candidate list. One-point exchange steps are continued until there is no improvement in the model or until every candidate point has been evaluated. A series of two-point exchanges then begins, in which the *pair* of points drawn from the candidate list that decreases the determinant of the covariance matrix the most is added, and the pair of points from the augmented model that increases the determinant of the covariance matrix the most is deleted. This process is repeated with more and more points in the exchange until typically 5-point to 10-point exchanges have been performed. Each step is taken to decrease the determinant of the covariance matrix the most by adding points, then increase it the least by dropping an equal number of points from the design.

The entire process is repeated a number of times (typically one to 10 times), each with a different randomly selected start point. The design with the smallest covariance matrix determinant is retained as the final design. Each experiment designed for this study involved four such replications.

To see the effect of these changes in the experiment design, consider the “hat matrix,” \mathbf{H} , which derives its name from the fact that a vector of individual response measurements, \mathbf{y} , can be transformed into a vector of corresponding response predictions, “ $\hat{\mathbf{y}}$,” as follows:

$$\hat{\mathbf{y}} = \mathbf{H}\mathbf{y} \quad (6)$$

It can be shown that the variance in predicted responses is likewise related via the hat matrix to the variance in individual response measurements, σ^2 , as follows:

$$\text{Var}(\hat{\mathbf{y}}) = \mathbf{H}\sigma^2 \quad (7)$$

The hat matrix is a function only of the design matrix:

$$\mathbf{H} = \mathbf{X}(\mathbf{X}'\mathbf{X})^{-1}\mathbf{X}' \quad (8)$$

The standard error in response predictions is just the square root of the prediction variance:

$$\sigma_{\hat{\mathbf{y}}} = \sqrt{\mathbf{X}(\mathbf{X}'\mathbf{X})^{-1}\mathbf{X}'\sigma} \quad (9)$$

Equation (9) tells us that the uncertainty in predictions made at the points used to fit the model only depends on two things, the intrinsic variability in the response estimates and the design of the experiment (i.e., the model being fit and the selection of independent variable levels in the test matrix). Since the researcher has complete control over \mathbf{X} , there is considerable latitude for improving the quality of response predictions by optimizing the design for this purpose, and this is a major mechanism by which MDOE experiment designs improve the quality of research results.

Equation (9) is especially useful because, for a unit standard deviation in response measurements, it allows us to know the distribution of unit standard prediction errors even before the experiment is executed. By setting $\sigma = 1$, we can use graphical representations of Eq. (9) to compare candidate experiment designs as in Fig. 4, where the distribution of unit prediction standard errors is displayed for two variations on the experiment design discussed above. The distributions on the left and on the right in Fig. 4 correspond to designs with different numbers of parameters. The reduction in average prediction variance reflects Eq. (2). These distributions of unit standard prediction errors allow rapid and convenient comparisons among candidate MDOE experiment designs. They can be displayed as a function of any two of the independent variables at a time, with the remaining variables held at specifiable levels. Changing the variables that are plotted and the levels of the other variables provides insight into the robustness of the experiment design. We seek designs for which the distribution of standard prediction errors is as low and as uniform as possible.

C. Variable Selection and Model Building

After identifying over 2500 D-optimal candidate points that each satisfied all factor constraints and were all distributed optimally throughout the inference space, with points near the constraint boundaries and other points distributed to minimize large holes in the inference space, the D-Optimal selection process identified a 124-point subset of these candidate points that most minimized the determinant of the covariance matrix for a third-order polynomial. This set of points is the smallest number that will fit the polynomial and still minimize uncertainty in the regression coefficients, while allowing up to 20 additional lack of fit degrees of freedom to fit unanticipated higher-order complexities in the data.

For each of the 33 body point zones, the USA/Boeing code was executed for each of 124 unique combinations of variables from Table 1. Various thermal and structural reentry load responses were recorded for each such case. Stepwise procedures were used to develop third-order polynomial models for each response variable as a function of all seven independent variables. These procedures are intended to identify a subset of the 104-terms in a full third-order polynomial in five numerical variables and two 2-level categorical variables, by discarded statistically insignificant terms from the full model. These are terms that are not large enough to be distinguished from zero with high confidence.

While the number of terms eliminated from the regression models in this way varied from response variable to response variable, and to a lesser degree from body point zone to body point zone, typically only around a quarter of the terms were retained. These were mostly first-order and second-order terms, with a few mixed cubic terms and an occasional full cubic term retained. This suggests that the responses being modeled were predominantly second order, with some slight additional complexity that could be addressed with a few higher-order terms. The elimination of so many terms enhanced the quality of predictions by Eq. (2), and also brought a certain level of clarity to the response models by reducing the clutter associated with superfluous terms in the response model. The truly important terms provided insight into which factors most significantly influenced each thermal and structural reentry load response.

Three stepwise procedures employed in various combinations in the model-building process will now be described briefly. These are known as (1) “Forward Selection,” (2) “Backwards Elimination,” and (3) “Stepwise Regression,” which is actually a combination of the first two methods. Standard textbooks on regression analysis^{6-8,10} can be consulted for a more detailed description of these methods, which are outlined here.

Forward Selection begins by assuming that only the intercept term is in the model. It then provisionally adds the one term that has the highest correlation with the response. An analysis of variance (ANOVA) is conducted for this two-term model. If the explained variance is increased by a specified threshold amount (indicating low probability that the added term is insignificant), the term is retained and the forward selection process continues. The next term to be provisionally entered is the one that has the highest correlation with the response after correcting for first term entered. This is called a “partial correlation,” and an ANOVA on this model indicates the probability that this term makes a significant change to the explained variance of the model. If so, this term is retained. The process continues until either all terms from the initial model have been included, or until the most significant remaining term does not cause enough of a change in the explained variance to satisfy the entry criterion.

Backward Elimination is similar to forward selection except that the process attempts to identify terms for the final model by working in the opposite direction. The backward elimination method begins with all terms from the

initial model included. The term having the weakest correlation with the response is provisionally rejected, and the impact on the explained variance of the model is assessed by ANOVA. If the rejection of this term produces a significant reduction in the variance explained by the model, it is retained and the process stops. Otherwise, the process continues until no terms in the model can be rejected without causing a significant reduction in the variance explained by the model, or until the only remaining term is the intercept.

Stepwise Regression is a combination of forward selection and backward elimination. It begins as a forward selection process and continues until the model contains the intercept and two regressors. Now backward elimination is applied to the three-term model, eliminating each regressor in turn to assess the reduction in the model's explained variance via ANOVA. When the backward elimination step is finished, the forward selection process is resumed with whichever remaining term causes the most significant increase in the explained variance. If such a term increases the explained variance by a user-specified threshold amount, it is retained and the backward elimination is initiated on the new model. This is continued until there are no candidate regressors significant enough to enter the model and none in the model that are so weak that they can be eliminated with no significant effect.

Proponents of stepwise regression note that the backward elimination step protects against multicollinearity, a condition that exists when two or more regressors are highly correlated and therefore do not each contribute independently to the model. If two regressors are highly correlated, adding one of them to the model may render the first one superfluous. In such a case it is better to eliminate that term.

It is common to apply more than one stepwise procedure to the same data set. For example, in this study backward elimination was applied first, to give all model terms a chance to be included. Stepwise regression was then applied to the surviving model terms, with one more applications of backward elimination as a finishing step, to reject the occasional high-order term that survives the first two applications without contributing significantly to the model.

To summarize, we have applied stepwise regression procedures to develop response surface models from a volume of data ample for this purpose but no larger, drawn from a large list of candidate data points that all satisfied complex constraints among the independent variables. The large candidate point list was designed to include a substantial fraction of candidates in the region of constraint boundaries in the seven-dimensional inference space, while providing uniform coverage throughout the rest of the inference space. This ensured a good representation for small cavities as well as large cavities. Points selected from the candidate list consisted of a relatively small subset with good D-optimal properties; that is, with a distribution within the inference space that minimized the uncertainty in regression coefficients by minimizing the determinant of the covariance matrix associated with the proposed third-order polynomial model. In short, many steps were taken through the MDOE process to ensure that the polynomial response models would faithfully reproduce the discrete values generated by the underlying debris impact assessment codes.

