Software Requirements

· · · ·	Appendix G			
NUCLEAR WASTE MANAGEMENT PROGRAM ational boratories	User's Manual Criteria Forr	n	I	rm Number NP 19-1-6 age 1 of 1
	ual contain as appropriate:			
1. Software Name:	CUTTINGS_S			
2. Software Version	5.10			
3. Document Versid	1.00			·····
	532340 Iser's Manual, all items shall be appropriately addressed by the	code spo	onsor s	o that "Yes"
or "N/A" may be checke	ed. Include this form as part of the User's Manual.		•	
	functional requirements (consistent RD) and system limitations?	\boxtimes	Yes	· .
	the mathematical model and , where applicable as based on code functionality?	\boxtimes	Yes	□ N/A
7. Physical and mat based on code fu	hematical assumptions, where applicable as nctionality?	\boxtimes	Yes	🗌 N/A
8. The capabilities a	nd limitations inherent in the software?	\boxtimes	Yes	
9. Instructions that of software?	describe the user's interaction with the	\boxtimes	Yes	
10. The identification	of input parameters, formats, and valid ranges?	\boxtimes	Yes	
11. Messages initiate the user can resp	ed as a result of improper input and how bond?	\boxtimes	Yes	
12. The identification	and description of output specifications and formats?	\boxtimes	Yes	,
13. A description of	any required training necessary to use the software?	\boxtimes	Yes	
14. The identification	of components of the code that were not tested?	\boxtimes	Yes	
	Sponsor (print) Signature	er.		10/22/03 Date
16. James Garner	(IW Gamer			10/22/03
Technical R	eviewer (print)	1		Date
	Manager (print) Signature	·		10/22/03 Date
18. Jennifer Long SCM Coo	rdinator (print) Signature			<u>10/22/03</u> Date

•

Key for check boxes above:

Check **Yes** for each item reviewed and found acceptable Check **N/A** for items not applicable, where applicable as based on code functionality

Dunagan, Sean

From: Sent: To: Subject: Hansen, Clifford W Monday, October 20, 2003 10:26 AM Dunagan, Sean; Chavez, Mario Joseph Sig. authority for CUTTINGS documents

Sean,

In my absence, you have permission to sign CUTTINGS documents as the code sponsor. You may sign the RD, VVP, VD, the ID, the I&C, and the criteria forms for the RD, VVP, VD and the ID.

Cliff Hansen Sandia National Labs Dept 6821 505-234-0103 (Carlsbad) 505-845-0285 (Albuquerque)

R- 10/21/2003

WIPP PA

User's Manual

for

CUTTINGS_S, Version 5.10

DOCUMENT VERSION 1.00

ERMS # 532340

OCTOBER 2003

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1.0 INTRODUCTION

This document is the User's Manual for the WIPP code called CUTTINGS_S, Version 5.10. Using repository-height data from SANTOS via BRAGFLO, CUTTINGS_S estimates the volume of wastes brought to the surface as a result of an inadvertent borehole drilled directly over the WIPP repository so as to penetrate a waste panel. The manual identifies the code, its authors and expert consultants (Section 1). It describes the code's WIPP-PA purposes and functions (Section 2), provides recommended user training (Section 3), outlines the code's theoretical basis and numerical methods (Section 4), its inherent capabilities and limitations (Section 5), describes user interactions (Section 6), input files (Section 7), error messages (Section 8), output files (Section 9), and useful reference materials (Section 10). It provides examples of relevant input, output, and debug files in its Appendices as well as a set of sample test problems to familiarize the user with the run sequence.

It is useful to remember that the present version of CUTTINGS_S's evolved in time through a sequence of WIPP PAs. Design criteria for each PA differed slightly from the set before, and careful readers will detect, especially in the appendices, terminology, test cases, and parameter values such as radioisotope decay data, that serve no apparent purpose in the *present* version of the code. Outdated but working subroutines were normally retained within the code so as to provide versatility.

1.1 Software Identifier:

CUTTINGS_S, a borehole waste-removal code for cuttings spallings.	s, cavings, and	d
CUSP		. •
5.10		
October 2003		
FORTRAN 77 for Open VMS 7.3-1, ES40.		
	spallings. CUSP 5.10 October 2003	CUSP 5.10 October 2003

1.2 Points of Contact:

Sponsor:

Cliff Hansen (505-234-0103) 4100 National Parks Highway Carlsbad, NM 88220 E-mail: <u>cwhanse@sandia.gov</u>

1.3 Code Overview

The CUSP code was written to calculate the quantity of material (in m³) brought to the surface from a radioactive waste disposal repository as a consequence of an inadvertent human intrusion through drilling. The code determines the amount of material removed from the repository by several release mechanisms, including cuttings, cavings and spallings. The CUSP code includes functionality to compute the radioactivity in the released material, including decay of the material to the time of intrusion.

2.0 FUNCTIONAL REQUIREMENTS

2.1 Functional Requirements

- R.1 CUTTINGS_S calculates the amount of repository material brought to the surface due to erosion of the borehole resulting from laminar flow in the drilling fluid.
- R.2 CUTTINGS_S calculates the amount of repository material brought to the surface due to erosion of the borehole resulting from turbulent flow in the drilling fluid.
- **R.3** CUTTINGS_S calculates the amount of repository material brought to the surface due to blowout of the borehole.
- **R.4** CUTTINGS_S calculates the amount of repository material brought to the surface due to gas erosion of the borehole.
- R.5 CUTTINGS_S calculates the amount of repository material brought to the surface due to a stuckpipe.
- R.6 CUTTINGS_S calculates model specific parameter values based on experimental data.
- R.7 CUTTINGS_S calculates the volume of spalled material using a pressure threshold and a distribution of spallings volumes (spall model 3).
- R.8 CUTTINGS_S determines the volume of spalled material using a set of distributions of spalled volumes, calculated for a set of reference repository pressures, by interpolating between distributions to account for current repository pressure (spall model 4).

2.2 **Performance Requirements**

There are no performance requirements for the CUTTINGS_S code.

2.3 Attribute Requirements

There are no attribute requirements for CUTTINGS_S code.

2.4 External Interface Requirements

- R.9 CUTTINGS_S utilizes routines from CAMDAT_LIB, CAMCON_LIB, SDBREAD_LIB, and CAMSUPES_LIB. Consequently it must be linked with these libraries.
- R.10 CUTTINGS_S requires one CDB input file from the BRAGFLO code, CUSP_INP\$BRAGCDB.
- R.11 CUTTINGS_S requires one input file containing preliminary data base information, CUSP_INP\$CDB.
- R.12 CUTTINGS_S requires one input file containing model and site dependent parameters and radionuclide properties, inventories, drilling procedures, and characteristics of the drilling fluid, CUSP_INP\$TXT0.
- R.13 CUTTINGS_S requires one input file identifying input sample vector values that will be used in the analysis, CUSP_INP\$TXT1.
- R.14 CUTTINGS_S generates one output file CUSP_OUT\$DBG, which contains information that is used for comparing with acceptance criteria, and is used only for testing purposes.
- R.15 CUTTINGS_S generates one output file CUSP_OUT\$NVERIFY, which contains information that is used in the functional testing for hand calculations, and is not used in production runs.
- R.16 CUTTINGS_S generates one output CDB (binary) file, CUSP_OUT\$CDB, containing output generated by the code. This output must conform to the format specified in the WIPP PA User's Manual for CAMDAT_LIB (4).
- R.17 If spall model 4 is used, CUTTINGS_S reads spall volume data from a text input file, CUSP_SPL4\$DAT.

2.5 Other Requirements

There are no other requirements for CUSP that require verification.

3.0 REQUIRED USER TRAINING AND/OR BACKGROUND

To exercise CUTTINGS_S, users should have (1) basic knowledge of open VMS, (2) basic facility with Digital Command Language, and (3) an overall understanding of Sandia's CAMDAT database, which is used in virtually every WIPP code (Rechard, 1992; and Rechard et al., 1993). User's should also have (4) access to the WIPP cluster of Compaq Alpha microcomputers with an open VMS 7.3-1 operating system or their functional equivalents.

To manipulate and/or interpret the results of CUTTINGS_S as it is exercised in WIPP PAs, user's should have (1) a basic understanding of the physics of radioactive decay, the mechanics of laminar and turbulent flow of non-Newtonian fluid suspensions in annular conduits, erosion scouring of matrixed materials, flow through permeable media, and of the large-scale deformation and failure of elastic, plastic, and viscoelastic materials (2) a basic understanding of partial-differential equations and integral calculus, as they apply in the mathematical formulation of physical principles and especially of conservation of momentum, energy, and mass*, (3) a generalist's conceptual understanding of numerical methods as they are applied to the numerical solution of the boundary-value problems of mathematical physics and chemistry, and (4) a basic overview understanding of the WIPP PA process, including conceptual models, scenarios, inventories, uncertainty sampling, input-data vectors, and a general familiarity with the files and functions of the WIPP codes that interface with CUTTINGS_S and especially GENMESH, MATSET, POSTLHS, and BRAGFLO (WIPP PA Dept [5 Volumes], 1992).

^{*} Strictly speaking, the decay of radioisotopes is not mass conserving. It requires loss of mass, although the amounts are insignificant for present purposes. Fluid and gas flows associated with CUTTINGS_S are mass conserving.

4.0 DESCRIPTION OF THE MODEL AND METHODS

The discussion of the model and methods given below is but a brief overview of the theory and methods employed by CUTTINGS_S. It is intended for users who have only a cursory interest in the mathematical and fluid dynamical details. A thorough, detailed treatment of the theory and numerics applied in CUTTINGS_S is given in Appendix A of this manual.

4.1 Description of the Model

CUTTINGS_S is a multi-faceted computational procedure to assess the effects of *direct removal* of wastes buried in a WIPP-like nuclear-waste repository. Direct removal is hypothesized to occur as the result of inadvertent penetration of the repository by a borehole drilled in connection with oil or gas exploration at some unknown time in the future (Berglund, 1993). The word "direct" refers to the fact that CUTTINGS_S removals from the repository to the surface occur *at the time* of drilling. Since drilling operations are normally completed on the time scale of weeks, removal of drilled wastes from the repository to the surface (i.e., the accessible environment) takes place entirely within a single timestep of the much-longer-term groundwater transport codes. Hence, the word "direct" should be taken to mean "from the repository to the surface within a single timestep of a typical WIPP groundwater code."

Removal is subdivided according to process, which results in three principal categories, cuttings, cavings, and spallings.

First, the borehole is assumed to penetrate directly through waste containers stored within the repository. The drill bit cuts directly through the waste containers, allowing those repository wastes located on a collision course with the drill bit (called cuttings) to be transported directly and immediately to the surface via the cooling fluid (muddy brine) that is circulated through the borehole during drilling. The volume of repository material removed as cuttings is the portion of the deformed repository cut by the borehole.

Wastes that originally bordered the borehole on its exterior are permitted to erode into the cooling fluid as a result of the fluid's frictional effects on the outer cylindrical surface of the borehole. If shear stresses in the cooling fluid at the wall exceed the shear strength of the matrixed materials located at the wall, erosion will occur and wall material (cavings) will be removed*. As a result, the borehole diameter may increase locally at the level of the repository. Because of that action, the volume of wastes removed from the repository can actually be larger than the volume of the borehole originally cut through the wastes. As the diameter of the borehole increases, the fluid shear stresses at the outer boundary decrease. If they decrease enough that the failure stress of the wall material equals or exceeds the fluid shear stress at the outer boundary of the flow region, then the caving process will terminate. CUTTINGS_S allows for both laminar and turbulent shear flow in the circulating fluids, which, in turn, leads to two

^{*} In tightly matrixed or crystaline media such as the WIPP-site halite formation, exploratory boreholes are usually unlined. Rather, the drilling fluid is changed to saturated salt solution to minimize dissolution at the walls, and the borehole is permitted to support itself until the drill bit reaches granular strata below the halite horizon.



different models to assess the effects of material removal due to erosion. If the cooling-fluid flow is laminar, CUTTINGS_S employs an analytical solution of the modeling equations to evaluate wall stresses and other flow parameters. The analytical solution takes the form of nonlinear integral equations, which CUTTINGS_S evaluates via an iterative numerical procedure. If the flow is turbulent, CUTTINGS_S turns to a simpler empirical model to evaluate the turbulent shear stresses. Transition to turbulence is assumed to occur at a Reynolds number greater than 2100. In nature, the transition to turbulence actually occurs over a range of Reynolds numbers and is not an abrupt process. Normally there is a transitional regime during which the flow is partly laminar and partly turbulent with active and chaotic transitions between the two states. CUTTINGS_S does not attempt to model the transition regime. Rather, it assumes the flow becomes fully and abruptly turbulent when the Reynolds exceeds 2100. CUTTINGS_S employs a corrective procedure to compensate for the effects of that assumption (see Appendix A for details). To date, all WIPP-like runs have required turbulent cooling-fluid flow.

The cuttings/cavings model described thus far is functionally similar to the CUTTINGS model used in previous WIPP PAs (WIPP PA Dept, 1992, Vol 1). However, an important new element has been added to the newer CUTTINGS_S code, namely, the inclusion of a third material-removal mechanism, removal in the form of spallings, which is discussed below.

If the repository is pressurized at the time of penetration, and if the pressure within the repository is high enough (i.e., above the hydrostatic pressure in the cooling-fluid column), repository materials originally surrounding the borehole, but interior to the repository-fluid interface, can be forced to deform inward toward the borehole so as to replace the materials already removed by action of the cutting bit and the erosive brine. The solid fraction of these materials are called spallings*. Depending on the parameters of the problem, spalling deformations can be rapid, or they can be a sluggish, more like a large-amplitude, viscoelastic flow. If repository materials are relatively impermeable, the spalled materials can deform slowly inward so as to press against the rotating drill string. If repository pressures are high enough, frictional forces against the drill string may be high enough to stop the drill from turning. In that event, drillers normally withdraw the drill (vertically) and redrill downward through the spalled materials that clog the borehole. On occasion, that procedure must be repeated more than once. If repository materials are more permeable, the pressurized gas may simply seep through the matrix carrying loose particles with it. Alternatively, if repository pressures are too low, the spallings part of CUTTINGS_S may do nothing at all. CUTTINGS_S may have to be applied more than once during an assessment run, at more than one intrusion time, since multiple borehole intrusions are permissible in WIPP PAs.

The total volume of repository material removed by CUTTINGS_S is the sum over all intrusions of cuttings, cavings, and spallings. Options cuttings and cavings are assumed to operate

^{*} If the repository is wet with brine at time of penetration, which possibility is allowed in WIPP PAs, three phases of material could, in principle, flow toward the borehole, namely: pressurized gases, pressurized brine, and solid wastes. CUTTINGS_S estimates only the amount of solids that are moved into the borehole. Pressurized brine that may be released into the borehole is treated by an independent WIPP code named DBR_BRAGFLO. The gas flow, although crucial to the overall deformational process, transports no radioisotopes and is therfore not assessed.

continuously, whereas spallings operates only when repository pressure conditions permit it (the cut off being 8MPa.). The amount of material removed by each process, during each hypothesized drilling intrusion is reported in CUTTINGS_S's output in terms of an equivalent area. Downstream WIPP codes, namely CCDFGF, use BRAGFLO's repository-height data to convert the equivalent areas to volumes of repository material, and then assign various sampled aged repository contents to convert these material volumes to releases. Thus, CCDFGF evaluates the release effects of striking contact-handled versus remote-handled wastes. The reader is referred to Appendix A for details.

The fluid-flow portion of CUTTINGS_S is based on a non-linear analytical integral solution for laminar helical flow of a non-Newtonian (Oldroyd-type) fluid in an annular pipe, and on empirical equations for turbulent flow in the same geometrical configuration. CUTTINGS_S acquires its parameter data from (i) the input CDB files provided by the WIPP codes MATSET and BRAGFLO, (ii) a user-supplied input text file, and (iii) the carefully controlled and QAed WIPP property database. Discrete values of sampled data are provided by the WIPP sampling code POSTLHS in cases where parameter uncertainty is judged to impact assessments of the repository's performance. CUTTINGS_S input includes the rotational speed of the drill string, cutting-bit diameter, the shear-rate-dependent viscosity and density parameters for the drilling mud, borehole roughness, the compacted repository thickness and porosity, and the effective failure shear strength of the compacted repository material. A few required parameters, namely drilling-mud flow rate and drill-collar diameter, are calculated internally by the code, based on present-day oil-drilling industry guidelines.

4.2 Description of the Modeling Methods

CUTTINGS_S does not endeavor to integrate conservation equations characterizing the flow on a point-for-point, time-for-time basis in search of a detailed space-time history of the removal process. Rather, removal is treated as instantaneous, that is, within a single timestep of the longer-term groundwater codes. Thus, numerical methods are used primarily to determine the limiting values of the physical parameters that represent the final configurations of the various (sampled) drilled systems, which is all that is required to assess radioactive waste removal due to the cuttings, cavings, and spallings that arise during drilling operations. These methods are fairly straightforward, iterative, recursive approximation techniques that involve systems of algebraic equations. They are discussed thoroughly in Appendix A (Sections 2.2.1.1, 2.2.2.1, and 2.4.1.2.1), which is highly recommended to readers who aim to understand this code.

5.0 INHERENT CAPABILITIES AND LIMITATIONS OF THE SOFTWARE

CUTTINGS_S is not presently put forward as a general borehole-mechanics code. Rather, it was developed with WIPP-like conditions and applications firmly in mind. To function properly, it requires inputs from other WIPP PA codes, specifically GENMESH, MATSET, POSTLHS, and BRAGFLO. Thus, although CUTTINGS_S is quite likely to be fully applicable to other kinds of boreholes, other geologies, and other parameter ranges, it is not presently recommended for applications outside the presently-conceived boundaries that define the WIPP, primarily because such applications have not yet been tested. Consequently, no claims are made that CUTTINGS_S should give reasonable results for parameter ranges that are atypical of the WIPP and its immediate environs.

All aspects of CUTTINGS_S are based on the present-day drilling methods and technologies that would apply at a WIPP-like site. CUTTINGS_S has been tested under those conditions and those conditions only. Hence, applications beyond the limits of those assumptions are not recommended at this time.

6.0 USER INTERACTIONS WITH THE SOFTWARE

Before launching into the exercising of CUTTINGS_S, the reader is reminded that CUTTINGS_S is normally exercised as one of a sequence of codes that includes GENMESH, MATSET, and POSTLHS, and requires input data files that originate with the code BRAGFLO. To assist the reader, the code sequence is depicted in figure 1, and examples of the input control files required by GENMESH and MATSET in preparation for a CUTTINGS_S run are given in Appendix B. POSTLHS requires an input control file, but it is empty.

6.1 Exercising CUTTINGS_S Interactively

CUTTINGS_S has been modified in comparison with its parent code, CUTTINGS, so as to remove the possibility of user-interactive runs.

The following is an example of how to execute the CUTTINGS_S code. The following example is the executable that was used in the 1996 CCA PA. To execute CUTTINGS_S from a command line, it is necessary to define a logical symbol that invokes its executable, by typing either of the following two lines at the VMS \$ prompt:

\$CUSP :== "\$WP\$PRODROOT:[CUSP.EXE]CUSPPA96.EXE" <CR>*

or

\$CUSP :== "\$WP\$TESTROOT:[CUSP.EXE]CUSP_TEST_NODBG.EXE"

No matter which definition is used, typing "\$CUSP" and striking the carriage-return key will result in invoking CUTTINGS_S's executable.

Like all WIPP codes, CUTTINGS_S must be made aware of the names of and paths to the suite of input files it requires in order to exercise. It must also be given user-selected names for the files to which it will write its output. In command-line execution mode, CUTTINGS_S will search for files by reviewing its internal list of logical symbols associated with input/output files. Thus, the user must inform CUTTINGS_S which of its list of internal logical symbols corresponds to which input or output file for that specific run. That information is normally transferred by typing (at the VMS \$ prompt that will appear after defining the executable logical) a line such as the following:

\$DEFINE CUSP_INP\$CDB XY:[ABC.DEF.GHI]JKL.MNO

^{*} In this user's manual, VMS commands will always be entered by striking the carriage-return key. That instruction will not be repeated hereafter.



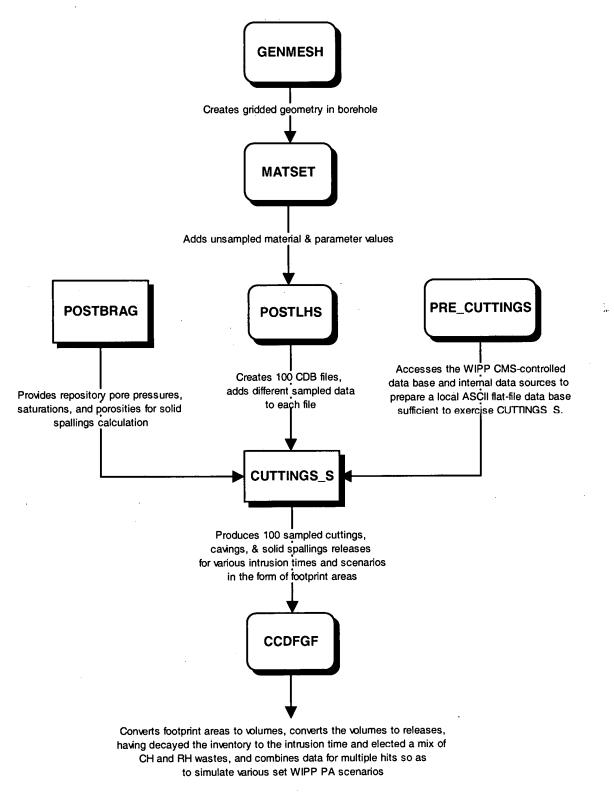


Figure 1: The code sequence for CUTTINGS_S in the 1996 CCA PA.

where XY:[ABC.DEF.GHI]JKL.MNO represents the specific input file the user wishes to associate with the input logical CUSP_INP\$CDB. The path need be included only if the file resides other then in the local working-level directory.

6.2 CUTTINGS_S's Input/Output File Structure

Up to 12 logical symbols may be associated with CUTTINGS_S's various input/output files. Of them, 6 are required and 5 are optional in regulatory runs. Before discussing the nature and function of the 12 possible input/output files, we first list them one-by-one, for ready reference, together with their corresponding logical symbols. They are:

File ID No.	Input/Output File Names	Associated Logical Symbol	Is the file Required or Not?
1. Binary inp	out CAMDAT file (from MATSET)	CUSP_INP\$CDB	Yes
repository	trol file (text) specifying //model parameters, initial es, generic radioisotope database.	CUSP_INP\$TXT0	Yes
	trol file (text) specifying drilling sion parameters	CUSP_INP\$TXT1	Yes
I. BRAGFL Regulator Test runs		CUSP_INP\$BRAGCDB	Yes No
	trol file for Spall Model 4 specifing and their volumes by vector	CUSP_SPL4\$DAT	Yes (spall model 4 only)
6. CUTTING	S_S's binary output .CDB file	CUSP_OUT\$CDB	Yes
	S_S's output debug text file	CUSP_OUT\$DBG	Yes
. Text file f	or verification testing	CUSP_OUT\$NVERIFY	No
). Radioisot output te	ope inventory history (in CURIES) ext file	CUSP_OUT\$IHISTO	No
(EPA 40	zed-release-history output text file CFR191, Subpt B) SCII format	CUSP_OUT\$NHISTO	No
Long Medium	story plotting text file for: half-life radioisotope decay half-life radioisotope decay half-life radioisotope decay	CUSP_OUT\$PLT1 CUSP_OUT\$PLT2 CUSP_OUT\$PLT3	No No No
12. Output t	ext file for transfer to	CUSP_OUT\$ICTRN	No

the WIPP code CCDFPERM

The first five files of the above list are CUTTINGS_S's input data and control files. They are described briefly in Section 7.0 of this user's manual. The text-format input files include specific examples printed out in full in the appendices of this user's manual. Appendix C is included to assist the user in understanding the format used in CUTTINGS_S's input control files (file numbers 1 and 2).

The last seven files of the above list are CUTTINGS_S's output files. They are described briefly in Section 9.0 of this user's manual. Of the required output files, only the debug file is created in text format. One example of a CUTTINGS_S output file is included in Appendix D. It is a debug file. It is annotated with italicized remarks that identify and explain the various groups of data that are reported therein.

6.3 Exercising CUTTINGS_S

Once the logical symbol CUSP has been defined (see Section 6.1), and the necessary input and output files have been properly assigned (see Sections 6.1 and 6.2), CUTTINGS_S may be exercised simply by typing the logical symbol:

\$CUSP

followed by a carriage return. If all goes well, CUTTINGS_S will run and issue a "normal completion" message. If errors are encountered, CUTTINGS_S will abort and issue an error message(s). See Section 8.0 for a listing of the possible error messages CUTTINGS_S can produce.

6.4 Using CUTTINGS_S's PRE_CUTTINGS Mode

Because interactions with the controlled WIPP database require large and unnecessary expenditures of input/output time during production runs, CUTTINGS_S includes a special operational mode designed to expedite data acquisitions during production runs. It is called PRE_CUTTINGS, and in it, the WIPP QAed property database is accessed but once during an entire production run. PRE_CUTTINGS reads a template file that interfaces the WIPP property database, extracting the required values and writing them to an ASCII file that can be read time and again by CUTTINGS_S without subsequent reference to the property database. The ASCII data file thus produced is under strict CMS control.

PRE_CUTTINGS's template file, which is shown in Appendix E, treats three data types. The first type is exemplified by CUTTINGS_S's code parameters, which are discussed in the theoretical Manual (see Appendix A). These values are essential to exercising CUTTINGS_S, but were not entered as part of the WIPP's CCA controlled database. They are discussed in detail in Appendix A. The second type are common property values as taken from the WIPP's CCA controlled database. For example, the datum BOREHOLE:L1 is called. In this example, the colon at mid word identifies the form as a WIPP controlled database property. BOREHOLE

is the material identifier and L1 is the parameter identifier. Using these two keywords, the WIPP property database is scanned until the numerical value associated with that specific parameter is located. CUTTINGS_S then writes to the new template the parameter identifier L1 and the corresponding numerical value that was found in the database.

The third data type in the input template file includes the 10 radioisotope decay chains and the 29 specific radioisotopes that have initial inventories, which are denoted by the word "SAVE." CUTTINGS_S tracks selected data for all of these radioisotopes. When CUTTINGS_S encounters the key word "GENERATE_RADIO", it sorts the list alphabetically and for each unique radioisotope, it queries the property database for atomic weight, half life, EPA release limit, initial contact-handled inventories, and initial remote-handled inventories. The code converts the half life in seconds to half life in years and calculates the specific activation energy, which is a function of atomic weight and half life. The code then writes all the necessary radioisotopic properties in tabular form. Appendix F is an example of the output ASCII text file from PRE_CUTTINGS. Appendix G is a sampled "COM" file to test this portion of the code. Because this functionality of the code is basically data in / data out, it was not assigned a specific test problem.