Figure 5 is a schematic representation of the relationship between a set of such discrete code outputs and the polynomial response model chosen to represent them. The difference between each output of the underlying debris impact assessment code and the corresponding polynomial response model prediction is a residual which gives some measure of how well the polynomial model represents the underlying code. Figure 6 displays the residuals for a representative example—structure temperature for body point zone 1602. The standard deviation of these residuals is 1.2° F. There are a total of 139 residuals in Fig. 6, of which the 124 blue points are model residuals and the 15 red points are confirmation point residuals. The model residuals might be expected to be relatively small because the least-squares fitting algorithm distributes the regression coefficients to ensure that this is so. The fact that the confirmation points residuals are of the same order as the model residuals suggests that the model results are transferable to points other than those for which the model was developed, implying a certain degree of robustness in the polynomial response model.

Models developed by these stepwise procedures were subjected to a wide array of model adequacy tests. The residuals were examined graphically as a function of predicted value, run number, and independent variable level to look for tell-tale patterns that would reveal a poor fit. Numerous model adequacy metrics were computed, including R-squared, adjusted R-squared, and Predicted R-squared statistics. The leverage of each point was computed as was its Cook's distance, two statistics indicating the susceptibility of the model to outliers. The magnitude of the residuals was examined to determine if the prediction accuracy was acceptable. The distributional properties of the residuals were also examined, to look for evidence of systematic lack of fit. Only after passing this battery of quality assurance tests was a polynomial model declared to adequately represent the underlying USA/Boeing code. It was then deemed suitable to use for propagating errors in the specification of cavity geometry into uncertainty in the thermal and structural reentry load estimates.

VI. Error Propagation

The purpose of approximating the underlying USA/Boeing debris impact assessment code with polynomial response functions was to facilitate error propagation, as will be discussed in this section. Geometry estimates describing the size and shape of debris impact cavities necessarily suffer from some degree of uncertainty. In addition to the usual experimental error in estimates of any experimental variable, there is also a significant component of uncertainty attributable to the approximations that are made in representing the cavities in the underlying damage assessment codes. Figure 7 provides an example of such an approximation.

Because of the inevitable error associated with these cavity size and shape approximations, it is important to understand how this error translates into uncertainty in the response estimates. That is, we wish to know how much uncertainty is introduced into the estimates of such system responses as structure temperature and structural margin of safety by the fact that the cavity size and shape estimates are imperfect. This is the central question for which the polynomial response model approximations were developed, as will be described in more detail in this section.

Consider a simple function of two random variables, x_1 and x_2 . Such a function can be represented as a Taylor series expansion about the means of those variables, which to first order is as follows:

$$y(x_1, x_2) = y(\bar{x}_1, \bar{x}_2) + \frac{\partial y}{\partial x_1}(x_1 - \bar{x}_1) + \frac{\partial y}{\partial x_2}(x_2 - \bar{x}_2) \quad (10)$$

or

$$y - \bar{y} = \frac{\partial y}{\partial x_1}(x_1 - \bar{x}_1) + \frac{\partial y}{\partial x_2}(x_2 - \bar{x}_2) \quad (11)$$

Following Coleman and Steele,¹¹ imagine now that there are n measurements of x_1 and x_2 , from which n estimates of $y(x_1, x_2)$ have been computed. We know the uncertainties in x_1 and x_2 , and we wish to propagate these uncertainties into an estimate of the corresponding uncertainty in y .

For all n sets of data, square the terms on the left and right side of Eq. (11), and average the squared values.

$$\begin{aligned} \frac{1}{n} \left[\sum_{i=1}^n (y_i - \bar{y})^2 \right] &= \left(\frac{\partial y}{\partial x_1} \right)^2 \times \frac{1}{n} \left[\sum_{i=1}^n (x_{1i} - \bar{x}_1)^2 \right] + \left(\frac{\partial y}{\partial x_2} \right)^2 \times \frac{1}{n} \left[\sum_{i=1}^n (x_{2i} - \bar{x}_2)^2 \right] \\ &+ 2 \left(\frac{\partial y}{\partial x_1} \right) \left(\frac{\partial y}{\partial x_2} \right) \times \frac{1}{n} \left\{ \sum_{i=1}^n [(x_{1i} - \bar{x}_1)(x_{2i} - \bar{x}_2)] \right\} \end{aligned} \quad (12)$$

By the definition of population variance, and with an obvious notational extension for the cross term on the right, in the large n limit this becomes:

$$\sigma_y^2 = \left(\frac{\partial y}{\partial x_1} \right)^2 \sigma_{x_1}^2 + \left(\frac{\partial y}{\partial x_2} \right)^2 \sigma_{x_2}^2 + 2 \left(\frac{\partial y}{\partial x_1} \right) \left(\frac{\partial y}{\partial x_2} \right) \sigma_{x_1 x_2} \quad (13)$$

The correlation coefficient for x_1 and x_2 is defined as

$$\rho_{x_1 x_2} = \frac{\sigma_{x_1 x_2}}{\sigma_{x_1} \sigma_{x_2}} \quad (14)$$

so that Eq. (13) becomes

$$\sigma_y^2 = \left(\frac{\partial y}{\partial x_1} \right)^2 \sigma_{x_1}^2 + \left(\frac{\partial y}{\partial x_2} \right)^2 \sigma_{x_2}^2 + 2 \left(\frac{\partial y}{\partial x_1} \right) \left(\frac{\partial y}{\partial x_2} \right) \rho_{x_1 x_2} \sigma_{x_1} \sigma_{x_2} \quad (15)$$

For independent x_1 and x_2 the correlation coefficient is zero, and for k independent variables Eq. (15) generalizes to

$$\sigma_y^2 = \left(\frac{\partial y}{\partial x_1}\right)^2 \sigma_{x_1}^2 + \left(\frac{\partial y}{\partial x_2}\right)^2 \sigma_{x_2}^2 + \dots + \left(\frac{\partial y}{\partial x_k}\right)^2 \sigma_{x_k}^2 \quad (16)$$

It is clear from Eq. (16) why it is so convenient to represent a complex computer code as a simple low-order polynomial function of the independent variables. The derivatives in Eq. (16) become trivial to evaluate in that case, and since the standard deviations for all x_i is assumed to be known, the corresponding response uncertainty in y is easily computed from Eq. (16). For example, if y represents reentry structure temperature as a function of cavity geometry variables that all have associated with them some uncertainty, then Eq. (16) can easily quantify the uncertainty in structure temperature attributable to the cavity geometry uncertainties. Table 2 lists the uncertainty specifications for debris impact cavity geometry variables assumed in this study.

Table 2. Uncertainty in cavity geometry variables.

Variable	Specified Uncertainty	Standard Error (“one σ ”)
Length	$3\sigma = 0.125$ in.	0.042 in.
Width	$3\sigma = 0.125$ in.	0.042 in.
Depth	$3\sigma = 0.125$ in.	0.042 in.
Entry angle	$2\sigma = 5^\circ$	2.5°
Sidewall angle	$2\sigma = 5^\circ$	2.5°

Figure 8 shows a typical polynomial response surface model for structure temperature as a function of cavity length and depth. The length and depth are displayed in coded units, where -1 and +1 correspond to the smallest and largest values of the indicated variables for which the regression model is valid. The corresponding uncertainty in structure temperature is displayed for the propagation of a two-sigma uncertainty in cavity geometry variables. So, by Fig. 8, the predicted temperature for a relatively long, deep cavity would have a two-sigma uncertainty of about $\pm 2.5^\circ$ F due to the errors in cavity geometry that are presented in Table 2. It is easy to examine the impact of errors of a different magnitude once the polynomial response and error propagation models have been created, and the sensitivity of the response uncertainty to errors in specific geometry variables can also be easily examined. Obviously, other insights can be obtained from response surface models of reentry loads and their corresponding uncertainties, as will be discussed in the next section.

VII. Results and Discussion

This paper is focused on the MDOE experiment design and analysis process used in this study to quantify the impact of cavity geometry error on estimates of thermal and structural reentry loads. Many other insights can be, and were, achieved from the polynomial response models developed in this study. In addition to revealing interesting patterns in the data that illuminated the underlying physical processes in play, the modeling process also revealed certain apparent anomalies in the parametric codes upon which they were based. These patterns and apparent anomalies are documented in detail in a comprehensive report¹² prepared by the Jacobs Technology Engineering and Science Contract Group in Houston. While a detailed exposition in this paper of all of these findings would be redundant, we site a few representative examples that bear on the response surface modeling activity and that therefore have some impact on the cavity geometry error propagation problem.