6.5 Using CUTTINGS_S's CHECK_OUT Mode

CUTTINGS_S has a special mode, called "CUSP\$CHECK_OUT," designed to check the agreement between Model 2's blowout calculation and corresponding experimental results. CHECK_OUT was added so the blowout model could be used for confirmation of experiential results. The CHECK_OUT mode is set by specifying "CUSP\$CHECK_OUT" as TRUE in DEC's Digital Command Language (DCL). Appendix H provides a listing of the DCL nomenclature required to exercise this operational mode. The CUTTINGS_S input file CUSP_INP\$TXT0 provides data for the experimental input, the geometry of the test fixture, and the properties of air. It is also shown in Appendix H. In CHECK_OUT mode, three functions can be performed, as follows:

- 1. The "EXPERIMENT" function is set by including the keyword EXPERIMENT in the CUSP_INP\$TXT1 text file. It allows CUTTING_S to be run as though data is being gathered from an experimental run.
- 2. The "VALIDATE" function is set by including the keyword VALIDATE in the CUSP_INP\$TXT1 text file. It allows CUTTINGS_S to be run over the total range of possible inputs and thereby checks the code for numerical instabilities.
- The "CALIBRATE" function is set by including this keyword CALIBRATE in the CUSP_INP\$TXT1 text file. This function searches for the optimum values of FSE & FGE (see Appendix A) to minimize the error between the experiential results and the computer model. The CALIBRATE function is illustrated as a test problem #6 in the Validation Document.

The input files for these three functions are also shown in Appendix H.

7.0 DESCRIPTION OF CUTTINGS_S'S INPUT FILES

CUTTINGS_S's input files are listed in Section 6.2 together with their corresponding CUTTINGS_S logical symbols. They are file numbers 1 through 4 in that listing.

File 1. CUSP_INP\$CDB is the principal input CDB file and is absolutely required by CUTTINGS_S. It is the file to which CUTTINGS_S's results will be transferred and it is normally provided by an upstream code such as MATSET. In production runs, sampled data from POSTLHS would be fed to this input file. In that case, there would be N input CDB files, one for each of the sampled values where N is the number of samples employed. In test-case mode or debug mode, it is inconvenient to open the editor and hand build an input CDB file for specific input values. However, even in test or debug mode, CUTTINGS_S expects to receive an input CDB file. As a convenience to users, an all-purpose, input CDB file for use with test problems and in debug runs is stored in the configuration-management-system (CMS) portion of the WIPP Alpha cluster. It is named: CUSP_TEST_1.CDB and can be called by typing:

\$LIBCUSP

\$CFE CUSP_TEST_1.CDB

File 2. CUSP_INP\$TXT0 is an ASCII input control file that contains various repository, drill geometry, and material input parameters required by CUTTINGS_S. File 2 also contains the required radioisotope inventory data for the scenario currently under study. In regulatory calculations, these data are taken directly from the controlled WIPP property database. File 2 also provides radioisotope data essential for (1) the Bateman (1910) decay calculation, for (2) calculation of radioactive release values in Curies, and for (3) normalizing calculated releases in terms of the Environmental Protection Agency's (EPA's) 40CFR191 release limits (WIPP PA Dept [Vol 1], 1992). A sample CUSP_INP\$TXT0 file usable for QA test and/or debug problems is available in the CMS portion of the WIPP Alpha cluster. It is named: CUSP_TEST_REPOSITORY.DAT, and it is printed in full in Appendix I of this user's manual. Most of its parameters are specific to the CUTTINGS_S model and to individual radioisotopic properties.

File 3. CUSP_INP\$TXT1 is an ASCII input control file that contains various drilling and intrusion input parameters required by CUTTINGS_S. A sample CUSP_INP\$TXT1 file usable for test and/or debug runs is available in the CMS portion of the WIPP Alpha cluster. It is named CUSP_TEST_#.INP where # is the QA test problem number and varies from 1 through 5. In a normal production run, this file would contain no specific input data other than an intrusion time. Instead, only character variables would appear, identifying the element property name on the input CDB file. Values would then be assigned from the secondary database (see item 4, below). Appendix J is sample printout of a typical production input file. For specific testing or

debugging, this file would have fixed specific numerical values. Appendix K is a listing of CUSP_TEST_2.INP.

This file also identifies which spall model will be used. At the end of the file the model is determined by entering MODEL3 or MODEL4, depending on which model will be used. This is followed by the parameters that are needed for the model. If model three is selected then three parameters will be needed. Model four uses three parameters and one dummy parameter. The parameter PTHRESH is not used in model four. An example of how to call each model and the parameters it requires is below.

MODEL3 PTHRESH WAS_AREA:PTHRESH VOLSPALL WAS_AREA:VOLSPALL REPPRES CAVITY_1 < MODEL4 REPPR_NM PRESGAS REPPRES CAVITY_1 < RNDSPALL DUM5:RNDSPALL PTHRESH DUM5:PTH

File 4. CUSP_INP\$BRAGCDB is the output CDB file from BRAGFLO as it is exercised in the undisturbed or disturbed scenarios. CUTTINGS_S extracts from it (1) the repository porosity at time of decommissioning, and (2) the repository porosity and (3) repository pressure at the specified intrusion time. In regulatory calculations, these data would be provided by a regulatory BRAGFLO run. For QA testing and debug purposes, a fixed BRAGFLO output CDB file has been stored in the CMS portion of the WIPP Alpha cluster. It is named CUSP_TEST_INP_BF.CDB, and it is the BRAGFLO output file used for test problem 1.

File 5. CUSP_SPL4\$DAT is a data file that contains information pertinent to spall model 4. The file supplies the code with pressures and their corresponding data to determine the spall volume. The first line of the file lists the number of vectors in pressure data set. The second line is the number of pressure data sets. The third line lists the different pressures in increasing order separated by blank spaces. The following lines list data relating to each pressure. The first data set is for the first pressure given in line three, the second data set is for the second pressure in line three, and so on. The data sets have three columns of data that are separated by blank spaces. The first column is the vector numbers, the second column is the time (not used), and the third column is the volume. Each of the data sets is followed by a blank line. Each data set must contain the same number of vectors and be in the same order. An example data file is shown below.

5		
2		
100	00000	20000000
1	0	1
2	0	1.25
3	0	1.5
4	0	1.75

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5	0	2.00
bla	nk	
1	0	10
2	0	12.5
3	0	15
4	0	17.5
5	0	20
bla	nk	

8.0 ERROR MESSAGES

Numerous procedural errors can cause CUTTINGS_S to abort. In general, these errors relate to the improper input of data. By noting where the code was when the abort occurred, and what the code was trying to read at the time, these problems can generally be resolved. However, there is not specific advice that can be offered to users in resolving errors beyond the general statement: locate the error and repair it. However, Appendix L, which was extracted primarily from the source code, contains useful operating notes and tips to the user. They are included purely because the code sponsor felt they might be useful to novice users.

To familiarize readers with CUTTINGS_S's error-message patterns, CUTTINGS_S's various error messages are listed below, section for section within the code.

SUBROUTINE ADDMEM(MORMEMR, MORMEMC)

CALL QAABORT(

- .\$ '(ADDMEM) REAL array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM) REAL array memory errors detected ***') CALL QAABORT(
- \$ '(ADDMEM) CHARACTER array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM) CHARACTER array memory errors detected ***') CALL QAABORT('(ADDMEM1) - REAL array size errors detected ***') CALL QAABORT('(ADDMEM1) - REAL memory errors detected ***') CALL QAABORT(
- \$ '(ADDMEM1) CHARACTER array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM1) CHARACTER array memory errors detected ***') CALL QAABORT('(ADDMEM2) - REAL array size errors detected ***') CALL QAABORT('(ADDMEM2) - REAL memory errors detected ***') CALL QAABORT(
- \$ '(ADDMEM2) CHARACTER array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM2) CHARACTER array memory errors detected ***')

CALL QAABORT('(ADDMEM3) - REAL array size errors detected ***') CALL QAABORT('(ADDMEM3) - REAL memory errors detected ***') CALL QAABORT(\$ '(ADDMEM3) - CHARACTER array size errors detected ***') CALL QAABORT(\$ '(ADDMEM3) - CHARACTER array memory errors detected ***')
SUBROUTINE BRGCDB(MAXC, MAXQ, IQ, LQ, Q, C, IERRDB)
IF(IERRDB.NE.0)CALL QAABORT('(BRGCDB) - reading input CAMDAT '// \$ 'sizing parameters ***') IF(IERRDB.NE.0)CALL QAABORT('(BRGCDB) - reading input CAMDAT '// \$ 'TITLE ***')
IF(IERRDB.NE.0)CALL QAABORT('(BRGCDB) - reading number of input'// \$ 'CAMDAT analysis variables ***')
SUBROUTINE DRILL(IWASFLG, IHIT, XTINT, DBDIAM, TUNITS, XNEWAREA)
CALL QAABORT('(DRILL) - Negative ROUTER %%%')
SUBROUTINE GETVAL(NOUTFL, NUQPRP, NELBLK, VARNAM, IASPRP, \$ XMATPR, NAMELB, NMATPR, XVAL)
CALL QAABORT('(GETVAL) - ILLEGAL I.C. interpolation '// + 'variable value range name ***')
INTEGER FUNCTION GINDX(RADNAM, IDMTRL)
INTEGER FUNCTION GINDX(RADNAM, IDMTRL) CALL QAABORT('(GINDX) - cannot locate nuclide index in '// \$ 'generic radioisotope data base (GRDB) ***')
CALL QAABORT('(GINDX) - cannot locate nuclide index in '//
CALL QAABORT('(GINDX) - cannot locate nuclide index in '// \$ 'generic radioisotope data base (GRDB) ***')
CALL QAABORT('(GINDX) - cannot locate nuclide index in '// \$ 'generic radioisotope data base (GRDB) ***') INTEGER FUNCTION GJNDX(RADNAM, NUCNAM) CALL QAABORT('(GJNDX) - cannot locate nuclide index in '//

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\$ 'ype using SDBREAD_LIB routine: SDBQERRY ***')

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PROGRAM CUSP_MAIN

IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - Setting up DYNAMIC'// \$ ' base arrays ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - Opening input '// \$ ' CAMDAT ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - Setting up DYNAMIC'// 'base arrays ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP MAIN) - Opening input '// ' CAMDAT ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// \$ ' input text file 1 ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// ' input parameter text file ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// ' input parameter binary file ***') \$ IF(IERRDB.NE.0) CALL QAABORT ('(BRGCDB) - Check BRAG CDB ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - Opening output'// \$ 'CAMDAT and SCRATCH files ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - Closing input '// \$ 'CAMDAT ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - Trouble closing'// \$ 'output cdb-scratch & CAMDAT files, etc... ***')

SUBROUTINE MEMADJ(MORMEM, MNUSED, Q)

CALL QAABORT('(MEMADJ) - Invalid REAL dynamic array '// 'indices detected ***')
CALL QAABORT('(MEMADJ) - REAL DYNAMIC MEMORY '//
'ALLOCATION ERRORS ***')
CALL QAABORT('(MEMADJ2) - Invalid REAL dynamic array '//
'indices detected ***')
CALL QAABORT('(MEMADJ2) - REAL '//
'DYNAMIC MEMORY ALLOCATION ERRORS ***')
CALL QAABORT('(MEMADJ3) - REAL '//
'DYNAMIC MEMORY ALLOCATION ERRORS ***')

+

SUBROUTINE MEMADJC(MORMEMC, MNUSED, C)

CALL QAABORT('(MEMADJ) - Invalid CHARACTER dynamic array '// 'indices detected ***')

CALL QAABORT('(MEMADJC) - CHARACTER DYNAMIC MEMORY '// + 'ALLOCATION ERRORS ***')

CALL QAABORT('(MEMADJC2) - Invalid CHARACTER dynamic array '//

'indices detected ***')
 CALL QAABORT('(MEMADJC2) - CHARACTER '//
 'DYNAMIC MEMORY ALLOCATION ERRORS ***')
 CALL QAABORT('(MEMADJC3) - ### CHARACTER ### '//
 'DYNAMIC MEMORY ALLOCATION ERRORS ***')

SUBROUTINE PARBINO

CALL QAABORT(' (PARBIN0) - PARBIN0 error input parameters ***')

SUBROUTINE PARTXT0

IF (IERR .GT. 0) CALL QAABORT(\$ ' (PARTXT0) - PARTXT0 error input parameters ***')

SUBROUTINE PREPRO(FIDB, FTXT0, FBIN0, FTXT1, FODB, FBRG)

IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening PRE_CUTTINGS'// \$ ' input template file ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening PRE_CUTTINGS'// ' input desc_0500 file ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening PRE_CUTTINGS'// \$ ' input text file 1 ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// 'output ascii SDB file ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// ' output binary SDB file ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// \$ ' input text file 0 ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// \$ ' input text file 1 ***') IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// ' output verify file ***') \$ IF(IERR.NE.0)CALL QAABORT('(CUSP_MAIN) - opening CUTTINGS'// \$ ' output plot file ***')

SUBROUTINE PROCDB(MAXC, MAXQ, IQ, LQ, Q, C) IF(IERR.NE.0)CALL QAABORT('(PROCDB) - reading input CAMDAT '// \$ 'sizing parameters ***') IF(IERR.NE.0)CALL QAABORT('(PROCDB) - reading input CAMDAT '// \$ 'TITLE ***') IF(IERR.NE.0)CALL QAABORT('(PROCDB) - reading number of input'// \$ 'CAMDAT analysis variables ***') SUBROUTINE PROSDB(MAXC, MAXQ, IQ, LQ, Q, C, IWASFLG, QAINFO) IF (IABORT .NE. 0) & CALL QAABORT('(CUSP_PROSDB) - finding data in the INGRES SDB'// & ' database ***') REAL*4 FUNCTION QUANKG(DT, LPOINT, NRILC, NLEFT, RADCHN, NGRAD, IDMTRL, HALFY, AWT, XMOLES) \$ CALL QAABORT('(QUANKG) - cannot find all radioisotopes in '// 'local chain in GRDB rad. names array ***') \$ SUBROUTINE REABINO(NNIC, RADCHN, IDMTRL, NUCNAM, MAPSAV, SAVNAM, HALFS, HALFY, AWT, \$ \$ ACTCNV, EPAREL, CIPMSQ, MAPINV, IWASFLG) CALL QAABORT(' (PARBINO) - PARBINO error input parameters ***') SUBROUTINE REASDB (DBNAME, CALC_NAME, IDMTRL, IDPRAM, MEDIAN, & IABORT) CALL QAABORT('(CUSP_PROSDB) - extracting '// ' idmtrl, idpram ***') \$ SUBROUTINE REATXTO(NNIC, RADCHN, IDMTRL, NUCNAM, MAPSAV, SAVNAM, HALFS, HALFY, AWT, \$ \$ ACTCNV, EPAREL, CIPMSQ, MAPINV, IWASFLG) IF (ICHECK .GT. 11111) & CALL QAABORT('(REATXT0) - read errors radionuclide input***') **Information Only**

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IF(IERR.GT.0)CALL QAABORT('(REATXT0) - read errors ***') CALL QAABORT('(REATXT0) - Repeating/Duplicate radio'// 'isotope names in chain ***') CALL QAABORT('(REATXT0) - Illegal names in "SAVE" '// 'radionuclides list %%%')	
SUBROUTINE REATXT1(IASPRP, NAMELB, NMATPR, XMATPR, & DIAMMOD, TUNITS, PORO, TCLOUT, PRESSR, & PARTDIA, SATGASX, CEMSIG, FGE, FSE, & FCE)	TINT, 1
CALL QAABORT('(REATXT1) - ILLEGAL drilling bit diamet'// & 'er specified (<=0) ***') IF(IERR.GT.0)CALL QAABORT('(REATXT1) - read errors ***')	
SUBROUTINE RECHEA(Q, C, IDEBLK, NUMPRP, IASPRP, \$ XMATPR, QAINFO, NAMELB, NMATPR)	. 1 ,
IF(IERR.NE.0) \$ CALL QAABORT('(RECHEA) - reading material ID and names array IF(IERR.NE.0) \$ CALL QAABORT('(RECHEA) - reading PROPERTY NAMES array *	
 IF(IERR.NE.0) CALL QAABORT('(RECHEA) - reading PROPERTY values ***') IF(IERR.NE.0)CALL QAABORT('(RECHEA) - reading QA array ***') IF(IERR.NE.0)CALL QAABORT('(RECHEA) - obtaining input CAMDA 'no. of time steps ***') 	
SUBROUTINE RGRDB(IDMTRL, HALFS, HALFY, AWT, ACTCNV, E	PANORM) 1
CALL QAABORT('(RGRDB) - read errors ***') CALL QAABORT('(RGDB) - counting entries in GRDB ***')	
SUBROUTINE SCASDB(FSDB, NENTRY, ID1, ID2, ID3, \$ ID4, ID5, ID6, ID7, IDENT)	1
CALL QAABORT('(SCASDB) - Bad SDB initialization ****') CALL QAABORT('(SCASDB) - Error calling SDBREAD_LIB routine'/ ': SDBCASES ***') CALL QAABORT('(SCASDB) - Error calling SDBREAD_LIB routine'/ ': SDBINFO ***')	
Information Only	

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CALL QAABORT('(SCASDB) - Error calling SDBREAD_LIB routine	·//
\$ ': SDBNINFO ***') CALL QAABORT('(SCASDB) - keyword IDMTRL missing on SDB '/	/
\$ 'header description ***') CALL QAABORT('(SCASDB) - keyword IDPRAM missing in SDB '//	
\$ 'header description ***')	
CALL QAABORT('(SCASDB) - keyword IDSCAL missing in SDB '// * 'header description ***')	, , <i>,</i>
CALL QAABORT('(SCASDB) - keyword MEDIAN missing in SDB '// \$ 'header description ***')	۱ ۰ ۰
CALL QAABORT('(SCASDB) - keyword LOWRNG missing in SDB 'header description ***')	2//
CALL QAABORT('(SCASDB) - keyword HIRNGE missing in SDB '//	,
\$ 'header description ***') CALL QAABORT('(SCASDB) - keyword UNITS missing in SDB '//	$\frac{\mu^{2}}{\mu^{2}}$
\$ 'header description ***')	
 \$ QAINFO, ISETUP) IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IEBB GT 0)CALL QAABOBT(
IF(N.NE.NGRAD)CALL QAABORT(· · · · · · · · · · · · · · · · · · ·
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(, NMATPR, 1
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for sizing parameters ***') SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB & IDMTRL, OUTMAT) IF (MULTP .NE. JULTP) CALL QAABORT(, NMATPR, 1
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for sizing parameters ***') SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB & IDMTRL, OUTMAT) IF (MULTP .NE. JULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 MULTP gt NULTP ***')	, NMATPR, 1
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for sizing parameters ***') SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB & IDMTRL, OUTMAT) IF (MULTP .NE. JULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 MULTP gt NULTP ***') IF (MULTP .GT. NULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 order of multiple hits wrong ***')	, NMATPR, 1
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for sizing parameters ***') SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB & IDMTRL, OUTMAT) IF (MULTP .NE. JULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 MULTP gt NULTP ***') IF (MULTP .GT. NULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 order of multiple hits wrong ***') IF (INUM .GT. NRMS) CALL QAABORT(& ' (SCATXT1) - SCATXT1 INUM gt NRMS ***')	, NMATPR, 1
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for sizing parameters ***') SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB & IDMTRL, OUTMAT) IF (MULTP .NE. JULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 MULTP gt NULTP ***') IF (MULTP .GT. NULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 order of multiple hits wrong ***') IF (INUM .GT. NRMS) CALL QAABORT(& ' (SCATXT1) - SCATXT1 INUM gt NRMS ***')	, NMATPR , 1
IF(N.NE.NGRAD)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for generic sizing ***') IF(IERR.GT.0)CALL QAABORT(\$' (SCATXT0) - SCATXT0 input text for sizing parameters ***') SUBROUTINE SCATXT1(IWASFLG, IASPRP, XMATPR, NAMELB & IDMTRL, OUTMAT) IF (MULTP .NE. JULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 MULTP gt NULTP ***') IF (MULTP .GT. NULTP) CALL QAABORT(& ' (SCATXT1) - SCATXT1 order of multiple hits wrong ***') IF (INUM .GT. NRMS) CALL QAABORT(& ' (SCATXT1) - SCATXT1 INUM gt NRMS ***') 99 IF(IERR.GT.0)CALL QAABORT(, NMATPR, 1

SUBROUTINE WRCHEA(IWASFLG, Q, CH, IASPRP, NMATF \$ QAINFO, LDUMP, RDUMP, KOMAT, NUCNAM, \$ SAVNAM, MAPSAV)	'R,	1	
IF(IERR.NE.0)CALL QAABORT(\$ '(WRCHEA) - ADDING "DENSITY" as a CAMDAT property ***') IF(IERR.NE.0)CALL QAABORT(
\$ '(WRCHEA) - ADDING "DOMEGA" as a CAMDAT property ***') IF(IERR.NE.0)CALL QAABORT(
\$ '(WRCHEA) - ADDING "ABSRO" as a CAMDAT property ***')			
IF(IERR.NE.0)CALL QAABORT(
\$ '(WRCHEA) - ADDING "TAUFAIL" as a CAMDAT property ***') IE(IERR NE 0) CALL: CAABORT(х ^а 2 т		
IF(IERR.NE.0)CALL QAABORT(\$ '(WRCHEA) - ADDING "VISCO" as a CAMDAT property ***')	• • • •		• •
IF(IERR.NE.0)CALL QAABORT(. :		ç İ
\$ '(WRCHEA) - ADDING "YLDSTRSS" as a CAMDAT property ***')	÷,		
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Trouble writing '// \$'CUTTINGS SDB QA record to CDB-scratch file ***')	• .		٠
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Trouble writing '//			- 1
\$'CUTTINGS QA record to CDB-scratch file ***')			
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Writing "HEADER" '// \$'section to the final output CAMDAT file ***')	•		
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//	:		
\$ 'NEW GLOBAL variable names (AREA_C) to CDB ***')			
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//	-		,
\$ 'NEW GLOBAL variable names (AREA_S) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//			
\$ 'NEW GLOBAL variable names (AREA_T) to CDB ***')			.,
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//			, te
\$ 'NEW GLOBAL variable names (VOL_C) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//	5. 1		
\$ 'NEW GLOBAL variable names (VOL_S) to CDB ***')			
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//			
* 'NEW GLOBAL variable names (VOL_T) to CDB ***')	•		,
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// \$ 'NEW GLOBAL variable names (DRILDIAM) to CDB ***')			
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//			
* 'NEW GLOBAL variable names (NUM_INTR) to CDB ***')			-
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// \$ 'NEW GLOBAL variable names (POROSITY) to CDB ***')			:
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//			
* 'NEW GLOBAL variable names (HFINAL) to CDB ***')			
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// \$ 'NEW GLOBAL variable names (PRESGAS) to CDB ***')			
IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'//			

'NEW GLOBAL variable names (PRESBRIN) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// 'NEW GLOBAL variable names (SATGAS) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// 'NEW GLOBAL variable names (SATBRIN) to CDB ***') S IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// NEW GLOBAL variable names (POROSITY) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// 'NEW GLOBAL variable names (HFINAL) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// 'NEW GLOBAL variable names (PRESGAS) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// 'NEW GLOBAL variable names (SATBRIN) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// \$' NEW GLOBAL variable names (NUCNAM C_) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// \$' NEW GLOBAL variable names (NUCNAM S_) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRCHEA) - Error writing'// \$' NEW GLOBAL variable names (NUCNAM T_) to CDB ***') SUBROUTINE WRPLFIL(IWASFLG, MAPSAV, HALFY, NUCNAM, RADHIS, 1 \$ TINT, XRAD1, XRAD2, XRAD3, TMPNM1, TMPNM2, TMPNM3) \$ CALL QAABORT('(WRPLFIL) - Exceeded maximum history plott'// \$ 'ing file "width" ***') C, TIME, HIFLAG, 1 SUBROUTINE WRTANA(IWASFLG, Q, \$ AINTC, DRILDIA, NUCNAM, MAPSAV, SUMCI, \$ XPORO, XPRGS, XPRBR, XSATG, KULTP, KH) IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (AINTC(1,KH)) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (AINTC(2,KH)) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (AINTC(3,KH)) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (VOL_C) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (VOL_S) to CDB ***')

- IF(IERR.NE.0)CALL QAABORT('(WRTANA) Error writing'//
- \$ 'NEW GLOBAL variable values (VOL_T) to CDB ***')

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IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (DRILDIA) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// 'NEW GLOBAL variable values (NUM_INTR) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// NEW GLOBAL variable values (POROF) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// 'NEW GLOBAL variable values (HFINAL) to CDB ***') S IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// NEW GLOBAL variable values (PRESGAS) to CDB ***') \$ IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// NEW GLOBAL variable values (PRESBRIN) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (SATGAS) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// 'NEW GLOBAL variable values (SATBRIN) to CDB ***') S IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// NEW GLOBAL variable values (POROF) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// 'NEW GLOBAL variable values (HFINAL) to CDB ***') S IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// NEW GLOBAL variable values (PRESGAS) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// NEW GLOBAL variable values (SATBRIN) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing'// \$ 'NEW GLOBAL variable values (SUMCI) to CDB ***') IF(IERR.NE.0)CALL QAABORT('(WRTANA) - Error writing a completed'// \$' time-step and advancing to next step /DBOSTEP ***')

INTEGER FUNCTION INEXT(INDEX)

WRITE (6,'(" INCORRECT WORD TYPE WAS FOUND, INEXT")') WRITE (6,'(" WORD CMD ",A, & "INTEGER ",I20, & "RNUMBER ",E16.8)') C18, NUM1, RNUM1

Information Only

STOP ' ERROR INEXT'

REAL*4 FUNCTION RNEXT(INDEX)

WRITE (6,'(" INCORRECT WORD TYPE WAS FOUND, RNEXT")')
WRITE (6,'(" WORD CMD ",A,
& "INTEGER ",I20,
& "RNUMBER ",E16.8)') C18, NUM1, RNUM1

STOP ' ERROR RNEXT'

SUBROUTINE ANEXT (INDEX, AVALUE, NP)

WRITE (6,100)

100 FORMAT(" NONE REAL NUMBER WAS FOUND, ANEXT") STOP ' ERROR ANEXT'

WRITE (6,130) 130 FORMAT(" NO END OF ARRAY, ANEXT") STOP ' ERROR ANEXT'

SUBROUTINE SOLID_OUT (DIA, SATBRIN, HF, PG, PARTDIA, PORO,

С

IF (PG .LT. PSUF) THEN WRITE(NOUTFL, & '(/" **** REPOSITORY PRESSURE LESS THAN ATMOSPHERIC")') STOP ENDIF

С

9.0 DESCRIPTION OF CUTTINGS_S'S OUTPUT FILES

CUTTINGS_S's output files are listed above in Section 6.2, together with their identification numbers and the logical symbols CUTTINGS_S associates with them. The output files are file numbers 7 through 13. Each of them is briefly described below. Of the required files, the only user-readable output appears in the debug file, an annotated example of which is given in Appendix D.

File 7. CUSP_OUT\$CDB is the principal CUTTINGS_S output CDB file. It contains the calculated total volume (in terms of an areal footprint) of all radioactive materials, isotope by isotope, brought to the surface by the combined actions of cuttings, cavings, and spallings. Being a CDB file, it is binary and therefore not directly human readable.

File 8. CUSP_OUT\$DBG is an ASCII output file that echoes the various input files, the input parameter values, and includes a summary of the output calculations given in file 7.

NOTE: There are three run-time debug flags that cause changes in the amount of output written to the DBG output file. They are defined as follows:

DEBUG	= <logical></logical>	for debugging dynamic memory allocation.
DBRAG	= <logical></logical>	for debugging the BRAGFLO CDB file
NOECHO	= <logical></logical>	to cancel echoing of input data to the DBG file

These run-time flags may be invoked and set by typing the variable, for example: DBSPALL on the command line. Alternatively, the user may type the line: \$DEFINE CUSP\$DPSPALL TRUE prior to the execution of CUTTINGS_S.

File 9. CUSP_OUT\$NVERIFY is an optional ASCII output file of intermediate results that is used during developmental phases to verify CUTTINGS_S's calculations against other independent calculations, normally performed by hand.

Files 10 through 11 are optional output files that are normally forwarded to system plotting packages for diagnostic and developmental purposes. See the summary table above for a listing of their logical symbols. Their use in regulatory calculations is not contemplated. Hence, these logicals will either remain undefined or will be defined as NULL or CANCEL, in which case the code will not write to them.

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APPENDIX A: A REFERENCE MANUAL ON THE THEORY AND NUMERICAL METHODS UNDERLYING THE WIPP CODE CUTTINGS_S (BY J. W. BERGLUND)

This appendix describes the theoretical basis of the CUTTINGS_S computer code which was written to calculate the quantity of solid radioactive material (in curies) brought to the surface from a radioactive waste disposal facility as a consequence of an inadvertent drilling intrusion. The code determines the amount of material removed by several release mechanisms, and decays the releases to the intrusion time. The report describes the background, assumptions, equations, and variables governing the releases and the methods used to solve them. The code was written specifically for the computations required for the Performance Assessment of the Waste Isolation Pilot Project (WIPP) but with due care can be used for other radioactive waste repositories.

1.0 Introduction

The Waste Isolation Pilot Plant (WIPP), located in southern New Mexico, is the first planned, mined geologic repository for transuranic wastes generated by U.S. defense programs. WIPP is currently being evaluated to assess compliance with the requirements of EPA 40 CFR 191 Subpart B. Briefly, this requirement, promulgated by the Environmental Protection Agency, limits the amount of radioactive material that can be released to the accessible environment over a 10,000-year regulatory period.

Of the possible pathways for release during this period, one of the most important is that caused by the inadvertent penetration of a waste storage room by an exploratory drill bit. The current performance assessment model relies on the assumption that future drilling techniques will be similar to those in use today. The validity of this assumption is unknown but is necessary to provide a basis on which predictions of release can be estimated. Assuming that current, standard drilling practices for gas and oil are used, mechanisms governing the direct removal of radioactive waste can be identified and governing equations developed and solved. The following report summarizes the current understanding of the processes related to the direct removal of wastes, presents the governing mechanisms, and describes the solution of the model equations.

1.1 Background

The WIPP repository will consist of a number of excavated waste disposal rooms located in bedded halite (salt) approximately 650m below the ground surface in southeast New Mexico. Most excavated rooms will be approximately 91 m long, 10 m wide and 4 m high. Transuranic waste packaged in 55-gallon drums or standard waste boxes will be placed in each room and backfilled primarily with crushed salt.

After the WIPP repository is filled with waste and sealed, the waste is expected to be slowly compacted vertically by salt creep from an original waste room height of 4 m to a compacted height of 0.5 to 2 m within 100 to 200 years. The overburden (vertical) stress acting on the waste will also increase to the lithostatic stress (~14.8 MPa) during this period. The waste in its unmodified form will consist of a mixture of contaminated organic (e.g., cloth, wood, rubber,

plastics) and inorganic (e.g. metals, glass) materials. After placement in the mined (salt) repository, the waste will be compacted by creep closure of the surrounding salt and in addition may become exposed to brine. The exposure of the metallic waste to brine is expected to cause corrosion of the metals and as a by-product will generate gas (H₂). Additional gas will be generated by the biodegradation of the organic materials in the waste inventory. The gas volumes generated by corrosion and biodegradation are expected to increase continuously for hundreds of years and the pore pressure may reach and possibly exceed the lithostatic overburden stress. During this time the repository is also expected to expand under the influence of the elevated gas pressure.

At some time within the 10,000-year regulatory period, it is probable that one or more exploration boreholes will be drilled into and through the vertically compacted waste and some of the waste will be carried to the surface as a direct result of the drilling process. The volume of waste removed to the ground surface will depend upon the physical properties of the compacted, decomposed wastes, the drilling procedures used, and the pore pressures encountered. Because of radioactive decay, the radioactivity of nuclides in the removed waste (in curies) will also depend upon the time of intrusion.

1.2 Current Drilling Practices

In considering the problem of the direct releases caused by drilling it is important to understand the procedures currently utilized in drilling a gas or oil well. The following description is tailored to practices in use near the WIPP site. Typically, independent drilling companies respond to requests for fixed price bids from oil companies to provide drilled and cased holes of a specified diameter to a specified depth at a location leased by the oil company. The bids are competitive and generally the low bidder is selected. As a result of this bidding process, profit margins for the winning driller are governed by the effectiveness of the driller in maintaining efficient drilling parameters and by solving unforeseen difficulties that may arise quickly and safely. The principal goal of the driller is to "make hole" and meet contract requirements. If this is accomplished the drilling operation will be profitable.

Within the Delaware Basin near the WIPP site, gas and oil wells are started by clearing the site, and drilling a shallow hole (~40') to house a conductor pipe. The conductor pipe is set in cement and serves to prevent surface sands from sloughing into the wellbore during later drilling. Below the conductor, drilling is continued to 300 - 600 feet (to top of salt section) using a large diameter (17-26") drill bit and another steel casing is set. This "surface" casing is cement grouted into the hole and a blowout preventer tree is attached to the casing at the surface. Drilling below the surface casing (in the salt section) uses a drilling mud that consists of a saturated brine solution to prevent dissolution of the salt. Drilling through the salt section proceeds at rates from 50 to 100 feet/hour depending on bit diameter. If higher rates are attempted by increasing the weight on bit (WOB) there is the danger of producing a crooked hole. While in the salt section, drilling mud (brine) is supplied from a large plastic lined reserve pit dug in the ground with a surface area of approximately 4000 ft². Drilling mud is pumped from the reserve pit down through the drillpipe and drillbit and up the annulus formed by the drillstring and drilled hole. The drilling mud and drill cuttings are returned directly to the reserve pit where the cuttings settle out. While drilling in the salt section no formal attempt is made to monitor the character of the cuttings or the fluid volume of the reserve pit. A gas analyzer is not attached to the returns until the hole is much deeper than the depth of the WIPP repository. The drillstring generally consists of a drillbit attached to approximately 20, 30 foot long drill collars followed by multiple sections of 30 foot long 4 1/2 " diameter drillpipe. One to four drillbit changes may be necessary when drilling through the salt section. This process requires the removal of the entire drillstring (tripping). Once the drill has passed the salt section (~4500 ft) the hole is again cased and set with cement. At this point, if the repository had been penetrated no further contact would occur between the drilling mud and the exposed WIPP waste.

2.0 Release Processes

Three separate physical processes are assumed to influence the quantity of waste brought to the ground surface as the result of the inadvertent penetration of the repository by an exploratory drill bit. These are:

*Cuttings - waste contained in the cylindrical volume created by the cutting action of the drill bit passing through the waste.

*Cavings - waste that erodes from the borehole in response to the upward-flowing drilling fluid within the annulus, and

*Spallings - waste introduced into the drilling fluid caused by the release of waste-generated gas escaping to the lower-pressure borehole. This requires a repository gas pressure that exceeds the hydrostatic pressure of the drilling mud.

In the above and in the following it is understood that usage of the word "waste" also includes any backfill or other materials that may have been added to the actual waste material.

Spallings can be further subdivided into three regimes that are dependent upon the state of waste permeability and gas pore pressure at the time of intrusion. These are:

**Blowout - the direct release of waste to the surface in waste decomposition gas that has cleared the borehole annulus of drilling mud and is flowing freely to the surface.

**Gas Erosion - low permeability waste that is pressed against the drillstring due to stresses from escaping decomposition gas and is eroded by the flowing drilling mud.

**Stuck Pipe - low permeability waste that is pressed against the drillstring sufficiently hard to prevent normal drilling. This occurs at high gas pressures.

These release processes and when they are active are depicted graphically in Figure 1.

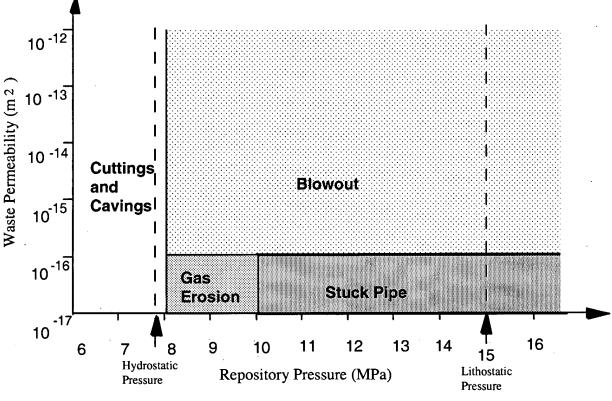


Figure 1. Release process zones

The release process zone boundaries are based on waste permeability and repository gas pressure and will be discussed in the model descriptions for the zones. Detailed descriptions of the release process zones, their principal assumptions, and mathematical formulation will be discussed in the following sections.

2.1 Cuttings

This is the simplest component of the direct release model and is assumed to occur for repository pressures less than the hydrostatic pressure of the drilling mud. This pressure is approximately 8 Mpa (Figure 1). For a gauge borehole, the volume of cuttings removed and transported to the surface is equal to the product of the drill bit area and the drill depth. Thus, to estimate the total solid volume of waste removed due to the cutting action of the drill bit , it is only necessary to know the compacted repository height (H), the porosity at the time of intrusion (φ), and the drill bit area (A).

$$Volume = AH(l - \varphi) \tag{1}$$

The cuttings volume calculated in this manner is a lower bound to the total quantity of waste removed by drilling. In the CUTTINGS_S code the waste is assumed to be uniformly distributed throughout the disposal region. Thus the actual computation for release requires only the drill bit area and the waste curie content per unit area. The actual computation of cuttings release in the CUTTINGS_S code is computed within the module for cavings. This description follows.

2.2 Cavings

The cavings component of direct surface release consists of that quantity of waste material that is eroded from the borehole wall by the action of the flowing drilling fluid after a waste disposal room is penetrated. The erosion process is assumed to be driven solely by the shearing action of the drilling fluid (mud) on the waste as it moves up the borehole annulus. For cavings it is further assumed that the repository is not pressurized by either brine or gas, i.e. the repository pressure is less than the hydrostatic pressure caused by a brine column to the surface (Figure 1). Gas pressurization effects are included in other release processes (spall) and will be discussed later.

In the annulus formed by the collars or drill pipe and the borehole wall, the flow of the drilling fluid has both a vertical and rotational component. Within this helical flow pattern, shear stresses are generated by the relative motion of adjacent fluid regions and by the action of the fluid on the borehole wall. In this model, it is assumed that if the fluid shear stress at the wall exceeds the effective shear resistance to erosion of the wall material (compacted repository wastes), erosion of the wall material will occur, increasing the diameter of the bored hole. The eroded material will then be passed to the surface in the flowing drilling fluid.

Flow in the annulus between the drillpipe and borehole wall is usually laminar (Darley and Gray, 1988, p243). Adjacent to the collars, however, the flow may be either laminar or turbulent as a consequence of the larger collar diameter and resulting higher mud velocities. For laminar flow, the analysis lends itself to classical solution methods. Turbulent flow, where the flow is assumed to be axial with a correction for the rotational component, requires a more approximate approach. A discussion of these two cases follows.

2.2.1 Laminar Flow

Below Reynolds numbers of about 2100 for Newtonian fluids and 2400 for some non-Newtonian fluids (Walker, 1976), experiments have shown that the flow of a fluid in a circular pipe or annulus is well behaved and can be described using a well-defined relationship between the velocity field and the fluid shear stress. This type of flow is called laminar.

Some of the early work on laminar helical flow of a non-Newtonian fluid in an annulus was performed by Coleman and Noll (1959), and Fredrickson (1960). The laminar helical flow solution procedure outlined below and used in the CUTTINGS_S code is, for the most part, an adaptation of methods described in a paper by Savins and Wallick (1966).

One of the principal difficulties in solving for the shear stresses within a helically flowing drilling fluid is the shear rate dependence of the fluid viscosity. This non-Newtonian fluid behavior necessitates choosing a functional form for the variation of viscosity with shear rate for the fluid. There are several functional forms for the viscosity of drilling fluids that can be assumed. For example, in the oil and gas industry the Bingham and power law models are often used to approximate the shear rate dependence of the fluid viscosity. A less common function is a form chosen by Oldroyd (1958) and used in the analysis by Savins and Wallick (1966). Oldroyd assumed that the viscosity varied according to the functional relation

$$\eta = \eta_o \left[\frac{1 + \sigma_2 \Gamma^2}{1 + \sigma_1 \Gamma^2} \right]$$
(2)

where σ_1 and σ_2 are constants, η_o is the limiting viscosity at zero rate of shear, η_{∞} (defined as $\eta_o(\sigma_2/\sigma_1)$) is the limiting viscosity at infinite rate of shear, and Γ is the shear rate. The viscous shear stress is described by $\tau = \eta \Gamma$.

Using the Oldroyd viscosity, Equation (2), the viscous shear stress can be illustrated graphically as in Figure 2a and 2b.

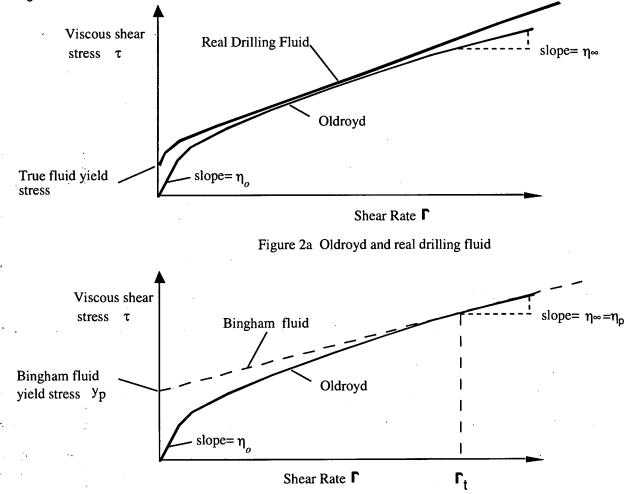


Figure 2b Oldroyd and Bingham fluid

This is a rate softening (pseudoplastic) model that has an initial slope of η_o and a limiting slope of η_∞ for large shear rates.

The Oldroyd model cannot account for drilling fluids that exhibit a yield stress. However, above a shear rate of zero, parameters can be chosen so that the model can be made to approximate the pseudoplastic rate response of many drilling fluids (see Figure 2).

Savins and Wallick (1966), expanding on the work of Coleman and Noll (1959) and Fredrickson (1960), showed that the solution for laminar helical flow of a non-Newtonian fluid in an annulus could be written in terms of three nonlinear integral equations.

$$F_{I} = \int_{\alpha}^{1} \left(\frac{\rho^{2} - \lambda^{2}}{\rho}\right) \frac{d\rho}{\eta} = 0$$

$$F_{2} = C \int_{\alpha}^{1} \frac{d\rho}{\rho^{3} \eta} - \Delta \Omega = 0$$

$$F_{3} = \frac{4Q}{\pi R^{3}} + 4 \left(\frac{RJ}{2}\right) \int_{\alpha}^{1} \left(\frac{\alpha^{2} - \rho^{2}}{\eta}\right) \left(\frac{\rho^{2} - \lambda^{2}}{\rho}\right) d\rho = 0$$
(3)

where α is the ratio of the collar radius over the cutting radius (R_i/R) (Figure 3), $\Delta\Omega$ is the drill string angular velocity, Q is the drilling fluid (mud) flow rate, r is the radial coordinate, and ρ is the non-dimensional radial coordinate representing the ratio r/R.

The unknown parameters λ^2 , RJ/2, and C are related to the fluid shear stresses through the relations

$$\tau_{r\theta} = \frac{C}{\rho^2}$$
$$\tau_{rz} = \frac{RJ}{2} \left(\frac{\rho^2 - \lambda^2}{\rho} \right)$$
$$\tau^2 = \tau_{r\theta}^2 + \tau_{rz}^2$$

(4)

where r, θ , and z represent radial, tangential, and vertical coordinates associated with the cylindrical geometry of Figure 3.

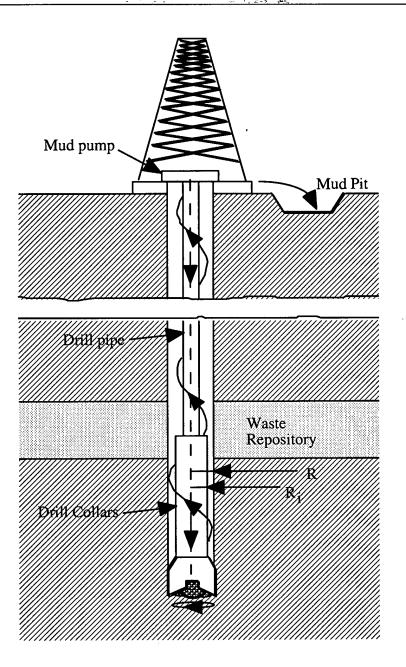


Figure 3. Detail of rotary drill string adjacent to repository.

The three nonlinear integral equations represented by (3) in general must be solved numerically. By expanding each of the integral equations into a Taylor series and retaining only the linear terms, a recursive solution procedure can be used (see Appendix AA)) to find the solution for the

unknowns $\delta \lambda^2$, $\delta \left(\frac{RJ}{2}\right)$, and δC . The three linear equations are

$$\frac{\delta F_{1}}{\partial \lambda^{2}} \delta \lambda^{2} + \frac{\delta F_{1}}{\partial C} \delta C + \frac{\delta F_{1}}{\partial \left(\frac{RJ}{2}\right)} \delta \left(\frac{RJ}{2}\right) = -F_{1}$$

$$\frac{\partial F_{2}}{\partial \lambda^{2}} \delta \lambda^{2} + \frac{\partial F_{2}}{\partial C} \delta C + \frac{\partial F_{2}}{\partial \left(\frac{RJ}{2}\right)} \delta \left(\frac{RJ}{2}\right) = -F_{2}$$

$$\frac{\partial F_{3}}{\partial \lambda^{2}} \delta \lambda^{2} + \frac{\partial F_{3}}{\partial C} \delta C + \frac{\partial F_{3}}{\partial \left(\frac{RJ}{2}\right)} \delta \left(\frac{RJ}{2}\right) = -F_{3}$$
(5)

The solution procedure consists of assuming initial values for λ^2 , $\left(\frac{RJ}{2}\right)$, and C and solving the three linear equations in (5) for the corrections $\delta\lambda^2$, $\delta\left(\frac{RJ}{2}\right)$, and δC . The unknowns λ^2 , $\left(\frac{RJ}{2}\right)$, and C are then replaced by $\lambda^2 + \delta\lambda^2$, $\left(\frac{RJ}{2}\right) + \delta\left(\frac{RJ}{2}\right)$, and $C + \delta C$. This recursive solution procedure is repeated until $\left|\delta\lambda^2\right|$, $\left|\delta\left(\frac{RJ}{2}\right)\right|$, and $\left|\delta C\right|$ are all less than some specified limit.

The coefficients of the unknowns $\delta \lambda^2$, $\delta \left(\frac{RJ}{2}\right)$, and δC in equations (5) are determined by differentiating equations (3). These derivatives are:

(8)

$$\frac{\partial F_{i}}{\partial \lambda^{2}} = -\frac{i}{a} \frac{1}{\eta \rho} \left[1 + \frac{(\rho^{2} - \lambda^{2})}{\eta} \frac{\partial \eta}{\partial \lambda^{2}} \right]^{d} \rho$$

$$\frac{\partial F_{i}}{\partial C} = -\frac{i}{a} \frac{1}{\eta \rho} \frac{(\rho^{2} - \lambda^{2})}{\eta} \frac{\partial \eta}{\partial C} d\rho$$

$$\frac{\partial F_{i}}{\partial \left(\frac{RI}{2}\right)} = -\frac{i}{a} \frac{1}{\eta \rho} \frac{(\rho^{2} - \lambda^{2})}{\eta} \frac{\partial \eta}{\partial \left(\frac{RI}{2}\right)} d\rho$$

$$\frac{\partial F_{2}}{\partial \lambda^{2}} = -C \int_{a}^{i} \frac{1}{\eta^{2} \rho^{3}} \frac{\partial \eta}{\partial \lambda^{2}} d\rho$$

$$\frac{\partial F_{2}}{\partial C} = \int_{a}^{i} \frac{1}{\eta^{2} \rho^{3}} \left[1 - \frac{C}{\eta} \frac{\partial \eta}{\partial C} \right] d\rho$$

$$\frac{\partial F_{2}}{\partial C} = -C \int_{a}^{i} \frac{1}{\eta^{2} \rho^{3}} \frac{\partial \eta}{\partial \lambda^{2}} d\rho$$

$$\frac{\partial F_{3}}{\partial C} = -4 \left(\frac{RI}{2}\right) \int_{a}^{i} \frac{\alpha^{2} - \rho^{2}}{\eta \rho} \left[1 + \frac{(\rho^{2} - \lambda^{2})}{\eta} \frac{\partial \eta}{\partial C} d\rho$$

$$\frac{\partial F_{3}}{\partial C} = -4 \left(\frac{RI}{2}\right) \int_{a}^{i} \frac{(\alpha^{2} - \rho^{2})}{\eta \rho} (\rho^{2} - \lambda^{2}) d\rho - 4 \left(\frac{RI}{2}\right) \int_{a}^{i} \frac{(\alpha^{2} - \rho^{2})}{\eta \rho} \frac{(\rho^{2} - \lambda^{2})}{\rho} \frac{\partial \eta}{\partial C} d\rho$$

From equations (4) and using the relation $\tau = \eta \Gamma$ the viscosity can be related to the shear rate function $Y(\Gamma)$ by the equation

$$\eta^{2}Y = 2\left[\left(\frac{RJ}{2}\right)^{2}\left(\frac{\rho^{2}-\lambda^{2}}{\rho}\right)^{2} + \frac{C^{2}}{\rho^{4}}\right]$$
(7)

where

For the Oldroyd viscosity function, equation (2), the unknown derivatives of the viscosity in equations (6) can be determined by using the chain rule of differentiation and equation (7).

 $Y = 2I^2$

$$\frac{\partial \eta}{\partial \lambda^{2}} = \frac{\partial (\eta^{2}Y)}{\partial \lambda^{2}} \frac{\partial \eta}{\partial (\eta^{2}Y)} = -4 \left(\frac{RJ}{2}\right)^{2} \left(\frac{\rho^{2} - \lambda^{2}}{\rho^{2}}\right) \frac{\partial \eta}{\partial (\eta^{2}Y)}$$

$$\frac{\partial \eta}{\partial C} = \frac{\partial (\eta^{2}Y)}{\partial C} \frac{\partial \eta}{\partial (\eta^{2}Y)} = \frac{4C}{\rho^{4}} \frac{\partial \eta}{\partial (\eta^{2}Y)}$$

$$\frac{\partial \eta}{\partial \left(\frac{RJ}{2}\right)} = \frac{\partial (\eta^{2}Y)}{\partial \left(\frac{RJ}{2}\right)} \frac{\partial \eta}{\partial (\eta^{2}Y)} = 4 \left(\frac{RJ}{2}\right) \left(\frac{\rho^{2} - \lambda^{2}}{\rho}\right)^{2} \frac{\partial \eta}{\partial (\eta^{2}Y)}$$
(9)

The derivative $\frac{\delta \eta}{\partial (\eta^2 Y)}$ can be determined by combining equations (2) and (8) and differentiating to

obtain

$$\frac{\partial \eta}{\partial (\eta^2 Y)} = \frac{\left(\frac{\sigma_2}{2} \eta_o - \frac{\sigma_1}{2} \eta\right)^2}{\left(\frac{\sigma_2}{2} \eta_o - \frac{\sigma_1}{2} \eta\right) \left[\eta^2 + 2(\eta - \eta_o)\eta\right] + (\eta - \eta_o)\eta^2 \frac{\sigma_1}{2}}$$
(10)

A cubic equation for η in terms of the unknown parameters λ^2 , RJ/2, and C can be written by combining equations (2), (7) and (8)

$$\eta^{3} - \eta_{o}\eta^{2} + \sigma_{I} \left[\left(\frac{RJ}{2} \right)^{2} \left(\frac{\rho^{2} - \lambda^{2}}{\rho} \right)^{2} + \frac{C^{2}}{\rho^{4}} \right] \eta - \sigma_{2} \left[\left(\frac{RJ}{2} \right)^{2} \left(\frac{\rho^{2} - \lambda^{2}}{\rho} \right)^{2} + \frac{C^{2}}{\rho^{4}} \right] \eta_{o} = 0$$
(11)

2.2.1.1 Solution Process for Laminar Flow

Laminar helical flow of an Oldroyd fluid in an annulus of a given size with a rotating inner core is based on the solution to the three integral equations (3) which have three unknown parameters λ^2 , RJ/2, and C. A recursive process for computing the unknowns based on a Taylor series expansion (equation 5) requires the evaluation of the coefficients of the three linear equations. These coefficients (equations 6 combined with 9, 10, and 11) involve integrals across the radius of the annulus which must be integrated numerically. Simpson's 1/3 rule (Salvadori, et al 1962) is used to perform these integrations by dividing the annular radius into ten equal intervals. The initial solution obtained from the recursive process assumes no erosion has occurred and thus corresponds to an annulus with an outer radius equal to one half the borehole diameter. For the specific case of borehole erosion, once a solution to the unknowns (λ^2 , RJ/2, and C) in the three integral equations in (3) is found, the shear stress in the fluid τ at the wall can be calculated by setting $\rho = 1$ in the equations in (4). By changing the outer radius of the hole and repeating the above process a new shear stress in the fluid at the wall can be calculated. By adjusting the outer radius appropriately the fluid shear stress can be forced to equal the repository effective shear resistance to erosion. That radius corresponds to the final eroded radius. The required outer hole radius is

determined by iteration as shown in Figure 4. The derivatives required for the iteration $\left(\frac{d\tau}{dR}\right)$ are found

numerically by determining the change in shear stress for a small assumed change in outer radius R.