The coefficient of variation (CV) is a metric that describes how well a response surface function represents the underlying process it models. It is computed by dividing the standard deviation of residuals by the mean response over the points for which the residuals were computed. Figure 9 presents the coefficients of variation for the response variables modeled in this study. Note the pronounced difference between coefficients of variation for the thermal responses (blue) and the structural responses (red). The CV values for the thermal responses are less than 2%, while none of the structural response CV values are less than 4%—twice the maximum thermal CV. This illustrates that it was much more difficult to model the structural responses than the thermal responses.

There are a number of plausible explanations for this. Some of the structural responses exhibited a certain degree of discontinuous behavior that is difficult to model well with a low-order polynomial, for example. The tiles are believed to remain in plane until stressed to a certain threshold level, beyond which they “pop” into an out-of-plane

configuration. In mathematical terms, the out-of-plane-deflection (OOPD) functions are not continuous with continuous derivatives, conditions that must be met to represent them well with a Taylor series corresponding to the polynomial models we have tried to fit. This may explain why the two OOPD responses in Fig. 9 have the greatest coefficients of variation—about 14% for the Mach 25 to Mach 4 regime of the reentry trajectory and about 8% for the terminal area energy management regime. These are four to seven times greater than the largest thermal CV value.

The CV for tile factor of safety (FS) was twice as large as the largest thermal CV, a fact that may be attributable to a difference in behavior as the cavity exceeds the threshold level beyond which the RTV bonding fails and the tile is lost. The true tile FS is believed to degrade as debris impact loads increase, until this threshold is reached. Beyond that point, the tile FS is essentially undefined—the concept of factor of safety for a tile that is already lost is meaningless. Unfortunately, the underlying code reports a numerical value of zero for FS in such circumstances, rather than “Not Applicable” or some other such non-numeric descriptor. The result of this is represented schematically in Fig. 10.

The red curve on the left of Fig. 10 represents the best third-order polynomial approximation to the data points that track the true behavior, represented by the black line. The discontinuity results in a very poor fit over the full range of cavity sizes. For example, in an effort to model both the zero and non-zero FS regions, the response model underestimates tile FS where it has meaningful values. This is followed by a region in which the tile has already failed, but the model is reporting non-zero FS values. Beyond that, the modeled FS values go negative, which is physically non-realizable. Finally, a cavity size is reached beyond which the model predicts that as the cavity size increases, the tile FS increases. Obviously these are all nonsensical results, caused by an attempt to use a single, simple polynomial model to span two regions in which radically different physical processes are in play.

The blue curve on the right of Fig. 10 represents how different response models would be fitted to the two regions if the boundary between them were known. For the FS = 0 region, the polynomial response function is a trivial 0th-order function (intercept only); viz., $y = b_0$, where $b_0 = 0$. Where FS is not zero, a conventional polynomial model would be expected to fit the data adequately.

The problem of finding the boundary between meaningful and meaningless FS values is not straightforward. One can simply fit the FS values over cavity dimensions known to be small enough for the tile not to fail, but there is always a danger of being too conservative in this approach. Generally, the behavior right up to the boundary is of interest. An alternative approach, proposed as a result of this study, is to design a *series* of MDOE experiments. In the first, distribute points throughout the entire inference space without regard to whether FS response values are meaningful or not. Analyze the data, noting where the response surface goes below some threshold value. Figure 11 is an example of such a result, where the red and blue regions denote the boundary where FS = 0. (Note that other boundaries might be superior for this type of analysis, such as the boundary beyond which the FS values decline below 1, the lowest level attributable to an undamaged tile, or below 1.4, the acceptable flight threshold.) Design a second experiment in which the data points are distributed only in the region where preliminary results indicate FS values of actual interest. Fit the data in this region. It may be necessary to iterate more than once, but with experience it should become possible to satisfactorily identify the boundary.

In the present study, tile FS models were simply not attempted in body point zones with an excessive number of zero-value FS points (where “excessive” was defined as “five or more”). However, it is still possible that the FS responses were sufficiently discontinuous for the cases in which zero-FS values were modeled, that the low-order polynomial models were still inadequate to properly capture all of the variation in FS over these regions. We conclude that when there are two regions in which the underlying physics is radically different, it is necessary to first identify the boundary between those regions, and then fit separate models in each of them.

Certain anomalies were discovered in the course of building the structural response polynomial models that have been attributed at least in part to the behavior of the Stress Assessor Tool. The Stress Assessor Tool failed to produce estimates of structure margin of safety and tile out of plane displacement (OOPD) for 12 of the 33 body point zones. Adequate third order polynomial fits could not be achieved for structure margin of safety in four body point zones despite data having been produced by the Stress Assessor Tool, and likewise adequate Mach 25 to Mach 4 OOPD models could not be generated in four other zones. TAEM OOPD models could not be generated in five additional body point zones.

Further investigation revealed certain anomalies within the Stress Assessor Tool. For example, in some body point zones there were no changes in structural responses despite large changes in input variables. This caused some of the structure response polynomial models to have only one or two significant terms, as there was not enough change attributable to other factors for the regression coefficients for those terms to be resolvable from zero. In addition, in certain other zones structure margins of safety were sometimes larger for damaged tile than undamaged tile. This is clearly a suspect result. There have been a number of improvements in the Stress Assessor Tool and

other elements of the debris impact assessment software since these observations were originally made; however, these Stress Assessor Tool anomalies may be responsible at least in part for the fact that the structural response coefficients of variation reported in Fig. 9 were so much greater than the thermal coefficients of variation.

The response surface models were fitted using coded variables that take on values of ± 1 at the extremes of the independent variable ranges. The coefficients of the first-order terms in these models therefore represent the change in response over half the independent variable range, and are independent of any physical units associated with the variables. That is, the coefficient of the first-order length term is the same in coded units whether the length is measured in physical units of inches or meters, say. This facilitates a convenient comparison of the relative effectiveness of each factor over the ranges each factor was modeled.

Figure 12 displays the first order coefficients for maximum RTV bondline temperature in seven body point zones. Note that in all seven zones, cavity depth and boundary layer transition time are the dominant factors. The positive depth coefficients simply mean that RTV bondline temperature increases with increasing cavity depth, and the negative coefficients for boundary layer transition time means that lower bondline temperatures are associated with longer transition times, suggesting that the longer the flow remains laminar, the lower the RTV bondline temperature will be.

Coefficient comparisons can be equally revealing for the information they provide on which factors do *not* dominate. For example, it is clear from Fig. 12 that, at least over the range tested, cavity width has a relatively small influence on RTV bondline temperature. This may suggest that the width should be examined over greater ranges the next time. Note, however, that comparisons of first-order coefficients, while generally useful and revealing, do not account for interaction terms or other higher order terms that might be significant in a given model. Such terms need to be examined before final conclusions are reached about the relative importance of various independent variables.

There were six responses modeled in each of 33 body point zones in this study, for a total of 198 polynomial response models. Each of them was subjected to at test to determine if a more satisfactory fit could be achieved by transforming the response model. Commercial software packages of the kind used in this study⁹ provide a direct test on the data for whether such a transformation would be helpful. This test is known as the Box-Cox transformation test,¹³ which examines scaled transformations involving a power transformation, y^λ , of the response variables, y .

A model is fitted for a range of candidate exponents, typically from -3 to +3. The residuals (or the logs of the residuals) are plotted against λ and the value of λ for which this function is a minimum represents a maximum-likelihood estimator of the exponent that produces the best fit in a least-squares sense. Figure 13 is a Box-Cox transformation plot for a structure temperature response model. Note that it has a minimum near $\lambda = 1/2$, corresponding to a suggested square root transformation. This implies that a better model can be achieved by fitting the square root of structure temperature than by fitting untransformed structure temperature.