Based upon the preceding equations a Fortran computer code was written to perform the necessary computations for a solution to the problem of laminar helical flow in an annulus. This code was partially verified by comparing its results against those published by Savins and Wallick (1966).

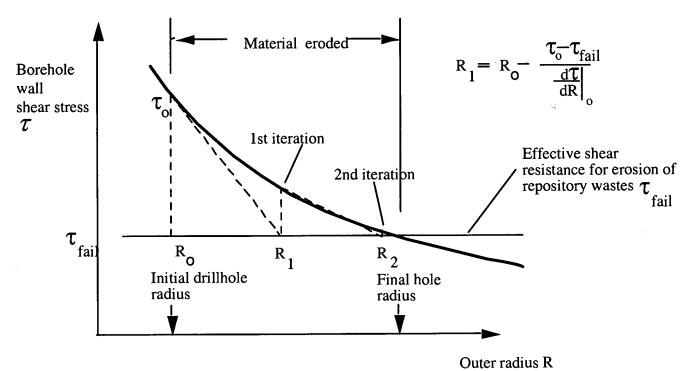


Figure 4. Iterate to find the final hole radius

The final eroded diameter is determined through an iterative process that equates the fluid shear stress adjacent to the waste to a measure of the erosion resistance of the waste. The erosion resistance is governed by the effective shear resistance to erosion, a parameter not directly related to the shear strength as normally determined for granular materials using conventional soil tests.

The effective shear resistance for erosion (τ_{fail}), equals the threshold value of fluid shear stress required to sustain general erosion at the borehole wall. Parthenaides and Passwell (1970), in discussing investigations on the erosion of seabed sediments and in channels, has noted that this effective soil shear resistance is not related to the soil shear strength as normally determined from conventional soil tests. The effective shear resistance for erosion based on seabed data, as determined by Parthenaides and Passwell (1970), is on the order of a few Pa and is thus smaller by several orders of magnitude than the macroscopic soil shear strength.

2.2.2 Turbulent Flow

For Newtonian fluids with Reynolds numbers greater than about 2100, flow in a circular pipe or annulus starts to become more or less random in character, which makes orderly mathematical analysis of the flow

difficult, if not impossible. With increasing Reynolds numbers, this random behavior increases until, at a Reynolds number of about 3000, the flow becomes fully turbulent. In fully turbulent flow, momentum effects dominate and the fluid viscosity is no longer important in characterizing pressure losses.

The Reynolds number (R_e) is defined as

$$R_e = \frac{\overline{\rho} \overline{V} D_e}{\overline{\eta}} \tag{12}$$

where D_e is the equivalent hydraulic diameter, $\overline{\rho}$ is the drilling fluid density, \overline{V} is the average axial fluid velocity, and $\overline{\eta}$ is the average fluid viscosity.

For Newtonian fluids, the value to use for the viscosity is clear since the viscosity is constant for all rates of shear. Non-Newtonian fluids exhibit a changing viscosity with shear rate and present a special problem in calculating R_e . For fluids that exhibit a limiting viscosity at high rates of shear (such as the Bingham model and in our case the Oldroyd model), it has been suggested (Broc, 1982) that the limiting viscosity $(\overline{\eta} = \eta_{\infty})$ be used in calculating the Reynolds number.

The Reynolds number for an Oldroyd fluid in an annulus can then be written as (Broc, 1982)

$$R_e = \frac{0.8165 D_e V \overline{\rho}}{\overline{n}} \tag{13}$$

where the equivalent hydraulic diameter is expressed as $D_e = 2(R - R_i)$ (see Figure 3).

The most important influence viscosity has on the calculation of pressure losses in fully turbulent flow of non-Newtonian fluids appears to be in the calculation of the Reynolds number. A far more important parameter is the surface roughness past which the fluid must flow. The Reynolds number, however, does have a role in determining the onset of turbulence. For Newtonian fluids this number is about 2100. For non-Newtonian, rate-thinning fluids, the critical value of R_e tends to be greater than 2100 but less than 2400 (Walker, 1976). For our purposes, a value of 2100 will be used to represent R_{ec} (the critical Reynolds number) for the Oldroyd fluid model. Since turbulent flow is more effective in generating fluid shear stresses at the borehole wall, this assumption is conservative.

There is a transition region beyond R_{e_c} before the development of fully turbulent flow. In this regime the flow has the character of both laminar and turbulent flow. However, since pressure losses increase rapidly in turbulent flow and affect borehole shear stresses more severely, it will be assumed that beyond R_{e_c} the flow is fully turbulent.

Turbulent flow is very complex and, thus, to characterize the turbulent flow regime, the great bulk of analysis has concentrated on empirical procedures. For axial flow in an annulus, the pressure loss under turbulent conditions can be approximated by (Broc, 1982)

$$\Delta P = \frac{2 f L \overline{\rho} V^2}{(0.8165) D_e} \tag{14}$$

where f is the coefficient of pressure head loss (Fanning friction factor) and L is the borehole length.

If the shear stress due to the flowing fluid is assumed to be uniformly distributed on the inner and outer surfaces of the annulus, it can be easily shown using Equation (14) that the shear stress is related to the average fluid velocity through the relation

$$\tau = \frac{f\bar{\rho}\,\overline{V}^2}{2(0.8165)} \tag{15}$$

The Fanning friction factor is empirically related to the Reynolds number and relative roughness by the equation (Whittaker, 1985)

$$\frac{1}{\sqrt{f}} = -4\log_{10}\left[\frac{\varepsilon}{3.72D} + \frac{1.255}{R_e\sqrt{f}}\right]$$
(16)

where $\varepsilon I D$ is the relative roughness. For circular pipes, D in this equation represents the inside diameter and ε is the absolute roughness or the average depth of pipe wall irregularities. In the absence of a similar equation for flow in an annulus, it is assumed that this equation also applies here, where D is the equivalent hydraulic diameter (D_e) as defined earlier and ε is the absolute roughness of the wasteborehole interface.

For laminar flow within the annulus both axial and rotational flow are modeled. Some of the available literature (Khader and Rao, 1974; Bilgen E., Boulos, R., and Akgungor A.C., 1973) indicates the importance of also accounting for drillstring rotation when the drilling mud flow within the annulus is turbulent. Consequently, to account for rotational flow in the turbulent regime, an axial flow velocity correction factor (rotation factor) is introduced into the above formulation that maintains eroded diameter compatibility across the laminar- turbulent flow transition. The rotation factor (F_r) is determined by

increasing (or decreasing) the axial velocity \overline{V} in equation 15 until the turbulent flow eroded diameter equals the laminar flow eroded diameter computed at the prescribed angular velocity at a Reynolds number of 2100. The rotation factor is defined as

$$F_r = \frac{\overline{V}_{2100}}{\overline{V}}$$

where \overline{V}_{2100} is the axial velocity required for eroded diameters to be the same for turbulent and laminar flow. This rotation factor is then used to modify the axial turbulent flow velocity at all other turbulent flow Reynolds numbers. The rotation factor is not used when computing the Reynolds number. Equation 15 then can be rewritten

$$\tau = \frac{f\overline{\rho}(F_r\overline{V})^2}{2(0.8165)}$$

(17)

2.2.2.1 Solution process for turbulent flow

Using a relative roughness and a calculated Reynolds number based on \overline{V} , a Fanning friction factor can be determined by iteratively solving Equation (16). The value of the shear stress acting on the borehole wall can then be determined from Equation (17). Using an iterative procedure similar to that for the laminar flow problem, the fluid shear stress can be forced to equal the repository shear resistance to erosion (τ_{fail}) to obtain the final eroded borehole radius (see Figure 4).

The equations governing cavings (and cuttings) based on laminar and turbulent flow form a portion of the Fortran computer code CUTTINGS_S (see figure 1). Using appropriately selected input based on the physical properties of the waste and other drilling parameters, this section of the code calculates the final eroded diameter of the borehole that passes through the waste. In the actual solution sequence employed in the fortran code the Reynolds number is calculated first to determine which solution regime (laminar or turbulent) should first be initiated. For Reynolds numbers initially less than R_{e_c} the code calculates the flow as laminar. Any increase in diameter of the borehole calculated during the laminar calculation will

cause the Reynolds number to decrease ensuring that the calculation remains laminar. If the initial Reynolds number is greater than or equal to R_{e_c} the turbulent formulation is used to calculate borehole erosion. When the turbulent calculation is complete, a check is again made to determine if the Reynolds number still exceeds R_{e_c} . If it does not, the laminar calculation is performed starting with a "critical" borehole radius. The critical borehole radius corresponds to a Reynolds number of R_{e_c} and is given by

$$R_{crit} = \frac{\overline{\rho}Q}{1286\,\pi\eta_{\rm m}} - R_i \tag{18}$$

2.3 Drilling parameters for cuttings and cavings

The drilling parameters chosen must reflect data typical of that valid near the WIPP repository. For drilling operations through salt in the Delaware basin , the drilling mud most likely to be used is a brine (Pace's letter on pages A-159 through A-164 of (Rechard et al., 1990)), with the density cut somewhat with an emulsified oil. The mud density and viscosity parameters required in the laminar and turbulent regime erosion calculations can be estimated based on the assumption of the use of such a brine-based drilling mud. For the 1992 PA the drilling mud properties used were (Sandia WIPP Project 1992, Table 4.4-1):

Parameter	Median	Range			Units		Distribution Type
Density 1211	1139	-	1378		kg/m ³		Constructed
Bingham Plasti Viscosity η_p	c 0.00917	0.005	-	0.03		Pa∙s	Constructed
Bingham Yield Stress y _p	4	2.4	-	19.2		Ра	Constructed

The values for plastic viscosity and yield stress correspond to parameters for a Bingham plastic fluid (Darley and Grey 1988, p188) from which the Oldroyd parameters are derived.

If in equation (2) it is assumed that $\eta_{\infty} = \eta_o \left(\frac{\sigma_2}{\sigma_1}\right) = \eta_p$ where η_p is the Bingham plastic viscosity, and

if $\eta_o = 2\eta_\infty$ (a trend based on other drilling muds (Darley and Grey 1988, p242)) the following relationship can be derived.

$$\sigma_2 = \frac{\eta_p \Gamma_i - y_p}{2\Gamma_i^2 y_p} \tag{19}$$

where I_r is the shear rate at which the Oldroyd model and Bingham model coincide (Figure 2b). This shear rate is chosen to be 1020 1/s which corresponds to a reading at 600 rpm on a direct reading viscometer (Darley and Grey 1988, p214).

For drilling through salt, the drilling speeds can vary from 40 to 220 rpm (Austin, 1983 and Pace's letter on pages A-159 through A-164 of (Rechard et al., 1990)). The most probable speed is about 70 rpm (Pace's letter on pages A-159 through A-164 of (Rechard et al., 1990)).

Mud flow rates are usually selected to be from 30 to 50 gallons/minute per inch of drill bit diameter (Austin, 1983) and usually result in flow velocities in the annulus between the drill collars and the borehole wall at or near the critical flow state (laminar-turbulent transition) (Pace's letter on pages A-159 through A-164 of (Rechard et al., 1990)).

The drill bit diameter is related to the total planned depth of the hole to be drilled. For gas wells in the 4000- to 10,000-foot range, it is likely that the drill bit used that passes through a waste room would have a diameter of 10.5 to 17.5 inches. The most probable drill diameter is 12.25 inches. The collar diameter is assumed to be 8 inches, a value consistent with current drilling practice in the Delaware Basin (Marquis, 1995). For the 1992 PA these parameters are presented in table 4.2-1 (Sandia WIPP Project 1992)

Parameter	Median	Range			Units	<u> </u>	Distribu	<u>ition Type</u>
Drill bit 0.355 diameter	0.267	-	0.444		m		Uniform	n
Drill string angular velocity	7.7	4.2	-	23		rad/s		Constructed
Drilling mud flowrate	0.09935	0.0745	-	0.124		m ³ /(s∙n	n)	Uniform

For turbulent flow, the shear stress acting at the borehole at the repository is dependent on the absolute surface roughness. The value chosen for the calculations exceeds that of very rough concrete or riveted steel piping (Streeter, 1958).

The amount of material eroded from the borehole wall is dependent upon the magnitude of the fluidgenerated shear stress acting on the wall and the effective shear resistance to erosion of the compacted, decomposed waste. In the absence of experimental data, the effective shear resistance to erosion of the repository material is assumed to be similar to that of a ocean bay mud (Parthenaides and Passwell, 1970), or a montmorillonite clay, (Sargunam et al., 1973). These values are on the order of a fraction to several Pa.

For the 1992 PA, the roughness and shear resistance parameters were given in table 3.4-1 (Sandia WIPP Project 1992).

Parameter	Median	Range)		Units	Distribution Type
Absolute roughness	0.025	0.01	-	0.04	m	Uniform
Erosion shear resistan	1 ce	0.05	-	10	Pa	Constructed

The lower bound to erosion resistance shown above is lower than that used in the 1992 PA and is based on the availability of additional data (Parthenaides and Passwell, 1970).

The hydrostatic pressure at the repository depth resulting from 1211 kg/m³ drilling mud corresponds to 7.78 MPa. In the CUTTINGS_S code repository pressures below 8 MPa are assumed to cause surface releases from cuttings and cavings. Above 8 MPa releases are caused by spall (see figure 1).

2.4 Spallings

Spallings comprises that quantity of solid waste that reaches the ground surface as the result of the high pressure gas in the waste. If the repository gas pressure exceeds the hydrostatic pressure of the column of drilling fluid, solids releases can occur from blowout, stuck pipe or gas erosion. Contaminated brine can also be expelled during a blowout event, a process not considered in the CUTTINGS_S code. Brine releases to the surface resulting from blowout are computed using a separate code.

2.4.1 Blowout (Solids Removal)

Blowout is assumed to occur when repository gas pressures exceed 8 MPa (~hydrostatic) and when waste permeabilities exceed 10^{-16} m² (see figure 1). For permeabilities that are lower than 10^{-16} m² the gas flow into the borehole is assumed to be too low to cause well blowout (Berglund 1994).

Two alternative computational models are used to compute the solid waste released to the surface resulting from the sudden penetration of a pressurized WIPP panel. The models assume waste strength based on brine saturation and are self limiting ie. neither model relies on driller response. Of the two models described below, model 2 is more consistent with data obtained from experiment.

2.4.1.2 Solids Removal Caused by Blowout (Model 1)

A borehole penetration of a WIPP panel saturated by decomposition gases at pressures above hydrostatic and with a waste permeability greater than 10⁻¹⁶ m² is likely to initiate a blowout of the well. The initial flow of gas into the borehole will be restricted somewhat by the presence of drilling mud and a backpressure may develop that could be greater than hydrostatic. As the gas flows into the borehole and rises due to buoyancy, the backpressure will diminish as the weight of the mud column decreases. With a decreasing backpressure the gas flowrate into the borehole will increase. It will also increase as the result of the removal of waste material and the generation of a larger surface through which the gas can flow. Eventually the borehole will be cleared of drilling mud and the backpressure will drop to a value governed by the rate of flow of a compressible fluid in the drillstring annulus.

In the development of a bounding model for waste release due to a blowout it is important to note that the blowout process occurs over a finite time (minutes) and consequently it is not necessary to consider the high transient gas flow rates that occur from the assumption of instantaneous pressures drops. This is especially true for waste with high permeability where pressure profiles in the waste change rapidly. After these initial transients the gas flowrates out of an exposed waste surface become relatively constant for volumes comparable to a WIPP panel. It is this relative constancy of flow that allows for the development of the bounding model described below.

During blowout and after the decay of initial flow transients the gas velocity (V) at an exposed waste surface remains relatively constant for long periods of time. If the gas velocity is great enough and there is little cohesive strength between waste grains, the waste grains at this free surface will be lofted into the flow of gas (entrained) and be carried into the borehole and to the ground surface. As the waste is removed and the exposed waste surface area increases, the resistance to flow up the borehole annulus increases causing a backpressure to develop in the evacuated volume. Consequently, the velocity of gas at this interface will decrease, eventually reaching a condition where waste is no longer entrained into the flowing gas at the waste surface. At this point an equilibrium condition will have been achieved and the volume associated with the dimensions of the exposed surface will be equal to a bounding waste release due to blowout.

The volume released is dependent on the shape of the void produced near the borehole. The shape is dependent on the waste properties and the character of the release process. A possible shape may be similar to the relatively flat ellipsoidal form shown in Figure 5.

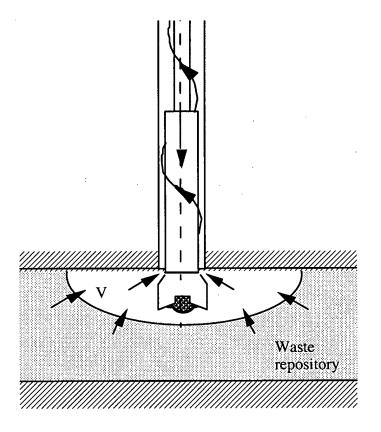


Figure 5 Potential Final Release Volume Configuration

A shape that provides a larger volume for the given exposed surface and one that is easy to calculate is the cylindrical volume shown in Figure 6.

The bounding blowout radius R_b (Figure 6) can then be written

$$R_{b} = \left[\left(\frac{D_{o}}{2} \right)^{2} - \left(\frac{D_{i}}{2} \right)^{2} \right] \frac{M_{I}c}{2HV_{e}}$$

where

 D_o and D_i are the borehole and collar diameters

- M_1 = the mach number for gas flow into the borehole entrance
- c = the gas sound speed at the borehole entrance
- H = the panel height at time of intrusion
- V_e = the entrainment velocity

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(20)

In this equilibrium condition the gas velocity V is equal to the entrainment velocity V_e and the solid waste volume released to the surface is

$$Volume = \pi R_b^2 H(l - \varphi) \tag{21}$$

where φ is the waste porosity at the time of intrusion. For the "actual" void shape (Figure 5), V_e is the gas velocity required to loft a waste particle vertically in a gravity field. This velocity is also used in the computational model where the void shape is assumed to be cylindrical.

As with cuttings (section 2.1), since the waste is assumed to be uniformly distributed throughout the disposal region, the actual computation for release requires only the area associated with R_b and the waste curie content per unit area. The panel height at the time of intrusion is related to the original height and waste porosity, and to the porosity at the time of intrusion.

$$H = \text{Initial height}\left(\frac{l - \text{initial porosity}}{l - \varphi}\right)$$

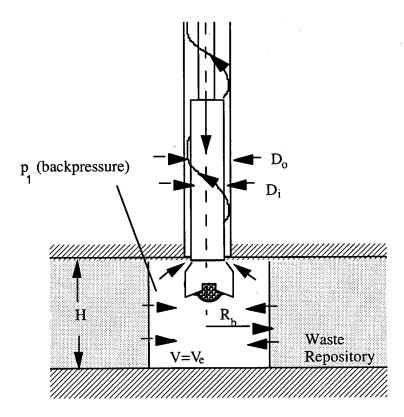


Figure 6 Idealized Cylindrical Final Volume Release Configuration

The borehole entrance mach number is computed based on flow in the borehole annulus and assumes that the borehole is free of drilling mud and undergoing full blowout. Compressible, isothermal flow in a channel (the borehole annulus) with friction is governed by the equation (Binder, 1958)

$$\frac{fl}{D} = \frac{1}{KM_1^2} \left[1 - \frac{p_2^2}{p_1^2} \right] - 2\ln\frac{p_1}{p_2}$$
(22)

where f=friction factor, D=pipe diameter, l = pipe length, K=ratio of specific heats of the gas, p₁ and p₂= inlet and outlet gas pressures, and M₁= inlet Mach number. The pipe diameter to use for our purposes is the effective hydraulic diameter computed from the annular area. Actually there are two annular areas that are considered for flow up the borehole; one area is adjacent to the drill collars and the other is adjacent to the drill pipe and consequently equation (22) has to be solved over these two regions. The pressure ratio is based on the exit pressure at the ground surface ($p_2 \ge$ atmospheric) and the borehole entrance pressure p₁ (backpressure).

The entrainment velocity V_e is the magnitude of the gas velocity at the exposed waste surface that will cause waste particles to be removed from the surface and transported in the gas to the borehole. For a horizontal waste surface this velocity is assumed to be the terminal velocity of the particle. The terminal velocity is the speed at which a particle will fall in a motionless gas of density ρ_g . For the present model, choosing this value to represent V_e is conservative since the terminal velocity is less than the surface normal gas velocity required to vertically support the particle if the surface is inclined.

The terminal velocity of a spherical particle can be found by equating the particle weight to the drag force acting on the particle when traveling at a velocity V_e . The relationship that results is (Cheremisinoff, 1984)

$$V_e^2 = \frac{4gd(\rho_s - \rho_g)}{3C_D\rho_g}$$
(23)

where g= acceleration of gravity, d= particle diameter, ρ_s = particle density, ρ_g = gas density, and C_D is the coefficient of drag. The coefficient of drag for a sphere is an empirical function of the particle reynolds number R_{ed} (Fox et. al. 1973)

where
$$R_{ed} = \frac{\rho_{g} V_{e} d}{\mu}$$
 and μ is the gas viscosity

For $R_{ed} < 0.4$ $C_D = \frac{24}{R_{ed}}$

For $0.4 < R_{ed} < 200000$ the data can be fit with the following function

$$C_{D} = 10^{(a+\bar{b}x+\bar{c}x^{2}+\bar{d}x^{3}+\bar{e}x^{4}+\bar{f}x^{5}+\bar{g}x^{6})}$$
(24)

where

(27)

 $\bar{a} = 1.3918$ $\bar{b} = -0.907723$ $\bar{c} = 0.136371$ $\bar{d} = 0.0165093$ $\bar{e} = -0.0285484$ $\bar{f} = 0.00933281$ $\bar{g} = -0.000897166$

and

 $x = \log_{10} R_{ed}$

For $R_{ed} > 200000$ $C_D = 0.2$

Equations (23) and (24) are solved simultaneously using an iterative process that converges to a value of V_a that satisfies both equations.

The velocity of gas flow (V) out of the cylindrical waste surface can be computed for various waste permeabilities and repository gas pressures using the equation

$$\frac{k}{2\mu}\nabla^2 p^2 = \varphi \,\frac{\delta p}{\partial t} \tag{25}$$

where k is the permeability, μ is the gas viscosity, ϕ is the waste porosity, $\nabla^2 = \frac{\dot{o}^2}{\dot{o}r^2} + \frac{l}{r}\frac{\dot{o}}{\dot{o}r}$, p is the

gas pressure, and r is the radial coordinate. The rate is dependent on the pressure boundary condition (p_1) (Figure 6) assumed in the evacuated volume. The gas velocity in the waste at any radius r and time t is a function of the pressure gradient

 $V=v(R_b,t)$

$$v(r,t) = \frac{k}{\mu} \frac{\delta p(r,t)}{\partial r}$$
(26)

and the gas velocity at the free surface is

Solving equations (20), and (22-27) simultaneously and using stabilized flowrates out of the waste the bounding volumes of waste removed due to blowout can be obtained.

As observed earlier, the blowout process occurs over a finite time and the gas flowrate out of the free surface (r=R_b) tends to stabilize and remain constant for long periods of time. The time chosen to measure this "stable" flow condition has been chosen to be 700s after penetration. This is the minimum time necessary to replace the drilling mud in the annulus volume at the maximum expected drilling mud flowrate (50 gal/minute/inch of borehole diameter). This time is a measure of the minimum time necessary to generate full blowout.

The nonlinear partial differential equation for pressure in the waste (equation 25) can only be solved using numerical approximations (this is done in a code called GASFLOW) but the resulting solution is time consuming since it must be solved many times using different boundary conditions (p1 and Rb). This process can be simplified by analytically solving the steady state form of equation 25 and by using a set of correction factors (calculated in GASFLOW) to obtain the free surface gas velocity at 700s. The steady

> $\nabla^2 p^2 = 0$ (28)

The general solution to equation 28 is

where

In these equations R_a is assumed to be the radius of an equivalent WIPP panel (~61m) and p_i is the far field gas pressure (ie. the initial pressure in the panel). The steady state gas velocity at the free surface is

$$V(R_b) = \frac{k}{\mu} \frac{dp(R_b)}{dr} = \frac{k}{\mu} \frac{C_2}{2R_b (C_1 + C_2 \ln R_b)^{0.5}}$$
(30)

 $p^2 = C_1 + C_2 \ln(r)$

 $C_2 = \frac{p_1^2 - p_{\infty}^2}{ln\frac{R_b}{R_{\infty}}}$

 $C_1 = p_1^2 - C_2 \ln(R_h)$

Equation 30 generally gives velocities lower than those that would be calculated from equation 26 using the GASFLOW code and setting the time to 700s. To be conservative in computing waste releases, the results from equation 30 should be at least as great as the velocities computed by GASFLOW at 700s. This is accomplished by multiplying the velocity calculated by equation 30 by a factor obtained from specific GASFLOW computations. For example, assuming the repository is pressurized with hydrogen gas (μ =9.3x10⁻⁶ Pa·s) the following velocity ratios

$$V_{R}$$
 = velocity ratio = $\frac{Velocity from GASFLOW @ 700s}{Steadystate velocity from equation 30}$

are obtained for different radii Rb and waste permeabilities k.

let $p = 8x10^6$ Pa, $p_1 = 1x10^6$ Pa, $\phi_0 = 0.19$, R = 61m, then

k (m²) velocity velocity velocity velocity velocity ratio V_Rratio V_Rratio V_Rratio V_R

 $R_{b}=0.1 m$ $R_{h}=0.3m$ 1x10⁻¹²0.558 0.897 0.92 0.734 0.836 1x10⁻¹³1.46 1.29 1.23 1.18 1.38 1x10⁻¹⁴2.75 2.19 1.78 1.53 1.12 1x10⁻¹⁵6.98 1.87 4.64 3.08 2.25

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state form of equation 25 is

(29)

(31)

 $R_b=8mR_b=3mR_b=1m$

1x10 ⁻¹⁶ 20.5	12.23	6.68	3.90	2.79

equation 30 thus becomes

$$V(R_b) = \frac{k}{\mu} \frac{V_R C_2}{2R_b (C_1 + C_2 \ln R_b)^{0.5}}$$
(32)

A more complete set of tables for the velocity ratio data is shown in appendix AB. Data from these tables can be used to generate an equation for V_R as a function of the waste permeability. The fit corresponds to data values which produce the largest velocity ratio for a particular permeability. This process is conservative. See appendix AB. The resulting relationship for the velocity ratio is

$$V_{R} = -12593 - 5040\xi - 806.18\xi^{2} - 64.455\xi^{3} - 2.5771\xi^{4} - 0.04125\xi^{5}$$
(33)

where

$$\xi = \log_{10} k$$

The terminal velocity described above (equations 23 and 24) is based on no cohesive strength between waste particles. Strength can be included in the model by changing the gravity term to that of an effective acceleration of gravity.

2.4.1.2.1 Effect of Strength on Solids Removal caused by Blowout

The model for solids removal during blowout is based on the lofting of particles from a horizontal surface of the waste as the result of the flow of escaping gas from the interior of the waste. Waste is lofted into the flowing gas if the velocity of the gas is equal to the terminal velocity of the waste particle. The equation for terminal velocity (equation 23) can again be written as

$$V_e^2 = \frac{4gd(\rho_s - \rho_g)}{3C_D \rho_g}$$
(34)

where g= acceleration of gravity, d= particle diameter, ρ_s = particle density, ρ_g = gas density, and C_D is the coefficient of drag.

The downward force resisting the drag force caused by the flowing gas is the weight of the particle. If there are additional forces acting to resist lofting such as tensile strength an effective gravity force can be computed and used to replace the gravity force in the terminal velocity equation.

Consider a material that has a tensile strength σ . This parting stress on the particle level can be written

$$\sigma = \frac{Parting \text{ force}}{Particle \text{ projected area}} = \frac{mg_t}{\pi R^2} = \frac{\frac{4}{3}\pi R^3 \rho_s g_t}{\pi R^2} = \frac{4}{3}\rho_s g_t R$$
(35)

where m is the particle mass, R is the particle radius, ρ_s is the particle density and g_i is the effective acceleration of gravity necessary to generate a particle weight equal to the parting force.

Solving equation (35) for g, obtains

$$g_t = \frac{3}{4} \frac{\sigma}{\rho_s R} \tag{3}$$

The effective acceleration of gravity that accounts for particle weight and strength can then be written

$$g_{eff} = (g + g_t) = g + \frac{3}{4} \frac{\sigma}{\rho_s R}$$
(37)

This effective acceleration of gravity can be used in the equation that computes the particle terminal velocity (equation 23) thus accounting for strength. A model for computing the parting stress based on waste saturation and particle size is presented in appendix AD. Additional contributions to strength, such as from cementation between particles, can also be included in the parting stress σ .

2.4.1.2.2 Solution Process for Solids Removal caused by Blowout (Model 1)

Utilizing Boyles Law which relates gas pressure to gas density it is possible to reduce equations 20, 22, 23, 24, and 32 to two simultaneous nonlinear equations in the unknowns R_b and the backpressure p_1 . A solution process similar to that described in appendix AA was used to solve this system but it often failed to converge to the solution and was therefore considered unsuitable for use in a code where the input parameters were to vary over considerable ranges. To create a more robust solution to the system of equations, they are solved sequentially starting with equation 22.

As indicated earlier, flow in the borehole annulus is assumed to occur adjacent to both the drill collars and the drillpipe. Consequently, there are two annular areas that must be considered for flow up the borehole and equation (22) has to be solved over these two regions. The pressure ratio is based on the inlet and exit pressures for the two annuli with pressure compatibility maintained at the drillpipe / drill collar connection.

Utilizing conservation of mass, equation 22 can be divided into the two equations

$$\frac{fl_{coll}}{D_{coll}} = \frac{1}{KM_{I}^{2}} \left[1 - \frac{p_{i}^{2}}{p_{I}^{2}} \right] - 2\ln\frac{p_{I}}{p_{i}}$$
(38)

$$\frac{fl_{drill}}{D_{drill}} = \frac{1}{KM_{j}^{2} \left(\frac{p_{j}}{p_{i}}\right)^{2} \left(\frac{D_{coll}}{D_{drill}}\right)^{4}} \left[1 - \frac{p_{exit}^{2}}{p_{i}^{2}}\right] - 2\ln\frac{p_{i}}{P_{exit}}$$
(39)

where p_{exit}= exit pressure

l_{drill}= length of drillpipe l_{coll}=collar length

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(36)

D_{coll}=equivalent diameter of annulus opposite collars (based on annulus area) D_{drill}=equivalent diameter of annulus opposite drillpipe (based on annulus area) p_i=intermediate pressure

Eliminating M_1^2 from equations 38 and 39 results in a single nonlinear equation in the dependent variables p_i and p_{exit} .

$$F(p_i, p_{exit})=0$$
(40)

A sequence of solutions to equation 40 for various backpressures p_1 are obtained by systematically scanning over possible values of p_i and p_{exit} and using the additional constraint (Binder, 1958 p 64) that dp

either $\frac{dp}{dl}$ becomes unbounded at the pipe exit or $p_{exit} = p_{atm}$. Using this method results in a discrete set of

 p_1 , M_1 pairs that satisfy equations 38 and 39.

Three equations remain to be satisfied; equations 20, 23 and 32.

For each p_1 , the gas density can be determined from Boyles law and the terminal velocity V_e computed from equations 23 and 24 again using an iterative process. For the corresponding mach number in the p_1 , M_1 pair it is now possible to compute R_b utilizing equation 20. The velocity V at the waste surface is then obtained from equation 32 using p_1 and R_b . For a valid solution the waste surface velocity V must equal the terminal velocity V_e . New p_1 , M_1 pairs are selected sequentially until the V, V_e solutions approach equality and then linear interpolation is used for the final adjustment.

2.4.1.3 Solids Removal Caused by Blowout (Model 2)

Results from steady state flow experiments through granular material in a cylindrical geometry (Lary R. Lenke, et al. 1996) indicate that a porous pattern of channels is formed adjacent to the "borehole". This channeling process does not conform to conceptual model 1 as described above. As a consequence of these experiments a new conceptual model has been defined (model 2) which more closely corresponds to the physical processes that appear to be occurring. The new conceptual model is described below.

Conceptual model 2 is based on a number of assumptions. These are:

1. After intrusion by a drillbit, the pressure gradients associated with the flow of gas toward the borehole fracture the porous (waste) material sufficiently to permit the escaping gas to flow within the fractures rather than through the porous matrix (Figure 7). Consequently, the intrinsic permeability of the matrix no longer restricts gas flow and the gas pressure at the borehole entrance can be assumed to be the initial pore pressure.

2. The gas flow velocity up the borehole is governed by the isothermal flow of gas in a long tube of a given cross-sectional area, tube roughness, and gas pressure at the borehole entrance. (For sufficiently high borehole entrance pressures the flow velocity tends to choke or become self limiting.)

3. The mass flowrate of gas in the fractures at any radial cross-section away from the borehole is equal to the mass flowrate up the borehole.

4. Radial gas flow within the fractures of the waste matrix erode and widen the fractures. Erosion is assumed to occur if the fracture gas velocity exceeds a velocity v_e . (The higher gas velocities in fractures near the borehole create wider channels near the borehole.)

5. The fracture erosion velocity v_e is related to the terminal velocity of a waste particle at the fracture surface and to the cohesive strength afforded by moisture and cementation in the matrix. Experiments will determine the effectiveness of cohesive strength and gravity on erosion and will provide experimentally determined effectiveness factors F_{se} , F_{ge} , and F_{ce} which are related to these factors.

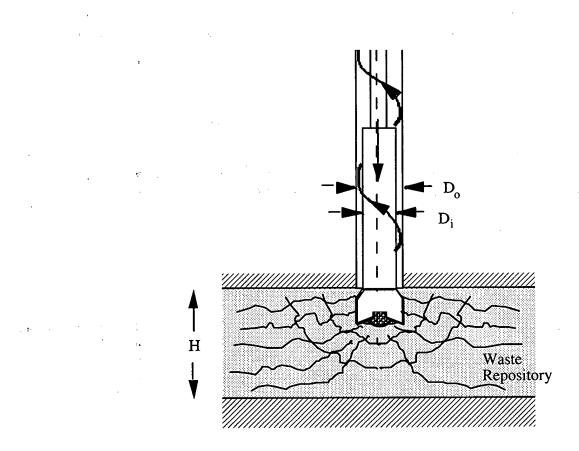


Figure 7 Fractured Waste Matrix after Penetration

2.4.1.3.1 Computation of Release Volume (Model 2)

Since the gas pressure is assumed to be constant throughout the fractured, porous matrix, conservation of mass flow requires (Figure 8)

(41)

$$2\pi r V H = A_{borehole} V_{borehole}$$

or

$$V = \frac{A_{borehole}V_{borehole}}{2\pi r H}$$

where V is the average gas velocity within the matrix at a radius r from the borehole and H is the height of a repository room at the time of intrusion. The borehole annulus area is denoted as A_{borehole} and the average gas velocity up the borehole annulus is denoted V_{borehole} .

The local gas velocity within the fractures (Figure 8) can be written as $v = \frac{V}{\varphi_f}$ where φ_f is the fracture porosity. Thus the fracture porosity associated the erosion fracture velocity v, is

$$\varphi_{fe} = \frac{V}{v_e} = \frac{A_{borehole} V_{borehole}}{2\pi r H v_e}$$
(42)

The solid volume eroded from the fractures is

$$V_{ol} = 2\pi H (1-\phi) \int_{0}^{r_{o}} \varphi_{fe} r dr = \frac{A_{borehole} V_{borehole}}{v_{e}} (1-\phi) \int_{0}^{r_{o}} dr$$

$$\tag{43}$$

or

$$V_{ol} = \frac{A_{borehole} V_{borehole} r_o}{v_e} \left(1 - \phi \right) \tag{44}$$

where ϕ is the matrix porosity at the time of intrusion

Equation (44) indicates that the solid volume of material removed (spalled) due to a borehole penetration is a function of the flowrate up the borehole, the physical extend of a repository room, the waste porosity, and a velocity above which erosion occurs in waste fractures. The gas velocity up the borehole is a known function of the gas pressure at the borehole entrance.

Erosion Velocity (within fractures) 2.4.1.3.2

The fracture erosion velocity v_{a} is the average threshold gas velocity in a fracture required to "erode" the fracture walls. It is assumed to be related to the terminal velocity of a waste particle and to the cohesive strength between particles caused by pore water and cementation.

Equations 23 and 37 relate the gas velocity required to loft a waste particle to the weight of the particle and to the tensile strength required to separate two particles that are bonded by pore water and cementation. The resulting equation

 $v_e^2 = \frac{4g_{eff}d(\rho_s - \rho_g)}{3C_D\rho_e}$ depends on an effective gravity term

$$g_{eff} = (g + g_t) = g + \frac{3}{4} \frac{\sigma}{\rho_s R} = g + \frac{3}{4} \frac{\sigma_p}{\rho_s R} + \frac{3}{4} \frac{\sigma_c}{\rho_s R}$$
 where the parting stress (equation

37) is assumed to have contributions resulting from pore water σ_p and intergranular cementation σ_c . For the channeling model, since particle erosion is occurring in the channels rather than lofting, correction factors are applied to the parting stresses and acceleration of gravity. Thus the effective gravity term becomes

$$g_{eff} = gF_{ge} + \frac{3}{4} \frac{\sigma_{p}}{\rho_{c}R} F_{se} + \frac{3}{4} \frac{\sigma_{c}}{\rho_{c}R} F_{ce}$$
(45)

where $F_{g_{g}}$ is the gravity effectiveness factor and $F_{g_{g}}$ and $F_{c_{g}}$ are the stress effectiveness factors. These constants are determined by experiment. Since the fracture erosion velocity v_{g} is dependent upon a coefficient of drag C_{p} which is in turn a function of v_{g} (see equation 24) an iterative solution process is used to converge to a solution for v_{g} .

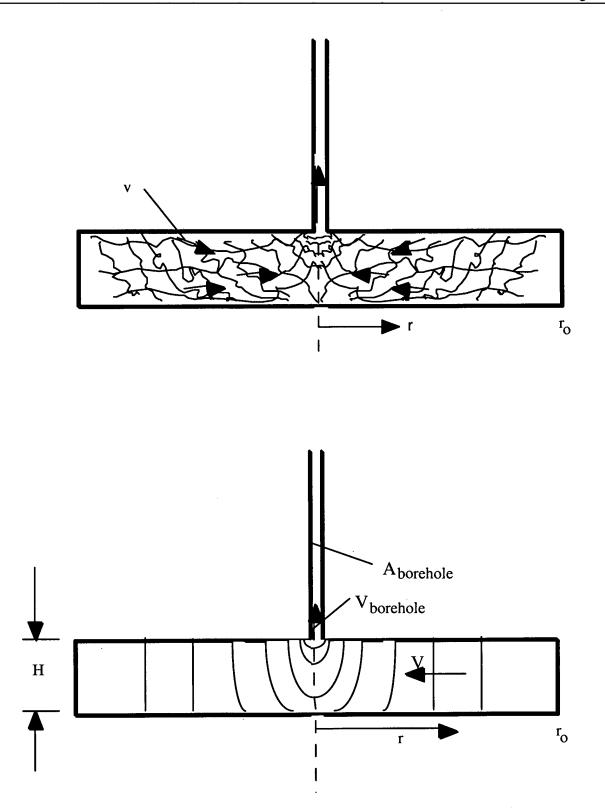


Figure 8 Flow Model

2.4.1.3.3 Solution Process for Solids Removal caused by Blowout (Model 2)

The solution to equation 44 requires 5 quantities. The value r_o is chosen as the equivalent radius of a repository room based on area. For a WIPP room $r_o \sim 17$ m. The borehole area $A_{borehole}$ is detemined by the annulus area or (Figure 7)

$$A_{borehole} = \frac{\pi}{4} \left(D_o^2 - D_i^2 \right)$$

To determine V_{borehole} requires the solution of equation 22. As indicated earlier, flow in the borehole annulus is assumed to occur adjacent to both the drill collars and the drillpipe. Consequently, there are two annular areas that must be considered for flow up the borehole and equation (22) has to be solved over these two regions. The pressure ratio is based on the inlet and exit pressures for the two annuli with pressure compatibility maintained at the drillpipe / drill collar connection.

Utilizing conservation of mass, equation 22 can be divided into the two equations

$$\frac{fl_{coll}}{D_{coll}} = \frac{J}{KM_1^2} \left[J - \frac{p_i^2}{p_j^2} \right] - 2\ln\frac{p_j}{p_i}$$
(46)

$$\frac{fl_{drill}}{D_{drill}} = \frac{1}{KM_{j}^{2} \left(\frac{p_{1}}{p_{i}}\right)^{2} \left(\frac{D_{coll}}{D_{drill}}\right)^{4}} \left[1 - \frac{p_{exit}^{2}}{p_{i}^{2}}\right] - 2\ln\frac{p_{i}}{p_{exit}}$$
(47)

where p_{exit}= exit pressure

 $\label{eq:ldrill} \begin{array}{l} l_{drill} = length \ of \ drillpipe \\ l_{coll} = collar \ length \\ D_{coll} = equivalent \ diameter \ for \ collar \ annulus \ (based \ on \ annulus \ area) \\ D_{drill} = equivalent \ diameter \ for \ drillpipe \ annulus \ (based \ on \ annulus \ area) \\ p_i = intermediate \ pressure \\ M_i = inlet \ Mach \ number \end{array}$

In addition to equations 46 and 47 the flow requires the additional constraint (Binder, 1958 p 64) that either $\frac{dp}{dl}$ becomes unbounded at the pipe exit or $p_{exit} = p_{atm}$. A recursive process based on Taylor series expansions (appendix AA) is used to obtain the solution to the three governing equations. Seed values

(48)

used as initial guesses to the solution are obtained by systematically scanning over possible values of p_i and p_{exit} for a fixed backpressure (p_1 = waste gas pore pressure).

The borehole velocity is related to the inlet Mach number by the relation

 $V_{borehole} = CM_{\mu}$

where $C = (KRT)^{\frac{1}{2}}$ is the local sound speed, *T* is the absolute temperature and *K* and *R* are gas constants. For hydrogen gas K=1.41 and R=4116. $\frac{Nm}{kg.s}$ (Obert, 1948, p541)

2.4.1.3.4 Determination of the factors F_{aa} and F_{sa}

The gravity effectiveness factor F_{go} and the stress effectiveness factors F_{so} and F_{co} are required in equation (45) to compute an effective gravity term for use in model 2 (Section 2.4.1.3.1). Laboratory scale blowout experiments (SNL Contract AL 7022) have been performed which produce data on the volume of material ejected under various conditions of borehole diameter, gas pressure, waste particle size, etc. For arbitrarily selected values of the parameters F_{go} , F_{so} and F_{co} various degrees of fit to the data can be achieved. This process is conducted within the CUTTINGS_S code prior to performing any predictions of full scale response. In the present analysis a best fit is achieved when the sum of the squares of the differences in computed volumes and experimental volumes is a minimum for all data points.

Defining the error as E

$$E = \sum_{l}^{n} \left[V_{ei} - V_{mi} \left(F_{se}, F_{ge}, F_{ce} \right) \right]^{\circ}$$

where

 V_{ei} is the volume of waste simulant ejected in experiment *i*

 $V_{mi}(F_{se}, F_{ge}, F_{ce})$ is the computed volume of waste simulant ejected (calculated by CUTTINGS_S computer code) *n* is the number of experiments

A best fit is achieved when F_{ae} , F_{se} and F_{ce} are chosen such that E is a minimum.

For the set of *n* data points the computations defined by equation (48) are made within a module of the CUTTINGS_S computer code by scanning over a range of possible values of F_{po} , F_{so} , and F_{co} .

2.4.2 Stuck Pipe

If the waste has a low permeability when penetrated by a drillbit, the gas flow into the drilling mud may cause waste failure adjacent to the borehole (Berglund, 1993) and jam the drillbit preventing further drilling. Prior to becoming completely stuck the driller will notice an increase in torque on the drillstring and a decrease in the rotational speed. When sticking occurs the driller will usually initiate a cleanout procedure wherein the drill bit is raised and lowered repeatedly into the sticking formation to clear the obstruction. This process can be continued for as much as 12-24 hours if it is shown to be effective. After

this time the problem must be solved by weighting up the mud, spot sealing with cement or setting casing (Short, 1982).

During the cleanout procedure waste will be transported to the surface with each thrust of the drillbit into the obstruction. The quantity of waste removed is related to the maximum carrying capacity of the drilling mud and can be estimated based on the observation that for drill cutting loadings above 5% in the drilling mud (Darley & Gray, 1988, p259) tight hole conditions or stuck drillpipe may occur when circulation is stopped for any reason. Thus the maximum solid waste removal rate would consist of 5% of the drilling mud flowrate. The total quantity of solid waste transported to the surface can thus be computed as the waste removal rate multiplied by the cleanout time. The range of releases possible is based on variations in drillbit diameter (10.5 to 17.5 inches), duration of the cleanout procedure (12 to 24 hours) and the drilling mud flowrate (30 to 50 gallons/minute/inch of drill diameter). The releases are based on

$$V_{S}=0.05QD_{O}T$$

(49)

14

where

 V_S = Solid waste volume brought to ground surface Q=Drilling mud flow rate/drillbit diameter T=Cleanout duration D_o = Drill bit diameter

and varies between 43 to 238 m³ of solid waste.

The lower limit of repository gas pressure at which sticking would occur (10MPa) is based on a drillstring power of 800hp and a coefficient of friction between the waste and drillcollars of 0.3. For these conditions the drill string angular velocity would decrease by more than 50% from the normal operating range alerting the driller to sticking conditions.

2.4.3 Gas Erosion

This again occurs when the waste has a low permeability and the gas flow to the borehole is very low and either is not detected by the driller or is allowed to trickle slowly along the drillstring and be released at the surface. The waste fails adjacent to the borehole perhaps causing some pipe sticking but the driller is able to continue and does not detect the unusual nature of the drill cuttings being brought to the surface. The flow of gas from the waste to the borehole generates a stress state in the waste adjacent to the borehole that depending on waste strength and the magnitude of the pressure gradient impresses the failed waste against the drillstring causing a continuous process of gas assisted erosion. As waste erodes more waste moves towards the drillstring in response to the gas pressure gradient and the process continues until sufficient gas has been released from the panel to preclude waste failure or until casing is set. Because the driller is either not aware of, or ignores the nature of the drill cuttings being removed in this case, the final volume of waste removed can be substantial.

For compacted waste with little or no strength waste failure will generally occur for all repository gas pressures that exceed the hydrostatic stress of the drilling mud. The failed waste is then transported to the accessible environment in the drilling mud. As with the sticking mode described above, the volume of waste removed can be computed based on the observation that above drill cutting loadings above 5% in the drilling mud (Darley & Gray, 1988, p259) tight hole conditions or stuck drillpipe may occur when circulation is stopped for any reason. Thus under these conditions the driller is not likely to remove waste to the surface at a rate faster than continuous drilling at the 5 percent limit.

The 5 percent cuttings loading will consist of both cuttings (from the hole bottom) and gas spallings. For a fixed mud flowrate the cuttings percentage will vary with the penetration rate. The penetration rate varies

between 50 - 100 ft/hr. For high penetration rates the cuttings percentage will be high leaving only a small amount for spallings to add up to the assumed 5 percent limit. For low penetration rates the cuttings percentage will be small and the spallings percentage correspondingly greater. The quantity of waste removed to the surface will be equal to (spall percentage) X (mud flowrate) X (drilling time). The drilling time is governed by the time required to drill from the elevation of the repository to the elevation at which casing is set which is below the Castile formation at 4500 ft.

The solid volume of waste brought to the surface can by readily computed based on available drilling parameters utilizing the following equation.

 $V_{s} = [0.05 QD_{o} - \pi (D_{o}^{2}/4.)R_{p}]\Delta/R_{p}$ (50)

where

 V_S = Solid waste volume brought to ground surface Q=Drilling mud flow rate/drillbit diameter Do=Drill bit diameter Rp=Penetration rate

Δ=Differential Drilling Depth (distance from repository depth to depth where casing is set)

Based on a differential drilling depth of 2350 ft, and placing the remaining variables at their extreme values the range of volumes of solid waste released to the surface ranges from 44 to 356 m³.

3.0 Radionuclide Chains and Decay

The radionuclide inventory contained in the waste decays with time and follows several decay pathways or chains until stable elements are attained. The number of atoms of each radioactive waste constituent thus varies with time and the activity of waste brought to the surface as the result of human intrusion must account for the initial radionuclide inventory, the decay pathways, the mass of waste released, and the time of intrusion.

The radioactive decay of a chain of radionucides is governed by the Bateman equations and is described in Kaplan, 1962.

Consider a chain of n radionuclides where N_1 , N_2 , N_3 , N_d N_n represent the number of atoms of each of the radionuclides in the repository and where N_n is stable. The differential equations that govern the decay and growth are

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$$\frac{dN_1}{dt} = -\lambda_1 N_1$$
$$\frac{dN_2}{dt} = \lambda_1 N_1 - \lambda_2 N_2$$
$$\frac{dN_3}{dt} = \lambda_2 N_2 - \lambda_3 N_3$$

(51)

(52)

$$\frac{dN_n}{dt} = \lambda_{n-1} N_{n-1}$$

It will be assumed that the initial number of atoms of radionuclide 1 in the repository is N_1^0 and the initial number of atoms of the daughter atoms are N_2^0 , N_3^0 , N_4^0 N_n^0 . The disintegration constants are λ_1 , λ_2 , λ_3 , λ_4 λ_{n-1} . The half lives of the radionuclides are related to the disintegration constants through the relation, $t_1 = \text{half life} = \frac{\ln 2}{\lambda}$. Solving the differential equations (7) sequentially

obtains

where t is the decay time at the time of intrusion.

 $N_n = \sum_{j=1}^n N_j^0 \prod_{i=j}^{n-1} \lambda_i \sum_{i=j}^n \left(\prod_{l=j+1}^n \frac{1}{\lambda_l - \lambda_i} \right) e^{-\lambda_i t}$

Additional details of this derivation are shown in Appendix AC. The activity (in curies) of each radionucide is related to the number of atoms through the relation (Sandia

WIPP Project. 1992, Table 15)
$$A_i = \frac{1.128 \times 10^{10} N_i}{\frac{t_1 M_i}{2}}$$
 where M_i is the gram molecular weight

of the ith radionuclide. The activity in curies of the ith radionuclide <u>released to the surface</u> resulting from a human intrusion (a_i) is determined by mutiplying the repository activity A_i by the percent of waste removed or

$$a_i = A_i \left(\frac{Volume \ Released}{Re \ pository \ Volume} \right) = A_i \left(\frac{Projected \ Area \ Re \ moved}{Re \ pository \ Area} \right)$$

Appendix AA

Recursive process for solving a system of non-linear equations (Salvadori, et al 1962)

Given a system of non-linear equations (in our case 3 equations) in the unknowns x,y,z of the form

$$F_{1}(x, y, z) = 0$$

$$F_{2}(x, y, z) = 0$$
(A1)
$$F_{3}(x, y, z) = 0$$

A recursive solution procedure can be developed by expanding the equations in Taylor series about an initial guess (x_0, y_0, z_0) for the solution.

$$F_{1}(x, y, z) = F_{1}(x_{0}, y_{0}, z_{0}) + \frac{\partial F_{1}(x_{0}, y_{0}, z_{0})}{\partial x}(x - x_{0}) + \frac{\partial F_{1}(x_{0}, y_{0}, z_{0})}{\partial y}(y - y_{0}) + \frac{\partial F_{1}(x_{0}, y_{0}, z_{0})}{\partial z}(z - z_{0}) + \dots = 0$$

$$F_{2}(x, y, z) = F_{2}(x_{0}, y_{0}, z_{0}) + \frac{\partial F_{2}(x_{0}, y_{0}, z_{0})}{\partial x}(x - x_{0}) + \frac{\partial F_{2}(x_{0}, y_{0}, z_{0})}{\partial y}(y - y_{0}) + \frac{\partial F_{2}(x_{0}, y_{0}, z_{0})}{\partial z}(z - z_{0}) + \dots = 0$$

$$F_{3}(x, y, z) = F_{3}(x_{0}, y_{0}, z_{0}) + \frac{\partial F_{3}(x_{0}, y_{0}, z_{0})}{\partial x}(x - x_{0}) + \frac{\partial F_{3}(x_{0}, y_{0}, z_{0})}{\partial y}(y - y_{0}) + \frac{\partial F_{3}(x_{0}, y_{0}, z_{0})}{\partial z}(z - z_{0}) + \dots = 0$$
(A2)

Retaining only the linear terms of the Taylor series and defining

$$\delta x = x - x_o$$

$$\delta y = y - y_o$$

$$\delta z = z - z_o$$
(A3)

three linear simultaneous equations result

$$\frac{\partial F_1(x_o, y_o, z_o)}{\partial x} \delta x + \frac{\partial F_1(x_o, y_o, z_o)}{\partial y} \delta y + \frac{\partial F_1(x_o, y_o, z_o)}{\partial z} \delta z + F_1(x_o, y_o, z_o) = 0$$

$$\frac{\partial F_2(x_o, y_o, z_o)}{\partial x} \delta x + \frac{\partial F_2(x_o, y_o, z_o)}{\partial y} \delta y + \frac{\partial F_2(x_o, y_o, z_o)}{\partial z} \delta z + F_2(x_o, y_o, z_o) = 0$$
(A4)
$$\frac{\partial F_3(x_o, y_o, z_o)}{\partial x} \delta x + \frac{\partial F_3(x_o, y_o, z_o)}{\partial y} \delta y + \frac{\partial F_3(x_o, y_o, z_o)}{\partial z} \delta z + F_3(x_o, y_o, z_o) = 0$$

Based on an initial trial solution (x_0 , y_0 , z_0) sufficiently "close" to the actual solution, the three linear equations (A4) can be solved for the increments δx , δy , and δz and new values for x, y, and z can be determined from equations (A3). This process can be repeated until $|\delta x|$, $|\delta y|$, and $|\delta z|$ are less than a specified limit.