The vertical blue line in Fig. 13 marks the current exponent ("1," so no transformation). The light red lines in this figure represent lower and upper limits on the 95% confidence interval for the optimum λ estimate. The transformation is generally applied with a value of λ within the interval that may have some physical significance, rather than the exponent for which the Box-Cox transformation plot achieves an absolute minimum.

Figure 14 shows an interesting pattern associated with these power transformations. In this figure, maximum RTV bondline temperature is plotted for each body point zone in increasing order of RTV temperature. The zones are color coded according to the power transformation that produced the best fit. Note that fits in the seven hottest zones were improved with a square root transformation ($\lambda = 1/2$) and fits in the five coolest zones were improved with an inverse square root transformation ($\lambda = -1/2$). For the mid-temperature zones, the fits were improved with a logarithmic transformation ($\lambda = 0$). It is not immediately obvious that a power transformation in which the exponent is zero corresponds to a logarithmic transformation, but in the limit as lambda approaches zero, y^λ approaches $\log(y)$. That is, for small lambda, y^λ plots against $\log(y)$ very nearly as a straight line.¹³

The meaning of the pattern in Fig. 14 is unknown, but the positive correlation between maximum RTV bondline temperature and the exponent of a power transformation that improves the fit seems unlikely to be coincidental with so many independently modeled zones involved. This pattern may lead to future insights into the underlying physical laws governing the bondline temperatures.

VIII. Summary of Results

The Modern Design of Experiments (MDOE) is a formal experiment design, execution, and analysis process originally introduced at Langley Research Center in 1997 to produce aerospace research results at the lowest cost consistent with specified maximum acceptable levels of uncertainty. It has recently been applied in a series of computational experiments in support of the Space Shuttle program. Complex computer codes developed at Johnson

Space Center by NASA and its contractor team are used to estimate thermal and structural reentry loads on the Shuttle Orbiter induced by ice and foam debris impact during ascent. Such debris can create cavities in the Shuttle Thermal Protection System. The computer codes are designed to forecast reentry temperatures and various structural safety factors associated with these cavities, which for a given location are a function of their size and shape, among other things.

The sizes and shapes are approximated to accommodate a code limitation that requires simple “shoebox” geometries to describe the cavities—rectangular areas and planar walls that are at constant angles with respect to vertical. These approximations induce uncertainty in the code results. MDOE methods were used to develop a series of 33 resource-minimal computational experiments—one for each body point zone on the Orbiter’s reentry windward side—designed to generate low-order polynomial graduating functions to approximate the more complex underlying codes. These polynomial functions were then used to propagate cavity geometry errors to estimate the uncertainty they induce in the reentry load calculations performed by the underlying code.

The effort reported here was a methodological study focused on evaluating the application of MDOE to this class of problem. The actual code that was executed during these computational experiments was an earlier version of the current operational Shuttle codes, which have since been improved and upgraded. Specific uncertainty results were therefore of secondary importance to the development of an MDOE-based methodology that could be applied to future operational codes in a rapid and low-cost way to assess the effects of cavity geometry uncertainty. Nonetheless, general trends in the relationships between responses and independent variables reported here are believed to apply to the current code as well as the earlier version used for this analysis.

The emphasis in this paper has been on the experiment design details, including computational resource savings and prediction quality enhancements achieved by reducing the number of variables as a result of preliminary MDOE screening experiments. Special experiment design considerations associated with complex, non-linear independent variable constraints have been described, as have the specific methods for developing MDOE test matrices and for generating response surface models with minimal uncertainty in the regression coefficients. Details for applying these methods to propagate uncertainty in cavity geometry specifications have been provided.

Key findings of this study are summarized as follows:

- 1) Preliminary MDOE screening experiments reduced the number of candidate factors from 10 to 7. This resulted in a reduction of 5,445 computer runs with an estimated savings of over 500 hours of CPU time.
- 2) Reentry loads can be adequately modeled as a second-order function of the cavity geometry and Shuttle state variables examined, augmented by a few third-order terms involving mostly wall angles and their interactions with other independent variables.
- 3) Thermal and structural reentry loads are influenced by numerous interactions among the cavity geometry and Shuttle state variables examined in this study.
- 4) MDOE models predicted structure temperature and RTV bondline temperatures within 2% of the underlying numerical codes. There was better agreement with the underlying code for thermal responses than for structural responses, with the difference attributed to a combination of discontinuous behavior in some structural variables and some anomalies in the underlying Structure Assessment Tool that have since been addressed.
- 5) While the approach used in this paper accommodates response functions that can be highly nonlinear in the independent variables, it is not suitable for functions that are discontinuous. One example may be tile out of plane deflection, in which relatively little deflection may occur until applied loads are of sufficient magnitude to cause a sudden dislocation of the tile. If this is the actual mechanism by which tiles achieve an out of plane deflection, then this process cannot be adequately modeled with simple low-order polynomials.
- 6) When there are two regions in which the underlying physics is radically different, it is necessary to first identify the boundary between those regions and then fit separate models in each of them. This is the case, for example, when cavity dimensions exceed the threshold beyond which the tile’s RTV bond fails and the tile is lost. Unacceptable results are likely to be achieved if an attempt is made to fit the same model across regions where there is a physically meaningful tile factor of safety and regions beyond which the tile has been lost. Tile Factor of Safety is undefined for cavities large enough to produce sufficiently high RTV bondline temperatures to lose the tile. Such circumstances require special procedures for estimating Tile FS uncertainty where this metric has physical meaning.
- 7) Larger cavity dimensions (length, width, and depth) increased reentry loads as expected. Reentry loads were also increased for steeper wall angles, a result attributed to the loss of less reflected thermal energy. Higher baseline heating and shorter boundary layer transition times also resulted in greater reentry thermal and structural loads.

- 8) Thermal loads were driven more by cavity depth and length, and Shuttle state variables (baseline heating and boundary layer transition time, a surrogate for surface roughness) than by cavity width or shape (entry angle and sidewall angle).
- 9) The thermal and structural response variables were relatively weak functions of cavity shape (entry angle, exit angle, and sidewall angle). Over the range of angles typically observed in practice, response variables were determined to be virtually independent of exit angle.
- 10) Independent variables that were the most sensitive factors in determining thermal loads were not always the most sensitive factors for determining structural loads.
- 11) The lowest structural margins of safety did not always correlated with the highest structure temperatures and the highest structural margins of safety did not always correlated with the lowest structure temperatures. This may reflect differences in structural geometry and materials properties as well as different methodologies and computer codes used to perform stress analyses for such diverse structures as the Forward Fuselage/Chine Assembly, the Mid-Fuselage Bottom Panel, the Mid-Fuselage/Wing Glove Sheet Stringer, the Elevon/Wing Honeycomb, the Mid-Fuselage/Wing Carry-Through Waffle, and the Aft-Fuselage Waffle Panel. Each structure has different geometries and different material properties.
- 12) Larger out of plane deflections correlated with smaller structure margins of safety, and smaller out of plane deflections correlated with larger structural margins of safety, for both the Mach 25 to Mach 4 portion of the reentry trajectory and the terminal area energy management portion.
- 13) Cavity surface area (length and width) had more influence on structural effects (margin of safety and out of plane deflection) than it had on thermal effects (maximum structure and RTV bondline temperatures).
- 14) Errors in the specifications of cavity geometry have a greater effect on the uncertainty in estimates of larger thermal and structural loads than on smaller loads.
- 15) The MDOE response models revealed certain patterns in the functional form of the response models that may provide further insights into the underlying physics.

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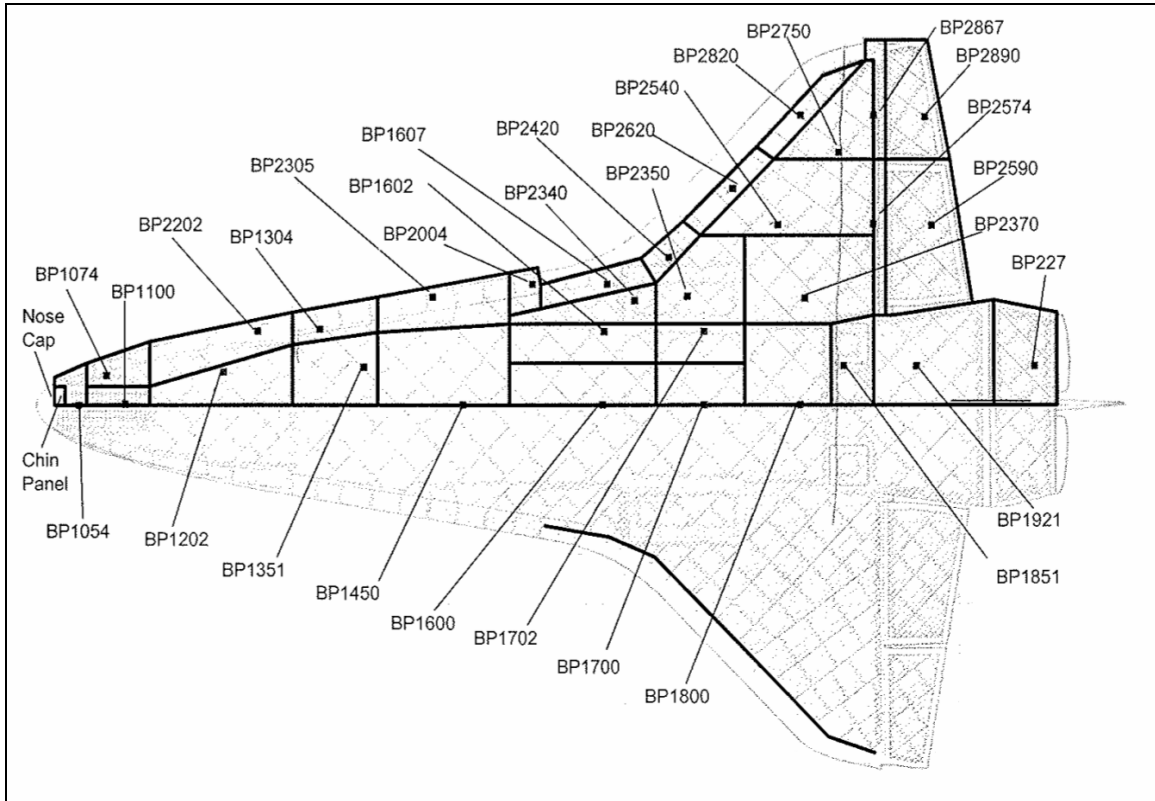


Figure 1. Body point zones on Shuttle Orbiter.

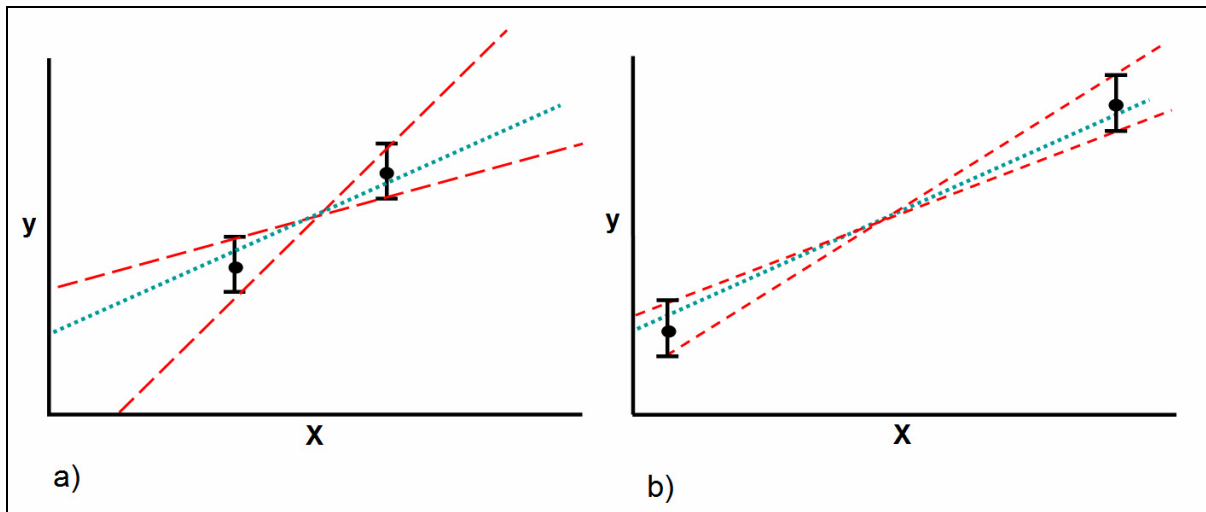


Figure 2. Uncertainty in slope and intercept estimates depends on fitted data site selection. a) Greater uncertainty. b) Less uncertainty.

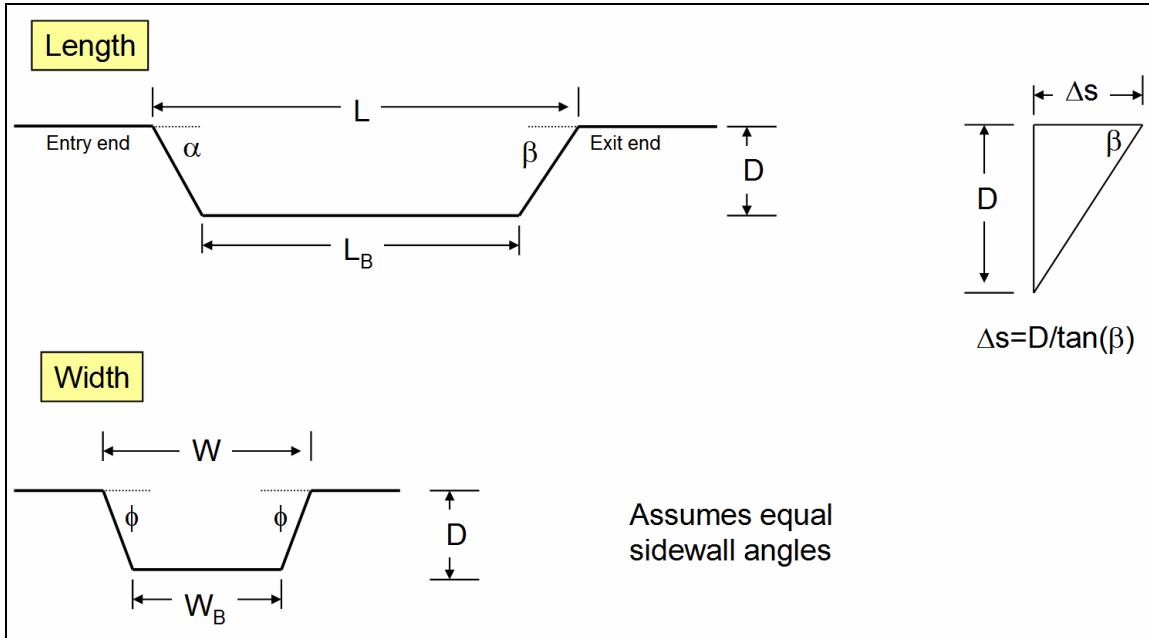


Figure 3. Cavity constraint geometry.

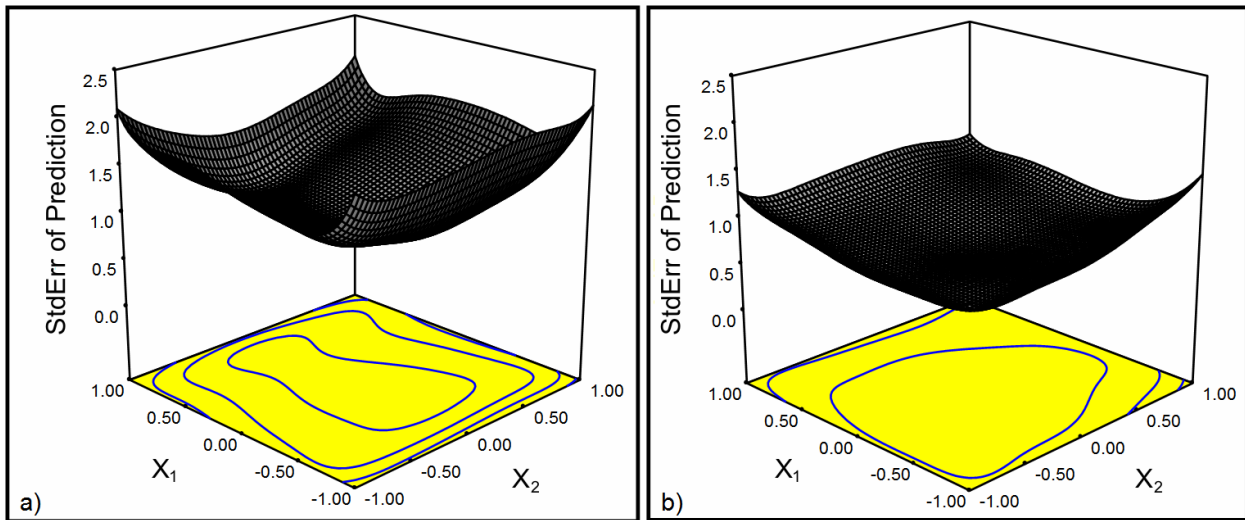


Figure 4. Distribution of unit standard prediction errors. a) D-optimal design with 10 independent variables, b) D-optimal design with 7 independent variables.