Appendix AB Velocity ratio factors

Calculation of blowout releases using the model of section 2.4.1 requires the value of the gas velocity escaping from a cylindrical waste surface 700s after the drill bit has breached the repository. Since the solution process for the final blowout radius is iterative, the required gas velocity has to be determined many times using a large number of blowout radii. Although this velocity can be computed using the finite difference code GASFLOW, the code is slow and the total computation time necessary would be prohibitive. Therefore rather than using the results of the GASFLOW directly in the computational model a corrected steady state model is used which can be solved quickly in closed form. To calculate conservative blowout releases the steady state model should generate releases at least as great as those that would be predicted using the computationally intensive GASFLOW code. The correction factors

required for this are determined by running both GASFLOW and the steady state model over a range of boundary and initial conditions and by computing the velocity ratios given by

 $V_{R} = velocity \ ratio = \frac{Velocity \ from \ GASFLOW @ 700s}{Steadystate \ velocity \ from \ equation \ 28}$

Tabulated values for V_R are given below for a number of cases which cover a wide variable range. By choosing the largest V_R value for a given permeability (indicated in each table) a fit to the data using permeability as the independent variable can be made.

k=10 ⁻¹⁶ m ²	p∞=8 MPa p ₁ =1Mpa			$p_{\infty}=8 \text{ MPa} p_1=1 \text{ Mpa} p_{\infty}=15 \text{ MPa} p_1=1 \text{ Mpa}$			ІМра
R _b (m)	$\varphi = 0.09$	<i>φ</i> = 0.19	<i>φ</i> = 0.29	$\varphi = 0.09$	<i>φ</i> = 0.19	$\varphi = 0.29$	
0.1	2.41	2.79	3.04	2.17	2.47	2.68	
0.3	3.20	3.90	4.40	2.76	3.30	3.69	
1.0	5.09	6.68	7.87	4.13	5.31	6.2	
3.0	8.82	12.23	14.8	6.81	9.32	11.2	
8.0	14.4	20.5	25.1	10.8	.15.3	18.7	
13.0	17.4	24.9	30.6	12.9	18.5	22.7	
20.0	19.0	27.4	33.7	14.1	20.3	24.9	

k=10 ⁻¹⁵ m ²	p∞=8 MPa p ₁ =1Mpa			p∞=	15 MPa p ₁ =1	Мра
R _b (m)	$\varphi = 0.09$	<i>φ</i> = 0.19	<i>φ</i> = 0.29	$\varphi = 0.09$	$\phi = 0.19$	$\varphi = 0.29$
0.1	1.69	1.87	1.99	1.56	1.71	1.82
0.3	1.96	2.25	2.45	1.76	2.00	2.16
1.0	2.52	3.08	3.48	2.17	2.6	2.91
3.0	3.55	4.64	5.46	2.89	3.71	4.31
8.0	5.07	6.98	8.42	3.94	5.35	6.41
13.0	5.86	8.21	9.99	4.47	6.21	7.51
20.0	6.25	8.85	10.8	4.71	6.62	8.07

k=10 ⁻¹⁴ m ²	p∞=8 MPa p ₁ =1Mpa			p∞=	15 MPa p ₁ =	1Mpa
R _b (m)	$\varphi = 0.09$	<i>φ</i> = 0.19	<i>φ</i> = 0.29	$\varphi = 0.09$	$\varphi = 0.19$	$\varphi = 0.29$
0.1	1.29	1.12	1.46	1.21	1.3	1.36
0.3	1.38	1.53	1.63	1.28	1.40	1.49
1.0	1.54	1.78	1.94	1.40	1.58	1.70
3.0	1.80	2.19	2.47	1.56	1.85	2.07
8.0	2.13	2.75	3.22	1.77	2.22	2.57
13.0	2.28	3.03	3.60	1.86	2.39	2.80
20.0	2.3	3.12	3.74	1.87	2.42	2.87

k=10 ⁻¹³ m ²	p∞=	p∞=8 MPa p ₁ =1Mpa			15 MPa p ₁ =1	Мра
R _b (m)	$\varphi = 0.09$	<i>φ</i> = 0.19	<i>φ</i> = 0.29	$\varphi = 0.09$	<i>φ</i> = 0.19	<i>φ</i> = 0.29
0.1	1.17	1.18	1.19	1.11	1.18	1.19
0.3	1.20	1.23	1.24	1.12	1.20	1.22
1.0	1.22	1.29	1.32	1.11	1.24	1.28
3.0	1.25	1.38	1.44	1.09	1.28	1.35
8.0	1.24	1.46	1.58	1.00	1.28	1.41
13.0	1.18	1.48	1.63	0.881	1.23	1.40
20.0	1.03	1.40	1.60	0.677	1.09	1.31

k=10 ⁻¹² m ²	p∞=8 MPa p ₁ =1Mpa			p∞=	15 MPa p ₁ =1	Мра
R _b (m)	$\varphi = 0.09$	$\varphi = 0.19$	<i>φ</i> = 0.29	<i>φ</i> = 0.09	<i>φ</i> = 0.19	φ=0.29
0.1	0.692	0.929	1.03	0.460	0.735	0.872
0.3	0.631	0.897	1.01	0.394	0.678	0.831
1.0	0.537	0.836	0.980	0.305	0.587	0.758
3.0	0.414	0.734	0.911	0.208	0.464	0.645
8.0	0.257	0.558	0.762	0.111	0.301	0.467
13.0	0.164	0.416	0.617	0.0646	0.200	0.337
20.0	0.0816	0.253	0.419	0.0292	0.106	0.198

The values of V_R and permeability k are thus

VR	k
33.7	10 ⁻¹⁶
10.8	10 ⁻¹⁵
3.74	10 ⁻¹⁴
1.63	10 ⁻¹³
1.03	10 ⁻¹²

These can be fit with the function

$$V_{\rm p} = -12593 - 5040\xi - 806.18\xi^2 - 64.455\xi^3 - 2.5771\xi^4 - 0.04125\xi^5$$

where

$$\xi = \log_{10} k$$

Appendix AC

Solution of the Bateman Equations

Consider a chain of n radionuclides where N_1 , N_2 , N_3 , N_d N_n represent the number of atoms of each of the radionuclides in the repository and where N_n is stable. The differential equations that govern the decay and growth are

(C2)

 $\frac{dN_{I}}{dt} = -\lambda_{I}N_{I}$ $\frac{dN_{2}}{dt} = \lambda_{I}N_{I} - \lambda_{2}N_{2}$ $\frac{dN_{3}}{dt} = \lambda_{2}N_{2} - \lambda_{3}N_{3}$ (C1) $\frac{dN_{n}}{dt} = \lambda_{n-I}N_{n-I}$

The initial number of atoms of radionuclide 1 in the repository is N_1^0 and the initial number of atoms of the daughter atoms are N_2^0 , N_3^0 , N_4^0 N_n^0 . The disintegration constants are $\lambda_1, \lambda_2, \lambda_3, \lambda_4$ λ_{n-1} . The half lives of the radionuclides are related to the disintegration constants through the relation, $t_{\frac{1}{2}}$ = half life = $\frac{\ln 2}{\lambda}$.

The Laplace transform of N_i and $\frac{dN_i}{dt}$ will be defined by the notation (Hildebrand, 1962) $L\{N_i\} = f_i(s)$ and $L\{\frac{dN_i}{dt}\} = sf_i(s) - N_i^0$

Taking the Laplace transform of both sides of equations C1 results in

$$sf_{1}(s) - N_{1}^{0} = -\lambda_{1} f_{1}(s)$$

$$sf_{2}(s) - N_{2}^{0} = \lambda_{1} f_{1}(s) - \lambda_{2} f_{2}(s)$$

$$sf_{3}(s) - N_{3}^{0} = \lambda_{2} f_{2}(s) - \lambda_{3} f_{3}(s)$$
....
$$sf_{n}(s) - N_{n}^{0} = \lambda_{n-1} f_{n-1}(s)$$

or solving for $f_1(s)$ $f_n(s)$

(C3)

(C4)

$$f_{1}(s) = \frac{N_{1}^{0}}{s + \lambda_{1}}$$

$$f_{2}(s) = \frac{N_{2}^{0} + \lambda_{1}f_{1}(s)}{s + \lambda_{2}} = \frac{N_{2}^{0}}{s + \lambda_{2}} + \frac{\lambda_{1}N_{1}^{0}}{(s + \lambda_{1})(s + \lambda_{2})}$$

$$f_{3}(s) = \frac{N_{3}^{0} + \lambda_{2}f_{2}(s)}{s + \lambda_{3}} = \frac{N_{3}^{0}}{s + \lambda_{3}} + \frac{\lambda_{2}N_{2}^{0}}{(s + \lambda_{2})(s + \lambda_{3})} + \frac{\lambda_{2}\lambda_{1}N_{1}^{0}}{(s + \lambda_{1})(s + \lambda_{2})(s + \lambda_{3})}$$

$$\dots$$

$$f_{n}(s) = \frac{N_{n}^{0}}{s + \lambda_{n}} + \frac{\lambda_{n-1}N_{n-1}^{0}}{(s + \lambda_{n})(s + \lambda_{n-1})} + \frac{\lambda_{n-1}\lambda_{n-2}N_{n-2}^{0}}{(s + \lambda_{n})(s + \lambda_{n-1})} + \dots$$

$$\cdot + \frac{\lambda_{n-1}\lambda_{n-2}\dots\lambda_{1}N_{1}^{0}}{(s + \lambda_{n})(s + \lambda_{n-1})\dots(s + \lambda_{1})}$$

The terms of equations (C3) can be rewritten in terms of partial fractions as

$$f_{1}(s) = \frac{N_{1}^{0}}{s + \lambda_{1}}$$

$$f_{2}(s) = \frac{N_{2}^{0}}{s + \lambda_{2}} + \frac{\lambda_{1}N_{1}^{0}}{(s + \lambda_{1})(\lambda_{2} - \lambda_{1})} + \frac{\lambda_{1}N_{1}^{0}}{(s + \lambda_{2})(\lambda_{1} - \lambda_{2})}$$

$$f_{3}(s) = \frac{N_{3}^{0}}{s + \lambda_{3}} + \frac{\lambda_{2}N_{2}^{0}}{(s + \lambda_{2})(\lambda_{3} - \lambda_{2})} + \frac{\lambda_{2}N_{2}^{0}}{(s + \lambda_{3})(\lambda_{2} - \lambda_{3})} + \frac{\lambda_{1}\lambda_{2}N_{1}}{(s + \lambda_{1})(\lambda_{2} - \lambda_{1})(\lambda_{3} - \lambda_{1})} + \frac{\lambda_{1}\lambda_{2}N_{1}}{(s + \lambda_{2})(\lambda_{3} - \lambda_{2})(\lambda_{1} - \lambda_{2})} + \frac{\lambda_{1}\lambda_{2}N_{1}}{(s + \lambda_{3})(\lambda_{1} - \lambda_{3})(\lambda_{2} - \lambda_{3})}$$

$$f_n(s) = \sum_{j=1}^n \frac{N_j}{(s + \lambda_n)} \prod_{i=j}^{n-1} \left(\frac{\lambda_i}{s + \lambda_i}\right)$$

Inverting equations (C4) back into the time domain obtains

 $N_1 = N_1^0 e^{-\lambda_1 t}$ $N_{2} = N_{2}^{0} e^{-\lambda_{1}t} + \lambda_{1} N_{1}^{0} \left[\frac{e^{-\lambda_{1}t}}{(\lambda_{2} - \lambda_{1})} + \frac{e^{-\lambda_{2}t}}{(\lambda_{1} - \lambda_{2})} \right]$ $N_3 = N_3^0 e^{-\lambda_1 t} + \lambda_2 N_2^0 \left[\frac{e^{-\lambda_2 t}}{(\lambda_3 - \lambda_2)} + \frac{e^{-\lambda_3 t}}{(\lambda_2 - \lambda_3)} \right]$ $+\lambda_1\lambda_2N_1^0\left[\frac{e^{-\lambda_1t}}{(\lambda_2-\lambda_1)(\lambda_3-\lambda_1)}+\frac{e^{-\lambda_2t}}{(\lambda_3-\lambda_2)(\lambda_1-\lambda_2)}+\frac{e^{-\lambda_3t}}{(\lambda_1-\lambda_3)(\lambda_2-\lambda_3)}\right]$

 $N_n = \sum_{i=1}^n N_j^0 \prod_{i=1}^{n-1} \lambda_i \sum_{i=1}^n \left(\prod_{i=i+1}^n \frac{1}{\lambda_i^{-1} - \lambda_i} \right) e^{-\lambda_i t}$

Appendix AD Parting Stress Based on Pore Water

Consider two spherical particles in contact with a small volume of liquid surrounding the point of contact (Figure D1). Assume that the surface tension is T and the pressure within the liquid is P_{e} and outside the liquid is P_{e} . For static equilibrium of a small element of the liquid surface

$$\left(P_c - P_o\right) ds' ds'' + 2T \sin\left(\frac{d\theta'}{2}\right) ds'' - 2T \sin\left(\frac{d\theta''}{2}\right) ds' = 0$$
(D1)

or

$$\left(P_{c}-P_{o}\right)ds^{'}ds^{''}+2T\left(\frac{d\theta^{''}}{2}\right)ds^{''}-2T\left(\frac{d\theta^{''}}{2}\right)ds^{'}=0$$
(D2)

thus

$$(P_c - P_o) ds' ds'' + T d\theta' ds'' - T d\theta'' ds' = 0$$
but
$$(D3)$$

$$ds' = r'd\theta'$$
 and $ds'' = r''d\theta'$

Therefore

$$P_c - P_o = T \left(\frac{l}{r} - \frac{l}{r} \right)$$
(D4)

The quantity $P_c - P_o$ represents the capillary pressure in the liquid volume. To physically separate the two spherical particles requires a force F

$$F = 2\pi r'' T - (P_c - P_o)\pi r''^2$$
(D5)

The parting stress is based on the projected area of a particle and the force F. It can be written

$$\sigma_{p} = \frac{F}{\pi R^{2}} = \frac{2\pi r^{"} T - (P_{c} - P_{o})\pi r^{"2}}{\pi R^{2}} = \frac{2r^{"} T - T\left(\frac{1}{r^{"}} - \frac{1}{r^{'}}\right)r^{"2}}{R^{2}}$$
(D6)

where from geometry (Figure D2)

$$r'' = -r' + \sqrt{r'^2 + 2r'R}$$
 (D7)

The above analysis assumes that only one fluid contact point is involved in the determination of the parting stress. For more densely packed material as many as three fluid contact points should be considered resulting in a greater stress. However to be conservative only one fluid contact point is assumed.

(D8)

Waste Saturation

The volume of liquid associated with each particle contact is

$$V_w = 2\pi w^2 h - 2V_R - V_c$$

Where V_B is the volume associated with a portion of the particle formed by a plane passing through the particle a distance h from the particle surface, and V, is the toroidal volume defined by an area A rotated around axis a'. The area of the latter is delineated by the straight line 2h and a portion of the circle of radius r' (Figure D3).

$$V_{R} = \frac{1}{3}\pi h^{2}(3R - h)$$

$$V_{c} = 2\pi A \left[r' + r' - \frac{2r}{3} \left(\frac{\sin^{3} \theta}{\theta - \sin\theta \cos \theta} \right) \right]$$
(D9)
(D10)

where

$$A = \frac{r^{\prime 2}}{2} (2\theta - \sin 2\theta) \tag{D11}$$

The liquid saturation is defined as the ratio of the liquid volume to the total void volume in a sample of material. Void volume is a function of particle shape, particle size distribution, and packing density. For spherical particles of a single diameter the maximum percent voids (void ratio) is 37.5% (Katz, HS, et al 1978). The number of single size particles (N) within a volume (V) of material at maximum packing density is

 $N = \frac{(1 - 0.375)V}{\frac{4}{3}\pi R^3}$ Within the volume V each particle is in contact with 12 surrounding particles. Thus the volume of liquid

associated with N particles is

$$V_{wt} = 6 V_w N$$

And the saturation (S) is

(D12)

(D13)

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$$S = \frac{V_{wt}}{0.375V} \tag{D14}$$

or

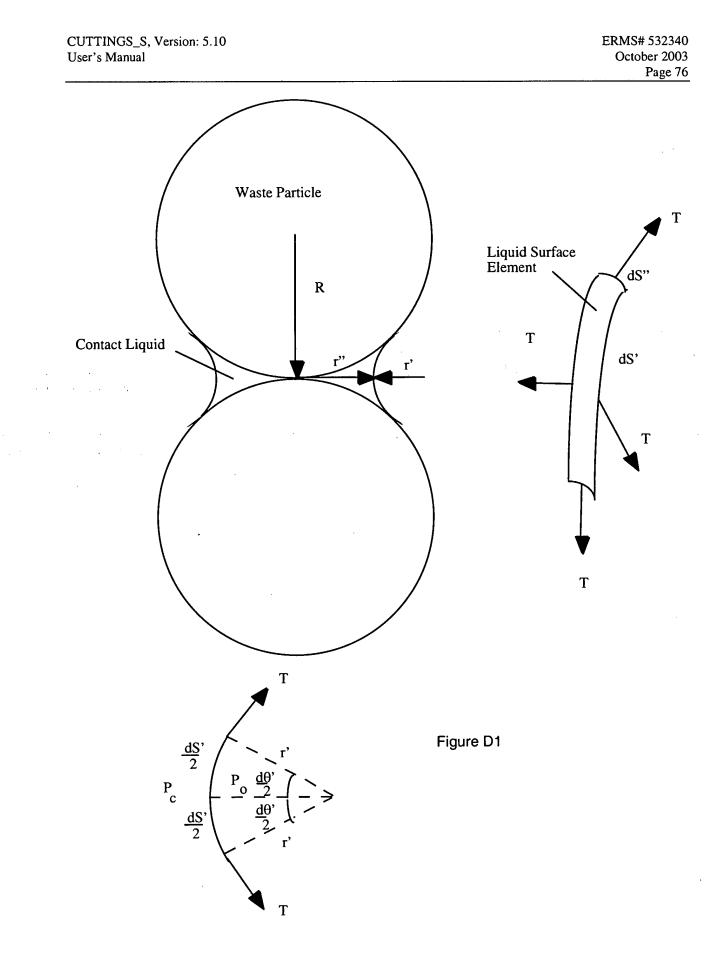
$$S = 2.39 \frac{V_w}{R^3} \tag{D15}$$

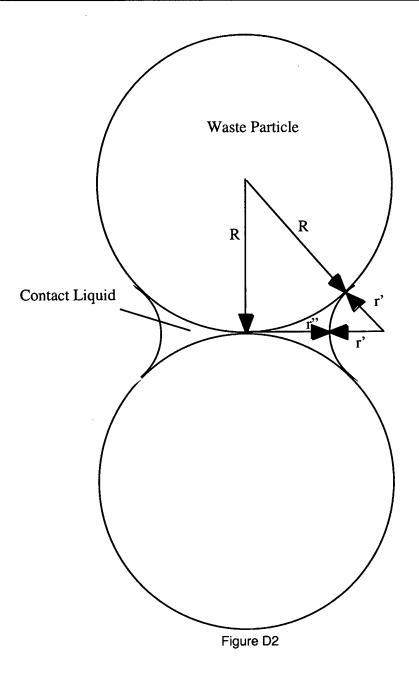
For a known saturation S and particle radius R it is possible to back solve for r' and the parting stress σ_p . This must be done numerically because of the complexity of the governing equations. Using equation (D8) and (D15) a variable Q can be defined thus

$$Q = V_w - \frac{SR^3}{2.39} = 2\pi w^2 h - 2V_R - V_c - \frac{SR^3}{2.39} = 0$$
 (D16)

The iterative numerical procedure used to determine the value of r' that satisfies equation (D16) is illustrated in Figure D4. The value of r' from the solution of equation (D16) can then be used to determine the parting stress using equations (D6) and (D7). The parting stress can then be used to compute the effective acceleration of gravity utilizing equation (37a).

Strength effects arising from moisture content of the waste can thus be included in the blowout model utilizing the waste saturation at the time of intrusion and the assumed waste particle radius. The waste particle radius is an unknown quantity which will have to be sampled.





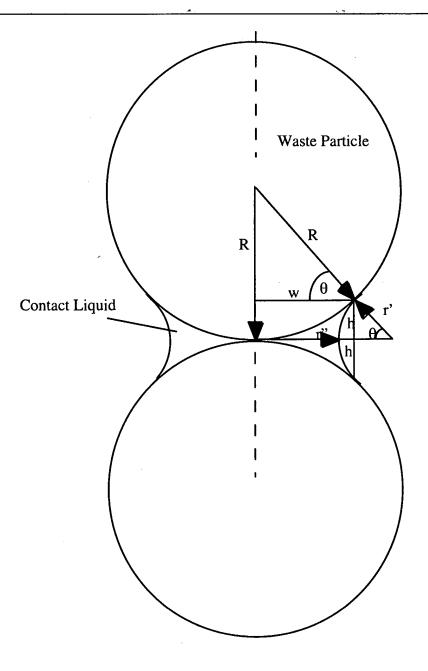


Figure D3

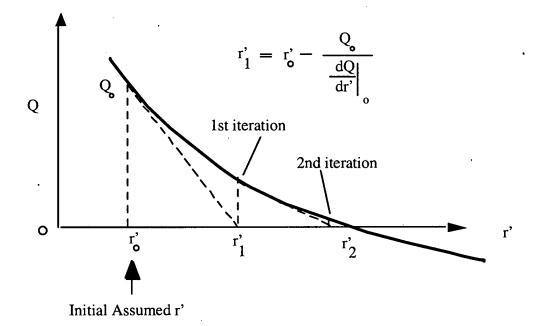


Figure D4

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1. 11.

APPENDIX B: SAMPLES OF INPUT FILES REQUIRED TO RUN GENMESH, MATSET, AND POSTLHS IN PREPARATION TO EXERCISING CUTTINGS_S

This appendix includes brief command files that exercise GENMESH, MATSET, and POSTLHS, as well as the actual input control files necessary to exercise the specific runs required prior to exercising CUTTINGS_S (see Figure 1 in Section 6).

CUSP_GENMESH.COM

\$! -----\$! MODULE: S! PURPOSE: \$! DATE: \$! AUTHOR: \$! -----\$! \$ SET NOON EXE_SRC :== \$WP\$PRODROOT: [GM.EXE]GM_PA96.EXE \$ \$! DEFINE GM_UIF\$INPUT GM_CUSP_CCA.INP \$ DEFINE GM_CDB\$OUTPUT GM_CUSP_CCA.CDB \$ \$ DEFINE GM_DBG\$OUTPUT GM_CUSP_CCA.DBG \$! EXE_SRC GM_UIF\$INPUT GM_CDB\$OUTPUT GM_DBG\$OUTPUT \$ \$! -----\$ EXIT

GM_CUSP_CCA.INP

```
1
! TITLE: CUSP Input GENMESH file
! ANALYST: Robert A. Cole
1
1
*SETup_grid
 DIMension= 3
 ORIgin= 0.0000E+00,
                          0.0000E+00
                 0.0000E+00,
 IJKmax=
          2,
                2,
ŧ
*GRID_spacing
 DEL,COORD=X,DEL= 1.0000E+00,INRANGE=
DEL,COORD=Y,DEL= 1.0000E+00,INRANGE=
                                      2
                                1,
                                      2
                                1,
                                      2
 DEL,COORD=Z,DEL= 1.0000E+00,INRANGE=
                                1,
T
*REGion
                      1,
                                               2, KRANGE=
 REGion=
           1, IRANGE=
                           2, \text{JRANGE} = 1,
      2
1,
1
*END
```

CUSP_MATSET.COM

```
$!
  EXE_SRC :== $WP$PRODROOT: [MS.EXE] MATSET_PA96.EXE
$
$!
  DEFINE MS_CDB$INPUT GM_CUSP_CCA.CDB
$
 DEFINE MS_UIF$INPUT MS_CUSP_CCA.INP
$
 DEFINE MS_SDB$INPUT WIPP::WIPP_DB
$
$
 DEFINE MS_SDB_CALC$INPUT CCA4
$
 DEFINE MS_CDB$OUTPUT MS_CUSP_CCA.CDB
$
 DEFINE MS_DBG$OUTPUT MS_CUSP_CCA.DBG
$!
  EXE_SRC MS_CDB$INPUT MS_UIF$INPUT MS_SDB$INPUT MS_SDB_CALC$INPUT -
Ś
         MS_CDB$OUTPUT MS_DBG$OUTPUT
$!
$
  EXIT
```

MS_CUSP_CCA.INP

```
1
!
   FILETYPE:
!
  ANALYSTS:
!
  DATE:
1
  PURPOSE:
1
1
1
1
*PRINT_ASSIGNED_VALUES
1
*HEADING
 TITLE, CUSP MATSET INPUT FILE
 SCALE, LOCAL
 SCENARIO, ALL
1
*UNITS=SI
1
*CREATE_blocks
 BLOCK_IDS=2,3,4
1
*RETRIEVE
 COORD, DIM=3, NAMES=X,Y,Z
1
! ... Define region names
 MATERIAL, 1=BLOWOUT, 2=BOREHOLE, 3=DRILLMUD, 4=WAS_AREA
!1...Define BLOWOUT property names
 PROPERTY, MATerial=BLOWOUT, NAMES=PARTDIA
!2...Define BOREHOLE property names
 PROPERTY, MATERIAL=BOREHOLE, NAMES=DIAMMOD, DOMEGA, TAUFAIL
1
```

1

The file that follows is an input control file to exercise POSTLHS in preparation to exercising CUTTINGS_S.