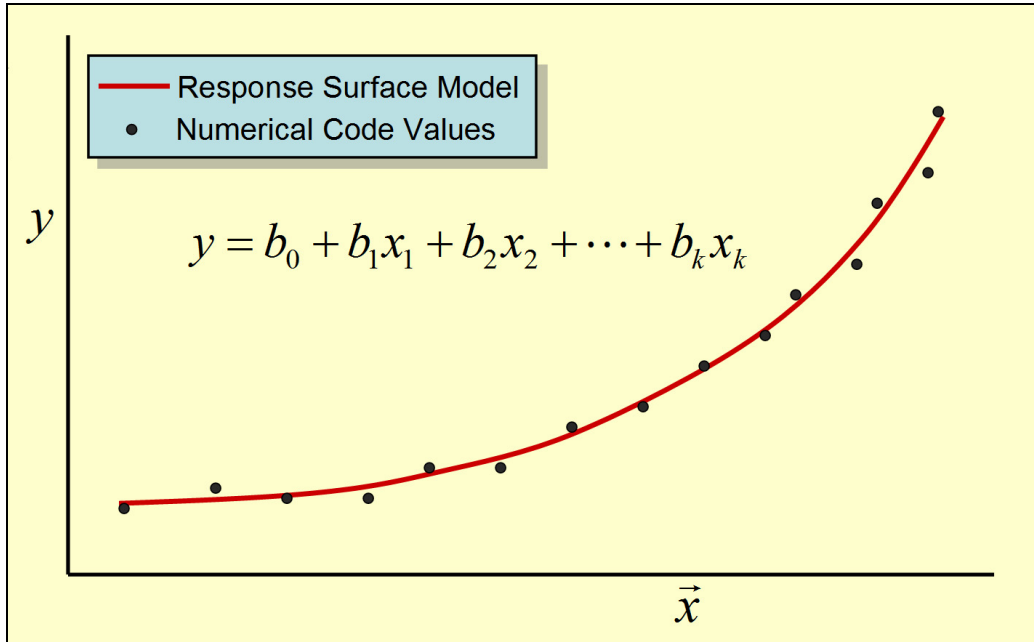


Figure 5. Polynomial response surface modeling approximation to numerical code.

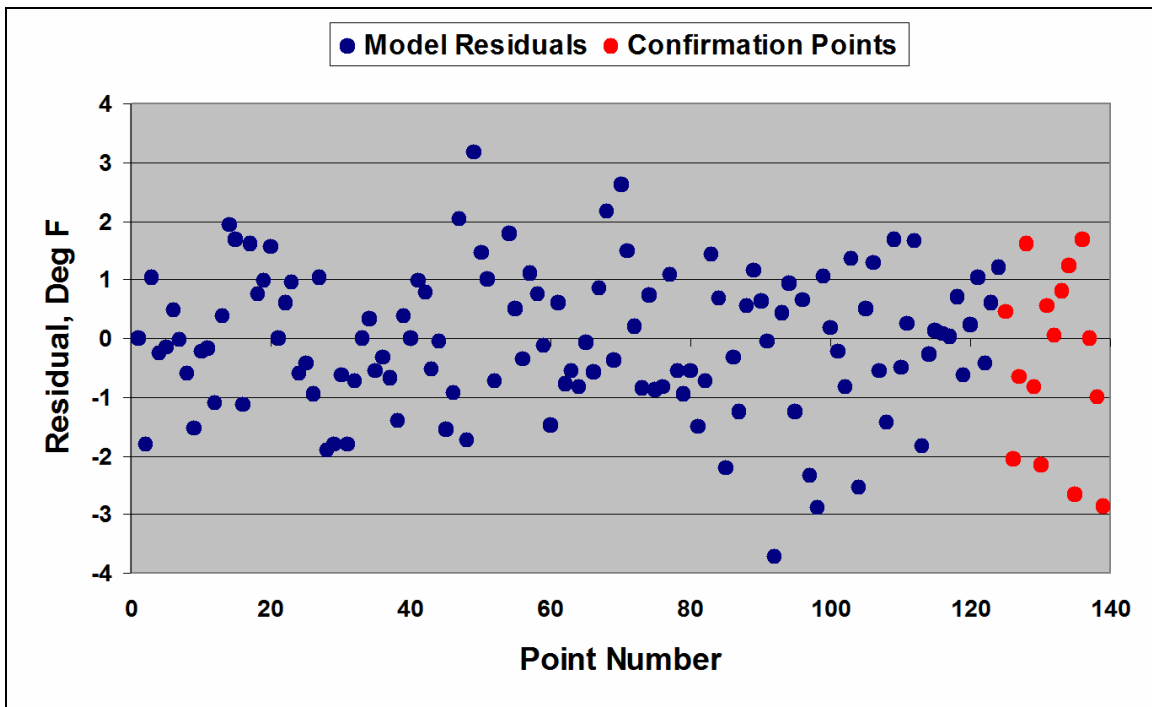


Figure 6. Example of residuals: Structure temperature for BPZ 1602. Standard deviation: 1.2° F. Confirmation point residuals comparable to design point residuals.

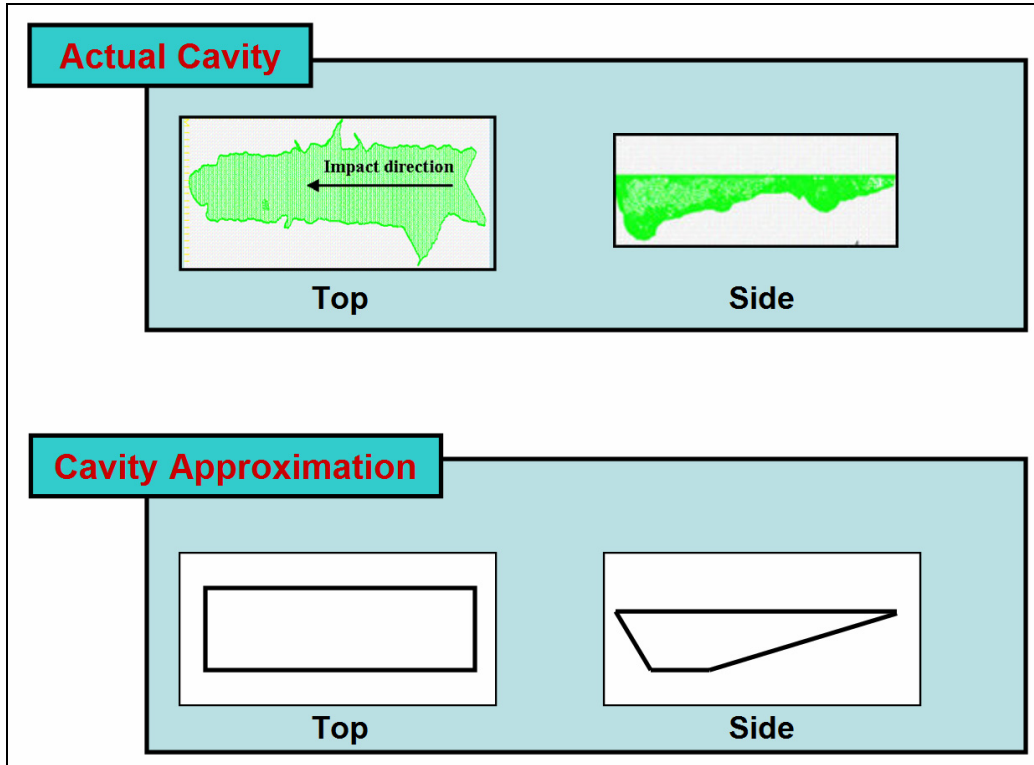


Figure 7. Simplified geometry introduces uncertainty.

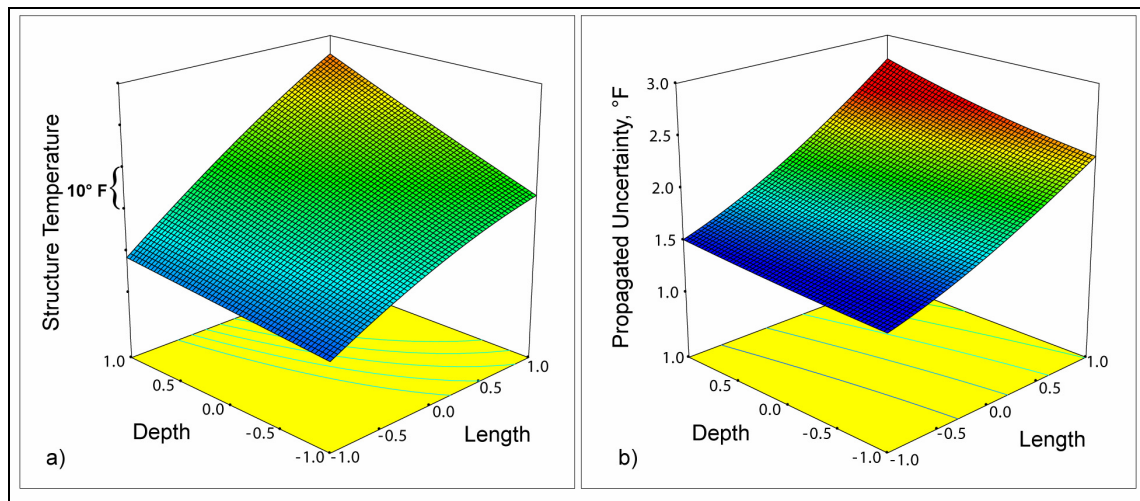


Figure 8. Structure temperature for representative cavity in BPZ 2540. a) Response surface for temperature. b) Response surface for uncertainty in temperature due to 2σ uncertainty in cavity geometry variables.

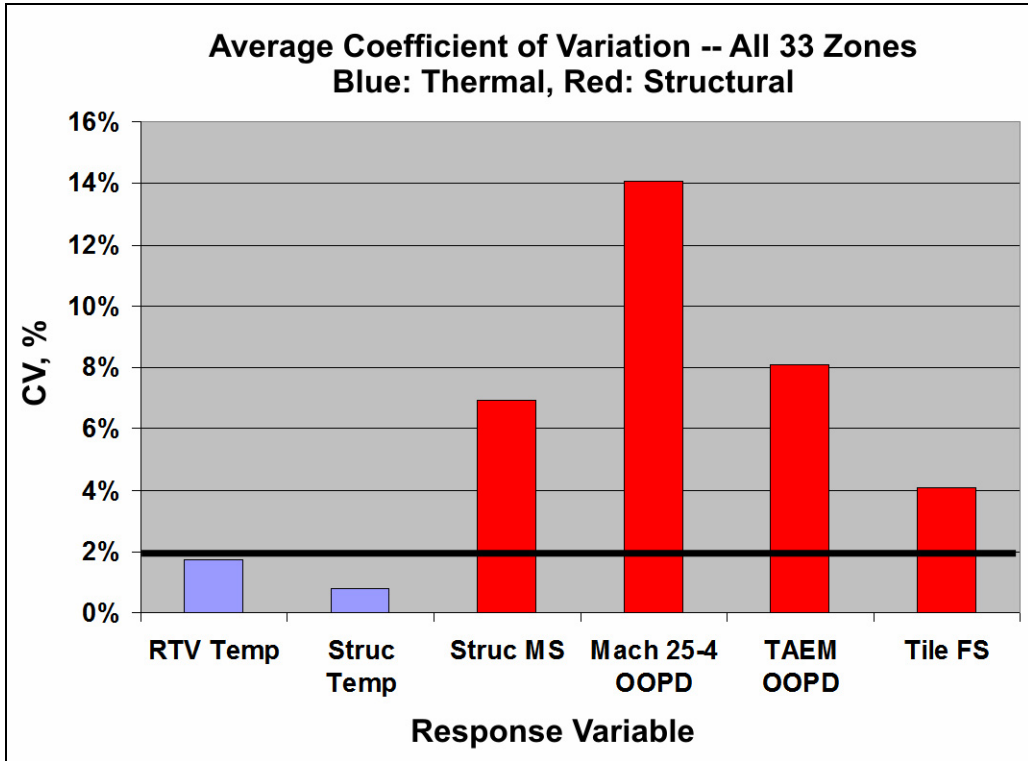


Figure 9. Comparison of average coefficients of variation for thermal and structural response variables across 33 body point zones. More variability in RSM predictions for structural responses than for thermal responses.

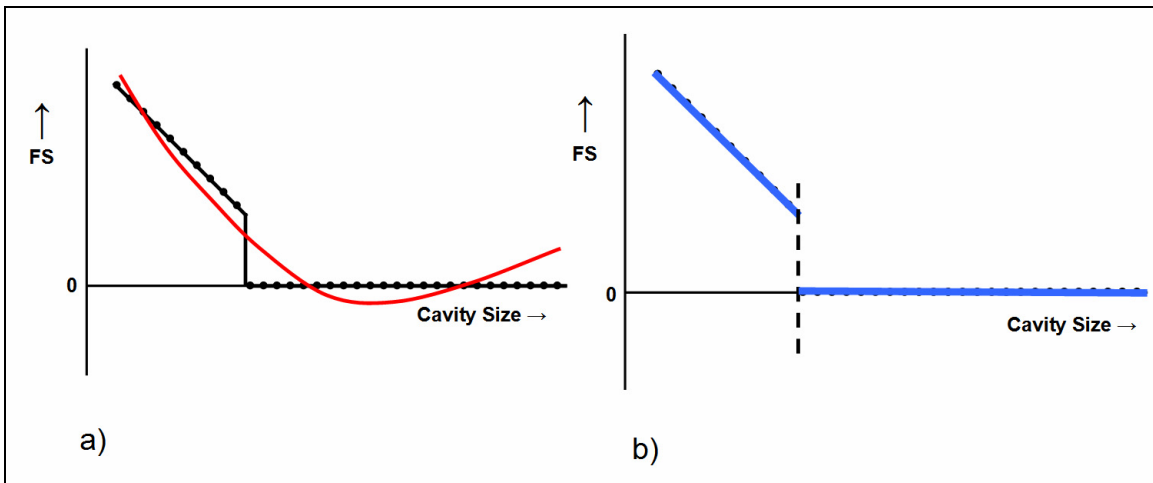


Figure 10. Fitting tile factor of safety with many zeros. a) One function spanning entire inference space. b) Separate functions for $FS = 0$ and $FS \neq 0$.