PLHS_CUSP_CCA.INP ------

TITLE: 1996 CCA Input File for Realization R1 for the PRELHS Code 1 ANALYST: Lanny Smith, May 3, 1996 1 1 DESCRIPTION: 1 1 WIPP 1996 CCA, Realization R1 PRELHS Input File 1 1 This input file to PRELHS is used to generate, as an output file, an LHS 1 input file containing all distribution information and execution options 1 required to create a sample for Realization R1 for the WIPP 1996 CCA 1 analysis. 1 1 !======= No Comments Allowed between *ECHO and *ENDECHO =============== * ECHOLHS TITLE 1996 CCA, Realization R1 Input File for the LHS Code NOBS 100 RANDOM SEED 238766283 CORRELATION MATRIX 3 18 19 -0.99 20 21 -0.99 28 29 -0.75 OUTPUT CORR HIST DATA * ENDECHO 1 !== PROPERTIES TO BE RETRIEVED FROM WIPP 1996 CCA DATABASE, 'CCA4' == 1 *RETRIEVE 1 MATERIALS, STEEL PROPERTIES, CORRMCO2 1 MATERIALS, WAS_AREA

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PROPERTIES, PROBDEG 1 MATERIALS, WAS_AREA PROPERTIES, GRATMICI Ţ MATERIALS, WAS_AREA PROPERTIES, GRATMICH 1 MATERIALS, CELLULS PROPERTIES, FBETA 1 MATERIALS, WAS_AREA PROPERTIES, SAT_RGAS 1 MATERIALS, WAS_AREA PROPERTIES, SAT_RBRN ! MATERIALS, WAS_AREA PROPERTIES, SAT_WICK 1 MATERIALS, CL_L_T1 PROPERTIES, PRMX_LOG I MATERIALS, CONC_T1 PROPERTIES, PRMX_LOG ! MATERIALS, ASPHALT PROPERTIES, PRMX_LOG ī MATERIALS, SHFT_DRZ PROPERTIES, PRMX_LOG 1 MATERIALS, SALT_T1 PROPERTIES, CUMPROB ! MATERIALS, SALT_T1 PROPERTIES, SAT_RGAS ļ MATERIALS, SALT_T1 PROPERTIES, SAT_RBRN ! MATERIALS, SALT_T1 PROPERTIES, PORE_DIS ! MATERIALS, S_HALITE PROPERTIES, POROSITY 1 MATERIALS, S_HALITE PROPERTIES, PRMX_LOG ï MATERIALS, S_HALITE PROPERTIES, COMP_RCK ! MATERIALS, S_MB139 PROPERTIES, PRMX_LOG 1 MATERIALS, S_MB139 PROPERTIES, COMP_RCK 1 MATERIALS, S_MB139

PROPERTIES, RELP_MOD ! MATERIALS, S_MB139 PROPERTIES, SAT_RBRN ţ MATERIALS, S_MB139 PROPERTIES, SAT_RGAS 1 MATERIALS, S_MB139 PROPERTIES, PORE_DIS ï MATERIALS, S_HALITE PROPERTIES, PRESSURE ï MATERIALS, CASTILER PROPERTIES, PRESSURE ī MATERIALS, CASTILER PROPERTIES, PRMX_LOG ! MATERIALS, CASTILER PROPERTIES, COMP_RCK ï MATERIALS, BH_SAND PROPERTIES, PRMX_LOG 1 MATERIALS, CASTILER PROPERTIES, GRIDFLO 1 MATERIALS, BLOWOUT PROPERTIES, PARTDIA 1 MATERIALS, BOREHOLE PROPERTIES, TAUFAIL 1 MATERIALS, CULEBRA PROPERTIES, MINP_FAC i MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ŗ MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ī MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL

PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX Ţ MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX i MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ï MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ļ MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL

PROPERTIES, PRMTZIDX ł MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX i MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ! MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ţ MATERIALS, GLOBAL PROPERTIES, PRMTZIDX ï MATERIALS, GLOBAL PROPERTIES, PRMTZIDX i MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1 MATERIALS, GLOBAL PROPERTIES, PRMTZIDX 1

*END

!

END OF APPENDIX B

APPENDIX C: GRAMMATICAL STRUCTURE OF CUTTINGS_S'S INPUT FILES NOS. 1 & 2

The grammatical structure of CUTTINGS_S's input control files (file numbers 1 and 2) follows the four rules stated below. In Section 6.2, these files are associated with the logical symbols: CUSP_INPT\$TXT0\$, and CUSP_INPT\$TXT0TXT1.

(1) Text files consist of keywords followed by (i) a real number, (ii) an integer, or (iii) a character variable. There is no required order for keywords.

(2) Text following an exclamation point (!) is treated as a comment. That is, it is printed out with the file and normally provides the user with useful explanatory information regarding the file, but it is ignored by the code when the file is read.

(3) Arrays can span more than one line of an input file. Arrays of numbers are terminated by any non-numeric character. Character arrays are terminated by any non-numeric, non-alphabetical character. These include the symbols: <>?:[]-+&%\$#@!.

(4) If input is specified as a character data field, numerical values will be extracted from the upstream (MATSET or POSTLHS) input CDB file. The character data field must have a colon, (:), separating the CAMDAT material name on the left from the CAMDAT property name on the right. No intervening blanks are permitted. Material and property names are restricted to a maximum of 8 characters.

END OF APPENDIX C

APPENDIX D: SAMPLE CUTTINGS_S DEBUG FILE FROM TEST -RUN PROBLEM #1

This file is taken from QA test problem number 1 (see Section 6.4) The file itself is annotated here and there with explanatory comments that describe the various collections of data. Explanatory comments are printed in italics to offset them from the normal content of the Debug file. The Debug file itself is described in Section 6.1 as file number 8.

Explanatory Comment: The data below are the standard WIPP title page and disclaimer, followed by dynamic memory allocation data. End comment. 06/04/96 09:52:46 CUSP_PA96 5.03 PROD PA96 05/23/96 CCCCC · υυ υU SSSSSS PPPPPP PPPPPP AAAAA 9999 6666 PP PP AA AA PP PP AA AA CC UU SS PP PP UU SS PP PP 9 6 9 CC υυ 96 CC υU 9 SSSSS PPPPPP SS PP SS PP PPPPPP PP CC ААААААА 99999 66666 тлт υU 9 66 CC υυ υU AA AA CC PP AA AA 9 6 6 CC υU υU SSSSSS 9999 6666 CCCCC υυυυυ AA PP PP AA Program for computing the quantity of radioactive material (in Curies) brought to the surface as cuttings/spall generated by a drilling operation that penetrates a rad-waste repository. CUSP_PA96 Version 5.03 PROD PA96 Built 05/23/96 Written by Jerry W. Berglund Sponsored by Robert A. Cole & (J. S. Rath) Run on 06/04/96 at 09:52:46 Run on ALPHA AXP BEATLE OpenVMS V6.1 Prepared for Sandia National Laboratories Albuquerque, New Mexico 87185-5800 for the United States Department of Energy under Contract DE-AC04-76DP00789 Disclaimer _____ This computer program was prepared as an account of work sponsored by an agency of the United States Government. Neither the United States Government nor any agency thereof, nor any of their employees, nor any of their contractors, subcontractors, or their employees, makes any warranty, express or implied, or assumes any legal liability or responsibility for the accuracy, completeness, or usefulness of any information, apparatus, product, or process disclosed or represents that its use would not infringe privately owned rights. Reference herein to any specific commercial product, process, or service by trade name, trademark, manufacturer, or otherwise, does not necessarily constitute or imply its endorsement, recommendation, or favoring by the United States Government, any agency thereof or any of their The views and opinions contractors or subcontractors. expressed herein do not necessarily state or reflect those of the United States Government, any agency thereof or any

of their contractors or subcontractors.

Program Name CUT_SPAL Program Version Program UpDated

CUSP\$EXE WP\$PRODROOT: [CUSP.EXE]CUSP_PA96.EXE

SUMMARY OF INPUT & OUTPUT FILES USED IN THIS RUN

CUSP_INP\$TXT0CUSP_TEST_REPOSITORY.DATCUSP_INP\$TXT1CUSP_TEST_1.INPCUSP_INP\$BRAGCDBCUSP_TEST_INP_BF.CDBCUSP_OUT\$DBGCUSP_OUT_TEST_1.DBGCUSP_INP\$CDBCUSP_TEST_1.CDBCUSP_OUT\$CDBCUSP_OUT_TEST_1.CDBCUSP_OUT\$NVERFYCUSP_OUT_TEST_1.VER

Debug flag DBRAG is ON

(SCATXTO)

!

Explanatory Comment: The folowing is an echo of data read from the repository/model input control file (logical symbol CUSP_INP\$TEXTO) End of Comment.

MODEL_DATA

!!!!	The reposit	following input variables are hard wired and are related to the cory modeling paper written by J. W. Berglund (see Appendix A)
! !	PR_MAX PR_MAX	<>The maximum pressure allowed by model 15.0E6
!	PR_MIN PR_MIN	
!	PE_MAX PE_MAX	
!	PE_MIN PE_MIN	
! !	PCUT PCUT	<> Pressure (Pa) which defines the border between the erosion phase and the blowout phase 8.0E6
! !	PGAS	<> Pressure (Pa) which defines the border between the gas erosion phase and the stuck-pipe phase
	PGAS	10.0E6
! !	KCRIT	<> Permeability which defines the phase border the blowout phase and the gas erosion/stuck pipe phase
	KCRIT	1.0E-16

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! ITER <> Number of iterations to solve for reynolds number ITER 1 Gravity effectiveness factor ! FGE <> 18.4 FGE <> Strength effectiveness factor ! FSE FSE 0.0 ! FCE <> Cementation effectiveness factor 0.0 FCE ! CEMENT Cementation stress <> CEMENT 0.0 <> Surface tension (N/N) ! SUFTEN SUFTEN 8.0E-2 ! Properties of the gas (Hydrogen) in the repository <<<<<<<<<<<<<< ł <> Ratio of Specific Heats ! KGAS 1.41E0 KGAS ! RGAS <> Gas Constant for Hydrogen 4116.0 RGAS <> Hydrogen viscosity ! VISC 0.92E-5 VISC ! PSUF <> Surface presure (Pa) PSUF 0.089465E6 ! Properties to do with the repository <<<<<<<<<<<<<<<< ! TREPO <> Temperature of repository (K) TREPO 300.0 <> Distance from repository depth to depth where casing ! DEPTH is set (m) 1 DEPTH 716.0 ! FLWCNST <> Percent volume of material that is carried by Drilling mud (unitless) 1 0.05 FLWCNST <> Height of repository at burial time (4m) ! HREPO 3.96 HREPO <> The equivalent radius of 1 panel ! RPANEL (910.0m**2) 60.87 RPANEL <> The equivalent radius of 1 room (11,640m**2) ! ROOM 17.1ROOM ! RHOS <> Waste particle density (kg/m**3) RHOS 2650.0 ! Variables to do with drilling <<<<<<<<<>> 1 ! L1 <> Collar Length (m) 182.88 L1! L2 <> Drill pipe length (m) 472.12 L2

```
! COLDIA
          <> Collar diameter (m)
                  0.2032004
  COLDIA
! PIPED
          <> Drill pipe diameter (m)
                  0.1143002
  PIPED
! ROUGHP
          <> Friction factor (unitless)
  ROUGHP
                        0.08
1
!
 APORO
          <> A constant in equation to determine permeability as a
               function of porosity
ţ
  APORO
                    1.0E-11
           <> N constant in equation to determine permeability as a
!
 NPORO
               function of porosity
1
                         4.6
 NPORO
1
 END_OF_MODEL_INPUT
÷
ļ
  RADIONUCLIDE_DATA
! GEOMETRY
                1.11520E+05
  INV_AR
  RHW_AR
                1.11520E+05
  WUF
                5.25
!MATERIAL
  OUT_MAT
                BOREHOLE
!REPOSITORY_TYPE
              WIPP
REP_NAME
REP_GEOLOGY HALITE
RADWASte_type CONtact_handled
!RADIOISOTOPE_chains
i
   chain1/chain2 from U234 & down are the same:
1
   (Both chains are required input)
1
                         . VZ
                                                                 \mathbf{N}
!
           \mathbf{X}
                    \backslash /
                                     \mathbf{V}
                                              \mathbf{X}
                                                        \mathbf{V}
            PU242
                    U238
                             TH234
                                      PA234M U234
                                                        TH230
                                                                 RA226
CHAIN1
                    PO218
                             PB214
                                      BI214
                                               PO214
                                                        PB210
                                                                 <
            RN222
CHAIN2
            PU238
                    U234
                              TH230
                                      RA226
                                               RN222
                                                        PO218
                             PO214
                                      PB210
            PB214
                    BI214
                                               <
! chain3/chain4 from PU239 & down are the same:
                    NP239
                             PU239
                                       U235
                                               TH231
                                                        PA231
                                                                 AC227
            AM243
CHAIN3
                                               PB211
                                                        BI211
                                                                 TL207
                                                                          <
                             RN219
                                       PO215
            TH227
                    RA223
                                                        AC227
                                                                 TH227
            CM243
                     PI1239
                             U235
                                       TH231
                                               PA231
 CHAIN4
                                                        TL207
            RA223
                     RN219
                              PO215
                                       PB211
                                               BI211
                                                                 <
  chain5/chain6 from U236 & down are the same:
!
                     CM248
                              PU244
                                       PU240
                                               U236
                                                        TH232
                                                                 RA228
 CHAIN5
            CF252
                                                                 BI212
            AC228
                     TH228
                             RA224
                                      RN220
                                               PO216
                                                        PB212
            PO212
                     <
```

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CHAIN6	CM244 RA224	PU240 RN220	U236 PO216	TH232 PB212	RA228 BI212	AC228 PO212	TH228 <		
CHAIN7	CM245 RA225	PU241 AC225	AM241 FR221	NP237 AT217	PA233 BI213	U233 PO213	тн229 рв209	<	
CHAIN8	CS137	BA137M	<					,	
CHAIN9	PM147	SM147	ND143	<					
CHAIN10	SR90	Y90	ZR90	< ,					
!	^	^	^	^	^	^	^		
N. Pl	P237 PA U242 PU	243 CF2 231 PB2 244 RA2 34 U23	10 PM14 26 RA22	7 PU238 8 SR90	PU239	PU240	CS137 PU241 TH232		
TABULAR	_DATA		,						
! ! Example (of how r	adioisot	ope data	are ent	ered:				
! !1st :	Line: Ra	dionucli			in colum clide nam				
!2nd	& 3rd li	ne	by	radionuc		е, ел;	ACZZJ)		
!Field !Field !Field !Field	d#2 Half d#3 Acti	vity Con Release entory	version	(Kg/Mole (Years) (Ci/Kg) (Ci) (Ci) (Ci)	HALFY	[REAL [REAL [REAL]]]]	x,1pe14.6 " " ")))
! !*PU241									
!xxxaaaaaa !AWT !EPAREL !	2.41	eeeeeeee 0000E-01 00000E+07	HALFY	1.4	eeeeeeee 139900E+0 930000E+0	1 ACT	CNV 1	eeeeeeeee 030000E+).000000E+	05
! <table_inp< td=""><td>UTS</td><td></td><td></td><td></td><td></td><td></td><td></td><td>,</td><td></td></table_inp<>	UTS							,	
*AC225 AWT EPAREL		230E-01	HALFY INVCHI		37909E-02		NV 5.	802700E+0	7
*AC227 AWT EPAREL		280E-01	HALFY INVCHI		77335E+01 00000E+00		NV 7.	232300E+0	4
*AC228 AWT	2.280	310E-01	HALFY	6.99	93710E-04 00000E+00	ACTC	NV 2.	.241700E+0	9
EPAREL *AM241 AWT	2.410	0000E+00	INVCHI HALFY	4.32	22347E+02	ACTC	NV 3.	.431200E+0	3
EPAREL *AM243 AWT		0000E+02	INVCHI HALFY		00000E+05 80313E+03		NV 1.	.992900E+0	2
EPAREL *AT217	1.000	000E+02	INVCHI	0.0	00000E+00)		<i>:</i>	
AWT EPAREL *BA137M)050E-01)000E+00	HALFY INVCHI		23547E-09 00000E+00		NV 1.	.609500E+1	. ว
AWT EPAREL		070E-01	HALFY INVCHI		51550E-06 00000E+00		NV 5	.382400E+1	1
*BI211 AWT	2.109	9870E-01	HALFY	4.0	49824E-06	ACTC	INV 4	.184000E+1	1



EPAREL *BI212	0.00000E+00	INVCHD	0.00000E+00		
AWT	2.119910E-01	HALFY	1.151253E-04	ACTCNV	1.464800E+10
EPAREL	0.000000E+00	INVCHD	· 0.000000E+00		
*BI213					
AWT	2.129940E-01	HALFY	8.679553E-05	ACTCNV	1.933800E+10
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*BI214					
AWT	2.139990E-01	HALFY	3.783638E-05	ACTCNV	4.415300E+10
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
*CF252					•
AWT	2.520820E-01	HALFY	2.638090E+00	ACTCNV	5.375899E+05
EPAREL	0.00000E+00	INVCHD	2.960000E+03		
*CM243					
AWT	2.430610E-01	HALFY	2.850088E+01	ACTCNV	5.160700E+04
EPAREL	1.000000E+02	INVCHD	0.00000E+00		
*CM244					
AWT	2.440630E-01	HALFY	1.811013E+01	ACTCNV	8.088299E+04
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*CM245					
AWT	2.450650E-01	HALFY	8.498927E+03	ACTCNV	1.716500E+02
EPAREL	1.000000E+02	INVCHD	0.00000E+00		•
*CM248					
AWT	2.480720E-01	HALFY	3.390698E+05	ACTCNV	4.250200E+00
EPAREL	1.000000E+02	INVCHD	0.00000E+00		•
*CS137					
AWT	1.369070E-01	HALFY	2.999975E+01	ACTÓNV	8.704300E+04
EPAREL	1.000000E+03	INVCHD	2.070000E+04		
*FR221	•				
AWT	2.210140E-01	HALFY	9.126364E-06	ACTCNV	1.772400E+11
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*ND143					
AWT	1.430000E-01	HALFY	3.168876E+30	ACTCNV	0.00000E+00
EPAREL	1.000000E+03	INVCHD	0.000000E+00		
*NP237	· · · · · · · · · · · · ·				
AWT	2.370480E-01	HALFY	2.139942E+06	ACTCNV.	7.047600E-01
EPAREL	1.000000E+02	INVCHD	6.670000E+01		
*NP239					0.0101007.00
AWT	2.390530E-01	HALFY	6.448664E-03	ACTCNV	2.319100E+08
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*PA231	o 010070- 01		2.0766108.04	1000101	4 7335008.01
AWT	2.310360E-01	HALFY	3.276618E+04	ACTCNV	4.722500E+01
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*PA233	0 000000000	113 T 132	7.392988E-02	ACTCNV	2.075100E+07
AWT	2.330400E-01 0.000000E+00	HALFY INVCHD	0.000000E+00	ACICINV	2.0/51001.0/
EPAREL *PA234M	0.0000000000000000000000000000000000000	INVCHD	0.0000000000000000000000000000000000000		
AWT	2.340430E-01	HALFY	2.224551E-06	ACTCNV	6.866600E+11
EPAREL	0.000000E+00	INVCHD	0.000000E+00	nerenv	0.0000002.12
*PB209	0.0000000000000000000000000000000000000	Invenie	0.000002.00		
AWT	2.089810E-01	HALFY	3.764625E-04	ACTCNV	4.544100E+09
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*PB210	0.000002.00				
AWT	2.099840E-01	HALFY	2.229938E+01	ACTCNV	7.634800E+04
EPAREL	1.000000E+02	INVCHD	0.00000E+00		
*PB211					
AWT	2.109890E-01	HALFY	6.863786E-05	ACTCNV	2.468600E+10
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*PB212					
AWT	2.119920E-01	HALFY	1.213680E-03	ACTCNV	1.389500E+09
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*PB214					
AWT	2.140000E-01	HALFY	5.095553E-05	ACTCNV	3.278500È+10
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*PM147					
AWT	1.469150E-01	HALFY	2.623513E+00	ACTCNV	9.275300E+05
EPAREL	0.00000E+00	INVCHD	5.250000E+02		
*PO212					1 7720007:00
AWT	2.119890E-01	HALFY	9.506629E-15	ACTCNV	1.773900E+20
EPAREL	0.000000E+00	INVCHD	0.00000E+00		

*P0213	· · · · · · · · · · · · · · · · · · ·		1 222227 12		1 0611007.10
AWT	2.129930E-01 0.000000E+00	HALFY	1.330928E-13 0.000000E+00	ACTCNV	1.261100E+19
EPAREL * PO214	0.0000000000000000000000000000000000000	INVCHD	0.00000E+00		
AWT	2.139950E-01	HALFY	5.206464E-12	ACTCNV	3.208700E+17
EPAREL	0.000000E+00	INVCHD	0.000000E+00	ACICINV	5.2007001117
*P0215	0.000000100	111100112	0.0000002.00		
AWT	2.149990E-01	HALFY	5.640600E-11	ACTCNV	2.947900E+16
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*P0216					
AWT	2.160020E-01	HALFY	4.753314E-09	ACTCNV	3.482000E+14
EPAREL	0.00000E+00	INVCHD	0.000000E+00		
*PO218					
AWT	2.180090E-01	HALFY	5.799044E-06	ACTCNV	2.827800E+11
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*PU238					
AWT	2.380500E-01	HALFY	8.774619E+01	ACTCNV	1.711500E+04
EPAREL	1.000000E+02	INVCHD	4.240000E+06		
*PU239			2.406445E+04	3 CUC187	6.214600E+01
AWT	2.390520E-01 1.000000E+02	HALFY ' INVCHD	2.408445E+04 3.920000E+05	ACTCNV	0.214000E+01
EPAREL * PU240	1.00000000000	INVCHD	3.92000E+03		
AWT	2.400540E-01	HALFY	6.537392E+03	ACTCNV	2.278100E+02
EPAREL	1.000000E+02	INVCHD	6.930000E+04	11010111	2.2,01002,02
* PU241	1.0000001.01	11110112	0.000002/01		
AWT	2.410570E-01	HALFY	1.439937E+01	ACTCNV	1.030000E+05
EPAREL	0.000000E+00	INVCHD	1.930000E+06		
*PU242					3
AWT	2.420590E-01	HALFY	3.869198E+05	ACTCNV	3.817100E+00
EPAREL	1.000000E+02	INVCHD	4.910000E+04		· · ·
*PU244					
AWT	2.440640E-01	HALFY	8.261261E+07	ACTCNV	1.773100E-02
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*RA223	0 0001000 01		2 1205225 02	N OMONTZ	E 120600E+07
AWT	2.230190E-01	HALFY	3.130533E-02 0.000000E+00	ACTCNV	5.120600E+07
EPAREL *RA224	0.00000E+00	INVCHD	0.0000000000000		
AWT	2.240200E-01	HALFY	1.001999E-02	ACTCNV	1.592700E+08
EPAREL	0.000000E+00	INVCHD	0.000000E+00	ACTONV	1.352,002,00
*RA225	0.0000002.00	1111 0112	0.0000002.00		
AWT	2.250240E-01	HALFY	4.052993E-02	ACTCNV	3.919900E+07
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
*RA226					
AWT	2.260250E-01	HALFY	1.599966E+03	ACTCNV	9.885801E+02
EPAREL	1.000000E+02	INVCHD	0,00000E+00		
*RA228			C COODEEE.00	1.000101	2 240000 0
AWT	2.280310E-01	HALFY	6.699955E+00	ACTCNV	2.340000E+05
EPAREL *RN219	0.00000E+00	INVCHD	0.00000E+00		
AWT	2.190090E-01	HALFY	1.254875E-07	ACTCNV	1.300800E+13
EPAREL	0.000000E+00	INVCHD	0.000000E+00	ACTONV	1.500000115
*RN220	0.0000000000000000000000000000000000000	11110112	0.0000002,000		
AWT	2.200110E-01	HALFY	1.761895E-06	ACTCNV	9.222600E+11
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*RN222					
AWT	2.220180E-01	HALFY	1.046997E-02	ACTCNV	1.538000E+08
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*SM147			4 4544555		0.0054000.05
AWT	1.470000E-01	HALFY	1.059355E+11	ACTCNV	2.295400E-05
EPAREL	1.000000E+02	INVCHD	0.00000E+00		
*SR90	8.990801E-02	HALFY	2.912197E+01	ACTCNV	1.365400E+05
AWT EPAREL	1.000000E+03	INVCHD	9.850002E+01		T.202-2001.00
*TH227	T.00000000403	TIANCIID	2.00000020100		
AWT	2.270280E-01	HALFY	5.124073E-02	ACTCNV	3.073200E+07
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*TH228					
AWT	2.280290E-01	HALFY	1.913051E+00	ACTCNV	8.195301E+05
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*TH229					