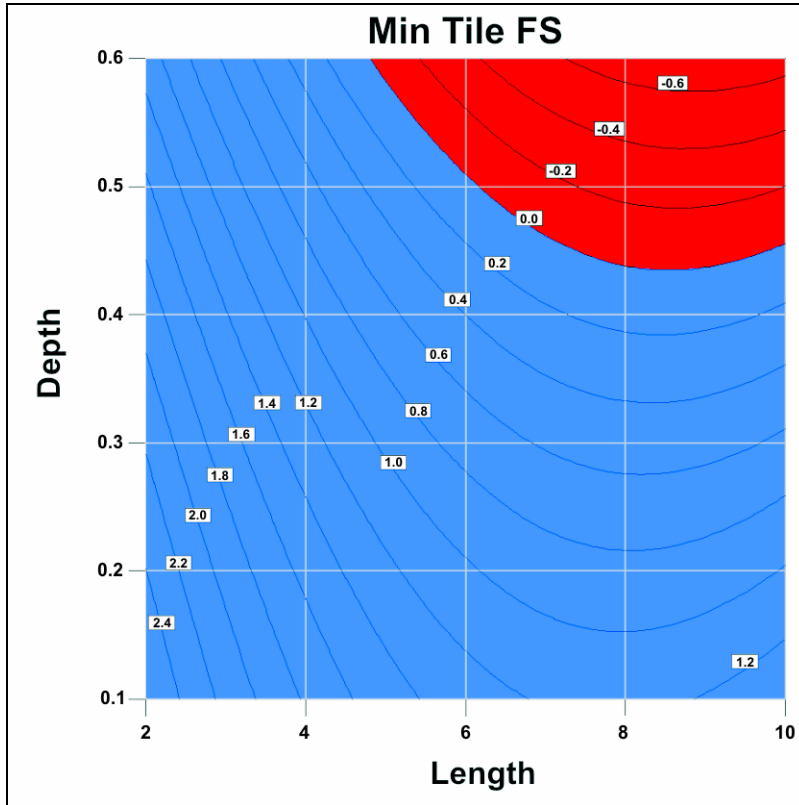


Figure 11. Tile factor of safety contours from data set with many zeros.

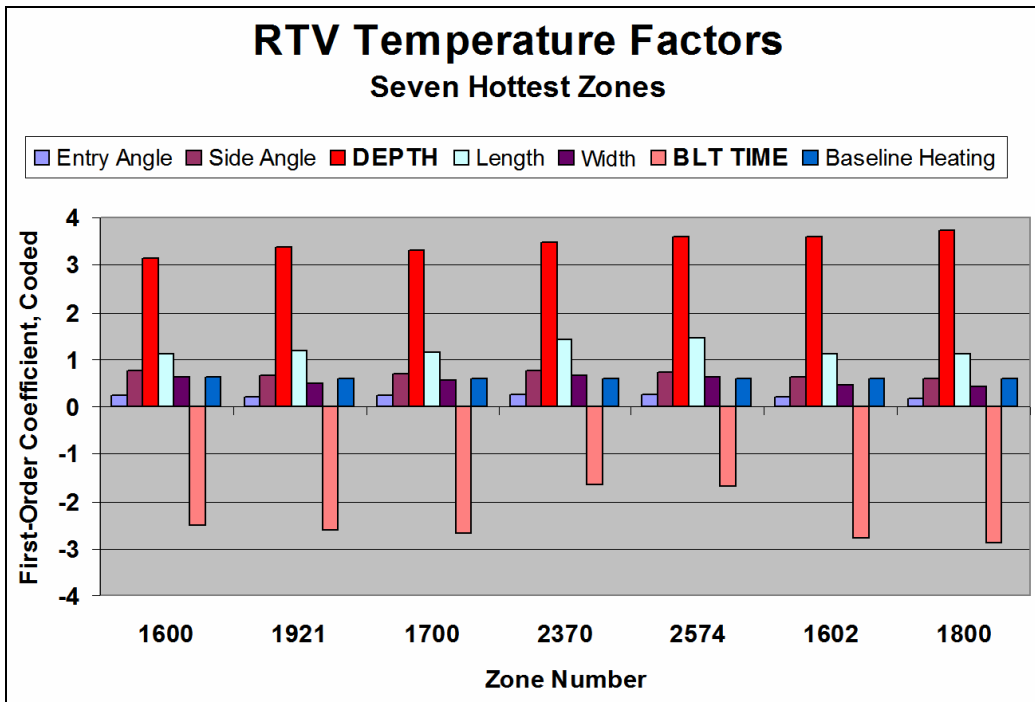


Figure 12. Maximum RTV bondline temperature factors, seven zones with highest average RTV Temperature. Depth and boundary layer transition time are the most sensitive factors.

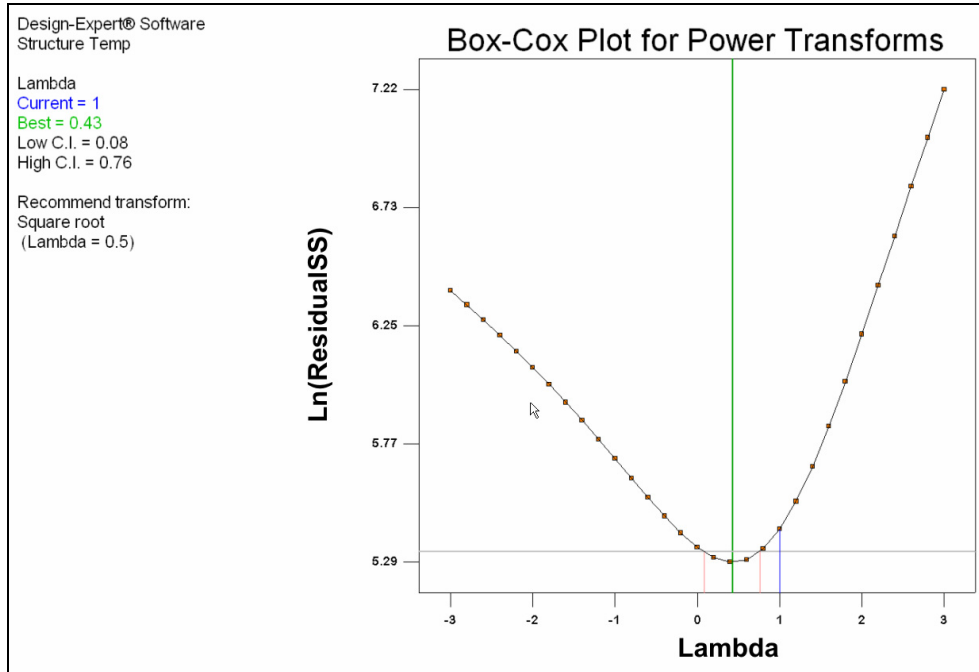


Figure 13. Example Box-Cox power transformation plot. Minimum occurs for lambda near 0.5, implying square root transformation.

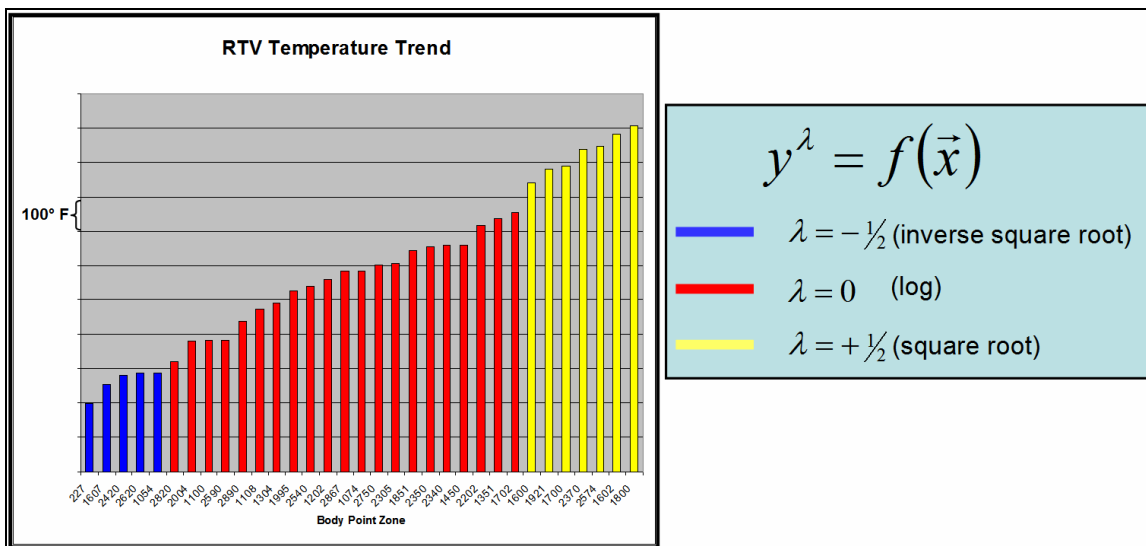


Figure 14. Power transformations to optimize the fit in maximum RTV bondline temperature models exhibit an interesting pattern.