			······			
AWT EPAREL	2.290320E-01 1.000000E+02	HALFY INVCHD	7.339118E+03 0.000000E+00	ACŤCNV	2.126900E+02	
*TH230						
AWT	2.300330E-01	HALFY	7.700370E+04	ACTCNV	2.018300E+01	
EPAREL	1.000000E+01	INVCHD	0.000000E+00			
*TH231						
AWT	2.310360E-01	HALFY	2.911247E-03	ACTCNV	5.315200E+08	i.
EPAREL	0.000000E+00	INVCHD	0.00000E+00			•.
*TH232						
AWT	2.320380E-01	HALFY	1.405080E+10	ACTĊNV	1.096500E-04	
EPAREL	1.000000E+01	INVCHD	6.110000E-01			
*TH234						
AWT	2.340440E-01	HALFY	6.597601E-02	ACTCNV	2.315200E+07	
EPAREL	0.00000E+00	INVCHD	0.00000E+00			
*TL207					1 0045005.11	
AWT	2.069770E-01	HALFY	9.069325E-06	ACTCNV	1.904500E+11	
EPAREL	0.000000E+00	INVCHD	0.000000E+00			
*U233			1 5050305 05		0. (202005:00	
AWT	2.330400E-01	HALFY	1.585072E+05	ACTCNV	9.678300E+00	
EPAREL	1.000000E+02	INVCHD	1.310000E+03			
*U234	0 0404105 01	175 T 137	2 445105D.05		6.247300E+00	
AWT	2.340410E-01	HALFY	2.445105E+05	ACTCNV	0.24/300E+00	
EPAREL *U235	1.000000E+02	INVCHD	0.00000E+00			
	2.350440E-01	HALFY	7.038074E+08	ACTCNV	2.161100E-03	
AWT		INVCHD	1.150000E+00	ACICINV	2.101100E-05	
EPAREL	1.000000E+02	INVCHD	1.1200005+00			
*U236 AWT	2.360460E-01	HALFY	2.341483E+07	ACTCNV	6.468300E-02	
EPAREL	1.000000E+02	INVCHD	0.000000E+00	ACICINV	0.4005004-02	
*U238	1.0000000000000	INVCHD	0.0000000000000000000000000000000000000			
AWT	2.380510E-01	HALFY	4.468115E+09	ACTCNV	3.361100E-04	
EPAREL	1.000000E+02	INVCHD	2.010000E-01	nerenv	5.5011002 01	
*Y90	1.000000102	Invenb	2.0100001 01			
AWT	8.990701E-02	HALFY	7.301091E-03	ACTCNV	5.446300E+08	
EPAREL	0.000000E+00	INVCHD	0.000000E+00			
*ZR90	0.0000001.00	111100112	01000002.00			
AWT	9.000000E-02	HALFY	3.168876E+30	ACTCNV	0.00000E+00	
EPAREL	1.000000E+03	INVCHD	0.000000E+00			
END_TABLES>						
!						
! END OF RAI	DIOISOTOPE_INPUT	ſ				· ·
!						
*********	*************	*******	*****	*******	* * * * * * * * * * *	
(SCATXT1)						
Explanatory	Comment: This	ends the	listing of data	read from	the repository/mod	lel
input contro	ol file and beg	ins the ec	ho of drilling/.	intrusion	data read from the	(input
control) fil	le whose logical	l symbol i	s CUSP_INP\$TXT1	•	End of Comment.	
1						
1	•					
! INTRUSION	N					
TINTR	100.0					
	BOREHOLE: TAUFAI	•				
	Sonenous:IAOFALI	J IFATI. valu	e extracted from	m CAMDAT *	**	
	NATEDIAL DOAD		C = BOREHOLE: TAU	FAIL · P	11200E+00)	
(CARDA)					· · · · · = • • • • • • • • • • • • • •	
	I MAIENIAD. PROPI					
TCLOUT V						•
	VAS_AREA:TCLOUT	.OUT(*) va	lue extracted f		***	•
*** CUTTING	VAS_AREA:TCLOUT SS variable: TCI	LOUT(*) va ERTY:VALUE	lue extracted f: = WAS AREA:TCL	rom CAMDAI	*** .48000E+04)	•
*** CUTTING	VAS_AREA:TCLOUT SS variable: TCI	LOUT (*) va ERTY : VALUE	lue extracted f: 2 = WAS_AREA:TCL	rom CAMDAI	.48000E+04)	•
*** CUTTING (CAMDA)	VAS_AREA:TCLOUT SS variable: TCI MATERIAL:PROPI	LOUT(*) va ERTY:VALUE	lue extracted f: 2 = WAS_AREA:TCL	rom CAMDAI	.48000E+04)	•
*** CUTTING	VAS_AREA:TCLOUT SS variable: TCI MATERIAL:PROPI	LOUT(*) va ERTY:VALUE	lue extracted f: 2 = WAS_AREA:TCL	rom CAMDAI	.48000E+04)	•
*** CUTTING (CAMDA)	WAS_AREA:TCLOUT SS variable: TCI MATERIAL:PROPI 0.006	LOUT(*) va ERTY:VALUE	lue extracted f: 2 = WAS_AREA:TCL	rom CAMDAI	*** .48000E+04)	
*** CUTTING (CAMDA) PARTDIA	WAS_AREA:TCLOUT SS variable: TCI MATERIAL:PROPI 0.006	LOUT (*) va ERTY: VALUE	lue extracted f: 2 = WAS_AREA:TCL	rom CAMDAI	*** .48000E+04)	•
*** CUTTING (CAMDAT PARTDIA ! PROPERTIES DNSFLUID I	VAS_AREA:TCLOUT SS variable: TCI MATERIAL:PROPI 0.006 DRILLMUD:DENSITY	ERTY : VALUE	: = WAS_AREA:TCL	rom CAMDAI OUT : é	.48000E+04)	•
*** CUTTING (CAMDAT PARTDIA ! PROPERTIES DNSFLUID I *** CUTTING	VAS_AREA:TCLOUT SS variable: TCI T MATERIAL:PROPI 0.006 DRILLMUD:DENSIT SS variable: DEI	ERTY:VALUE Y NSITY valu	: = WAS_AREA:TCL	rom CAMDAI OUT : é m CAMDAT *	***	
*** CUTTING (CAMDAT PARTDIA ! PROPERTIES DNSFLUID I *** CUTTING	VAS_AREA:TCLOUT SS variable: TCI T MATERIAL:PROPI 0.006 DRILLMUD:DENSIT SS variable: DEI	ERTY:VALUE Y NSITY valu	: = WAS_AREA:TCL	rom CAMDAI OUT : é m CAMDAT *	***	

DOMEGA BOREHOLE : DOMEGA *** CUTTINGS variable: DOMEGA value extracted from CAMDAT *** (CAMDAT MATERIAL: PROPERTY: VALUE = BOREHOLE: DOMEGA : 7.80000E+00) DRILLMUD:VISCO VISCO *** CUTTINGS variable: VISCO value extracted from CAMDAT *** (CAMDAT MATERIAL: PROPERTY: VALUE = DRILLMUD: VISCO : 9.17000E-03) YLDSTRSS DRILLMUD: YLDSTRSS *** CUTTINGS variable: YLDSTRSS value extracted from CAMDAT *** (CAMDAT MATERIAL: PROPERTY: VALUE = DRILLMUD: YLDSTRSS : 4.40000E+00) ABSROUGH WAS_AREA:ROUGH *** CUTTINGS variable: ABSRO value extracted from CAMDAT *** (CAMDAT MATERIAL: PROPERTY: VALUE = WAS_AREA: ROUGH : 2.50000E-02) ! ! BRAGFLOW ! Multiple hits (max of 10, 0 thru 9) must be entered in order is associated with the hit that CUTTINGS used for BRAG flow ! MHIT O properities I INTR_0 CAVITY_1 < CAVITY_1 NO_GRDVOL < . INTR_1 SIMPLE AVERAGE WILL BE USED INSTEAD OF WEIGHTED AVERAGE VOLUME . INTR_2 457 458 459 < INTR_3 460 461 462 < INTR_4 463 464 465 <

Explanatory Comment: The following is a summary of the data CUTTINGS_S read from the BRAGFLO output CDB file. Note that block number 7, which contains 21 elements, is the location of the repository. End of Comment.

element block names found

block name, S_HALITE DRZ_0 TRANS_0 S_MB139 S_ANH_AB S_MB138 CAVITY_1 CAVITY_2 CAVITY_2 CAVITY_2 CAVITY_3 CAVITY_4 IMPERM_2 CASTILER WAS_AREA REPOSIT DRZ_1 UNNAMED CULEBRA TAMARISK MAGENTA FORTYNIN DEWYLAKE SANTAROS BACKFILL	1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23	id

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SHFT_B_2	26		
SHFT_L_1	27		
SHFT_L_2	28 29		
SHFT_U_1 SHFT_U_2	29 30		
SHFT_LS1	31		
SHFT LS2	32		
SHFT_US1	33		
SHFT_US2	34		
PAN_S_1	35		
PAN_S_2	36	•	
BOREHOLE	37		
BRINESAL	38		
H2	39		
PB210	40		
RA226	41		
RA228	42 43		
TH229 TH230	43		
TH230 TH232	44		
PA231	46		
U233	47		
U234	48		
U235	49		
U236	50		
U238	51		
NP237	52		
PU238	53 54		
PU239 PU240	54		
PU240 PU241	56		
PU242	57		
PU244	58		
AM241	59		
CM244	60		
CM248	61		
CF252	62		
PB	63 64		
RA TH	65		
PA	66		
Ū	67		
NP	68		
PU	69		
AM	70		
CM	71		
CF	72		
TRANS_1	73		

number of elements per block

	••••••		
block 1	number, number 297	of	elements
	64		
2 3	• -		
3	16		
4	14		
5	14		
4 5 6	30		
7	21		
8	9		
9	12		
10	27		
11	275		
12	27		
13	0		
14	0		
15	0		
16	0		
17	0		

18	0			
18 19 20 21 22 23 24 25 26 27 28 29 30 31 32 33 34 35 36 37 38 40 41 42 43 44 45 46 47 48 49 50 51 52 53 54 55 56 57 58 59 60 61 62 63 64 65 66 65 66				
21 22	0			
23	0			
24	0			
26 27	0			
28	0			
29 30	0 0			
31	0			
32 33	0			
34	0			
36	0			
37 38	0 0			
39	0			
41	0			
42 43	0 0			
44	0			
45	0			
47 48	0 0			
49	0			
50	0			
52 53	0 0			
54	0			
55 56	0			×
57 58	0			
59	0			
61	0			
62 63	0			
64	0			
66	0			
67	0			
68 69 70	0			
71	0			
72 73	0 0 0 0 0			
	-			

attribute names found

i,	name
1	THICK
2	ELEVAT
3	DEL_X
4	DEL_Y
5	DEL_Z
6	GRIDVOL

element #	attribute value 1.2799E+11	
23	1.4830E+10 1.9030E+09	
4 5	2.3061E+08 2.9423E+07	
6 7 8	1.0210E+07 2.4135E+06	
8 9 10	1.8411E+06 9.8849E+06 1.9705E+06	
11 12	5.5664E+04 7.6922E+05	
13 14	4.0078E+04 1336.	
15 16 17	10.96 1336. 4.0078E+04	
18 19	7.6922E+05 2.1041E+05	
20 21	7.4433E+05 1.3151E+05	
22 23 24	1.3220E+04 1.7117E+05 3.6947E+06	
25 25	1.7519E+06 8.1208E+06	
27 28	2.8095E+07 2.7331E+08	
29 30 31	2.0035E+09 1.5118E+10 1.2885E+11	
32 33	4.5986E+10 5.3286E+09	
34 35	6.8375E+08 8.2858E+07	
36 37 38	1.0572E+07 3.6686E+06 8.6715E+05	
39 40	6.6150E+05 3.5516E+06	
41 42	7.0800E+05 2.0000E+04	
43 44 45	2.7638E+05 1.4400E+04 480.0	
46 47	3.936 480.0	
48 49	1.4400E+04 2.7638E+05	
50 51 52	7.5600E+04 2.6744E+05 4.7250E+04	
53 54	4750. 6.1500E+04	
55 56 57	1.3275E+06 6.2945E+05 2.9178E+06	
57 58 59	1.0095E+07 9.8200E+07	
60 61	7.1985E+08 5.4317E+09	
62 63 64	4.6295E+10 1.0117E+10 1.1723E+09	
65 66	1.5043E+08 1.8229E+07	
67 68	2.3257E+06 8.0709E+05	
69	1.9077E+05	

$\begin{array}{c} 70\\ 71\\ 72\\ 73\\ 74\\ 75\\ 76\\ 77\\ 78\\ 79\\ 80\\ 81\\ 82\\ 83\\ 84\\ 85\\ 86\\ 87\\ 88\\ 89\\ 90\\ 91\\ 92\\ 93\\ 94\\ 95\\ 96\\ 97\\ 98\\ 99\\ 100\\ 101\\ 102\\ 103\\ 104\\ 105\\ 106\\ 107\\ 108\\ 109\\ 110\\ 111\\ 112\\ 113\\ 114\\ 115\\ 116\\ 117\\ 118\\ 119\\ 120\\ 121\\ 122\\ 123\\ 124\\ 125 \end{array}$	1.4553E+05 7.8136E+05 4400. 6.0803E+04 3168. 105.6 0.8660 105.6 3168. 6.0803E+04 1.6632E+04 5.8836E+04 1.0395E+04 1.045. 1.3530E+04 2.9205E+05 1.3848E+05 6.4192E+05 2.2208E+06 2.1604E+07 1.5837E+08 1.1950E+09 1.0185E+10 1.2692E+09 1.4707E+08 1.8872E+07 2.2869E+06 2.9177E+05 1.0125E+05 2.3933E+04 1.2148E+09 1.4076E+08 1.8062E+07 2.1888E+06 2.7926E+05 9.6910E+04 2.2907E+04 1.2148E+09 1.4076E+08 1.8062E+07 2.1888E+06 2.7926E+05 9.6910E+04 2.2907E+04 1.2148E+09 1.4076E+08 1.8062E+07 2.1888E+06 2.7926E+05 9.6910E+04 2.2907E+04 1.2148E+09 1.4076E+08 1.8062E+07 2.1888E+06 2.7926E+05 9.6910E+04 2.2907E+04 1.2148E+09 1.4076E+08 1.8062E+07 2.1888E+06 2.7926E+05 9.6910E+04 2.2907E+04 1.2148E+09 1.4076E+08 1.8062E+07 2.1888E+06 2.7926E+05 9.6910E+04 2.2907E+04 2.2907E+04 3.5796E+07 4.3378E+06
116	1.4076E+08
117	1.8062E+07
118	2.1888E+06
119	2.7926E+05
120	9.6910E+04
121	2.2907E+04
122	2.4074E+09
123	2.7896E+08
124	3.5796E+07

171 $1.1406E+05$ 172 $5.2871E+05$ 173 $1.8291E+06$ 174 $1.7794E+07$ 175 $1.3044E+08$ 176 $9.8423E+08$ 177 $8.3887E+09$ 178 $5.6084E+10$ 179 $6.4987E+09$ 180 $8.3390E+08$ 181 $1.0105E+08$ 182 $1.2893E+07$ 183 $4.4742E+06$ 184 $1.0576E+06$ 185 $8.0677E+05$ 186 $4.3316E+06$ 187 $8.6348E+05$ 188 $2.4392E+04$ 189 $3.3707E+05$ 190 $1.7562E+04$ 191 585.4 194 $1.7562E+04$ 195 $3.3707E+05$	140 141 142 143 144 145 146 147 148 149 150 151 152 153 154 155 156 157 158 159 160 161 162 163 164 165 166 167 168 169 170	1.9016E+07 1.4348E+08 1.2229E+09 1.6628E+04 7.7077E+04 2.6666E+05 2.5941E+06 1.9016E+07 1.4348E+08 1.2229E+09 1.6628E+04 7.7077E+04 2.6666E+05 2.5941E+06 1.9016E+07 1.4348E+08 1.2229E+09 3.2953E+04 1.5275E+05 5.1410E+06 3.7686E+07 2.8436E+08 2.4236E+09 8.3326E+09 9.6554E+08 1.2390E+08 1.2390E+08 1.5014E+07 1.9156E+06 6.6475E+05 1.5713E+05
186 4.3316E+06 187 8.6348E+05 188 2.4392E+04 189 3.3707E+05 190 1.7562E+04 191 585.4 192 4.801 193 585.4 194 1.7562E+04	171 172 173 174 175 176 177 178 179 180 181 182 183 184	1.1406E+05 5.2871E+05 1.8291E+06 1.7794E+07 1.3044E+08 9.8423E+08 8.3887E+09 5.6084E+10 6.4987E+09 8.3390E+08 1.0105E+08 1.2893E+07 4.4742E+06 1.0576E+06
	186 187 188 189 190 191 192 193 194	4.3316E+06 8.6348E+05 2.4392E+04 3.3707E+05 1.7562E+04 585.4 4.801 585.4 1.7562E+04

210 211 212 213 214 215 216 217 218 219 220 221 222 223 224 225 226 227 228 229 230 231 232 233	8.7628E+05 4.5657E+04 1522. 12.48 1522. 4.5657E+04 8.7628E+05 2.3970E+05 8.4793E+05 1.4981E+05 1.4981E+05 1.4580E+11 1.6895E+10 2.1679E+09 2.6271E+08 3.3518E+07 1.1632E+07 2.7494E+06 6.3412E+04 8.7628E+05 4.5657E+04 1522.
234 235 236 237 238 239 240 241 242 243 244 245 244 245 246 247 248 249 250 251 252 253 255 255 256 257	12.48 1522. 4.5657E+04 8.7628E+05 2.3970E+05 8.4793E+05 1.4981E+05 4.6721E+09 5.4138E+08 6.9469E+07 8.4183E+06 1.0741E+06 3.7273E+05 8.8102E+04 4.6085E+05 7.1933E+04 2032. 2.8080E+04 1463.
258 259 260 261 262 263 264 265 266 267 268 269 270 271 272 273 274 275 274 275 276 277 278 279	2.8080E+04 7681. 2.7171E+04 4801. 7.5005E+04 1.6190E+06 7.6768E+05 3.5585E+06 1.2311E+07 1.1976E+08 8.7793E+08 6.6245E+09 5.6462E+10 1.9499E+05 4.2090E+06 1.9957E+06 9.2512E+06 3.2006E+07 3.1135E+08 2.2824E+09 1.7222E+10 1.4678E+11

280 281 282 283 284 285 286 287 288 289 290 291 292 293 294 295 296 297 298 299 300 301 302 304	1.9499E+05 4.2090E+06 1.9957E+06 9.2512E+06 3.2006E+07 3.1135E+08 2.2824E+09 1.7222E+10 1.4678E+11 6248. 1.3487E+05 6.3952E+04 2.9645E+05 1.0256E+04 2.9645E+05 1.0256E+06 9.9771E+06 7.3137E+07 5.5186E+08 4.7036E+09 1.1246E+04 6.0378E+04 1.2036E+04 340.0 4698. 244.8
304 305 306 307 308 309 310 311 312 313 314 315 316 317 318 319 320 321 322 323 324 325 326 327 328 329	8.160 6.6916E-02 8.160 244.8 4698. 1285. 4546. 803.3 1.8257E+04 9.8025E+04 1.9541E+04 552.0 7628. 397.4 13.25 0.1086 13.25 397.4 7628. 2087. 7381. 1304. 1046. 2.2568E+04 1697. 3.6639E+04
330 331 332 333 334 335 336 337 338 339 340 341 342 343 344 345 344 345 346 347 348 349	3.4631E+04 1.8593E+05 3.7065E+04 1047. 1.4469E+04 753.9 25.13 0.2061 25.13 753.9 1.4469E+04 3958. 1.4001E+04 2474. 3572. 1.9179E+04 3823. 108.0 1492. 77.76

3971.9810E+043984683.3993399.4001.5756E+044015.4510E+04
200 4002

420 421 422 423 424 425 426 427 428 429 430 431 432 433 434 435 435 436 437 438 439 441 442 444 445 445 447 448 445 445 451 455 455 455 457 458 459	$\begin{array}{c} 1.4170E-02\\ 1.728\\ 51.84\\ 995.0\\ 272.2\\ 962.8\\ 170.1\\ 221.4\\ 4779.\\ 2266.\\ 1.0504E+04\\ 3.6340E+04\\ 3.5352E+05\\ 2.5915E+06\\ 1.9554E+07\\ 1.6666E+08\\ 7301.\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 380.4\\ 12.68\\ 380.4\\ 12.68\\ 0.1040\\ 12.68\\ 0.1040\\ 0.1040\\ 0.1040\\ 0.1040\\ 0.1040\\ 0.104$
460 461 462 463 464 465 466 465 467 468 469 470 471 472 473 477 477 477 477 477 477 477 477 477 477 477 477 477 477 478 480 481 482 483 484 485 484 485 484 485 488 488 488 489	1.7474E+04 9.3820E+04 1.8703E+04 1.7474E+04 9.3820E+04 1.8703E+04 7065. 1248. 7065. 1248. 7065. 1248. 7065. 1248. 3.5067E+04 1625. 3.5067E+04 1528.3 528.3 1997. 1997. 80.75 131.1 125.5 125.5 248.7

560 561 562 563 565 565 566 567 568 569 570 571 572 571 572 573 574 575 576 577 578 579 579 580	1.5546E+06 1.5123E+07 1.1086E+08 8.3649E+08 7.1295E+09 2.2809E+10 2.6430E+09 3.3914E+08 4.1097E+07 5.2435E+06 1.8196E+06 4.3011E+05 3.2810E+05 1.7616E+06 3.5117E+05 9920. 1.3708E+05 7142. 238.1 1.952 238.1
581 582 583 584 585 586 587 588 589 590 591 593 593 593 593 593 595 594 595 597 598 590 597 598 590 601 602 603 604 605 607 608 609	7142. 1.3708E+05 3.7498E+04 1.3265E+05 2.3436E+04 3.0504E+04 6.5844E+05 3.1221E+05 1.4472E+06 5.0069E+06 4.8707E+07 3.5705E+08 2.6941E+09 2.2962E+10 7.8176E+09 9.0586E+08 1.1624E+08 1.4086E+07 1.7972E+06 6.2366E+05 1.4742E+05 1.1246E+05 1.2366E+05 1.4742E+05 1.2366E+05 1.2366E+05 3400. 4.6984E+04 2448. 81.60 0.6692
610 611 612 613 614 615 616 617 618 619 620 621 622 623 624 625 624 625 626 627 628 629	81.60 2448. 4.6984E+04 1.2852E+04 4.5464E+04 8033. 1.0455E+04 2.2568E+05 1.0701E+05 4.9603E+05 1.7161E+06 1.6694E+07 1.2238E+08 9.2339E+08 7.8702E+09 1.5911E+10 1.8437E+09 2.3658E+08 2.8669E+07 3.6577E+06

630	1.2693E+06
631	3.0003E+05
632	2.2888E+05
633	1.2289E+06
634	2.4497E+05
635	6920.
636	9.5627E+04
637	4982.
638	166.1
638 640 641	166.1 1.362 166.1 4982.
642	9.5627E+04
643	2.6158E+04
644	9.2533E+04
645	1.6349E+04
646	2.1279E+04
647	4.5932E+05
648	2.1779E+05
649	1.0096E+06
650	3.4927E+06
651	3.3977E+07
652	2.4907E+08
653	1.8794E+09
654	1.6018E+10
655	9.7490E+10
656	1.1297E+10
657	1.4496E+09
658	1.7566E+08
659	2.2412E+07
660	7.7774E+06
661	1.8384E+06
662	1.4024E+06
663	7.5294E+06
664	1.5010E+06
665	4.2400E+04
666	5.8592E+05
667	3.0528E+04
668	1018.
669	8.345
670	1018.
671	3.0528E+04
672	5.8592E+05
673	1.6027E+05
674	5.6696E+05
675	1.0017E+05
676	3.9824E+10
677	4.6145E+09
678	5.9213E+08
679	7.1755E+07
680	9.1549E+06
681	3.1770E+06
682	7.5095E+05
683	5.7286E+05
684	3.0757E+06
685	6.1313E+05
686	1.7320E+04
687	2.3934E+05
688	1.2470E+04
689	415.7
690	3.409
691	415.7
692 693 694	1.2470E+04 2.3934E+05 6.5470E+04 2.3160È+05
695	2.3160E+05
696	4.0919E+04
697	1.3038E+05
698	2.8143E+06
699	1.3344E+06

700 701 702 703 704 705 706 707 708 709 710 7112 713 714 715 716 717 718 719 720 721 722 723 724 725 726 727 728 729 730 731 732 733 734 735 737 738 739 741 745 746 747 748 745 755 753 755	6.1857E+06 2.1400E+07 2.0818E+08 1.5261E+09 1.515E+10 9.8146E+10 5.3259E+04 1.1496E+06 5.4510E+05 2.5268E+06 8.7418E+06 8.7418E+07 6.2339E+08 4.7039E+09 4.0092E+10 1.4495E+10 1.6796E+09 2.1552E+08 2.6117E+07 3.321E+06 1.1563E+06 2.7333E+05 2.0850E+05 1.1195E+06 2.2316E+05 6304. 8.7114E+04 4539. 151.3 1.241 151.3 4539. 8.7114E+04 2.3829E+04 8.4296E+04 1.4893E+04 1.9840E+05 9.1969E+05 3.1818E+06 3.0953E+07 2.2690E+08 1.7121E+09 1.4592E+10 2.5090E+09 2.9635E+08 2.5259E+09 4.3569E+10 5.0485E+08 2.5259E+09 4.3569E+10 5.0485E+09 6.4781E+08 7.8503E+07 3.4758E+06 8.2157E+05
746 747 748 750 751 752 753 754 755 756 757 758 759 760 761 762 763 764	2.9073E+08 2.9635E+08 2.5259E+09 4.3569E+10 5.0485E+09 6.4781E+08 7.8503E+07 1.0016E+07 3.4758E+06 8.2157E+05 6.2673E+05 3.3650E+06 6.7079E+05 1.8949E+04 2.6185E+05 1.3643E+04 454.8 3.729 454.8
765 766 767 768 769	1.3643E+04 2.6185E+05 7.1626E+04 2.5338E+05 4.4767E+04

	7712345677890123445678890123456789012333333333333333333333333333333333333	5.8 1.2 5.9 9.5 9.5 9.5 9.5 9.5 1.5 1.5 7.8 2.6 1.5 7.8 2.6 1.5 7.8 2.6 1.5 7.8 2.6 1.5 7.8 2.6 7.8 2.6 7.8 2.6 7.8 2.6 7.8 2.6 7.8 2.6 7.8 2.6 7.8 7.8 7.8 7.8 7.8 7.8 7.8 7.8	2577764992775639278 25777649927756392078 257776499278 257776499278 257776499278 2577775078 257776499278 2577789 2577789 2577789 2577789 25789 25789	32 ± 04 72 ± 06 72 ± 06 92 ± 06 92 ± 07 22 ± 08 22 ± 09 22 ± 10 52 ± 07 72 ± 06 32 ± 05 52 ± 04 32 ± 04 32 ± 04 32 ± 04 7 93 92 ± 10 52 ± 04 32 ± 04 7 93 ± 04 7 93 ± 04 7 93 ± 04 7 93 ± 04 7 93 ± 04 12 ± 04 32 ± 0	
number	of	elements	to	average	over
number	of	elements	to	average	over
number	of	elements	to	average	over
number	of	elements	to	average	over

unique property names

number of element to average over

Information Only

21

21

3

3

3

1

property	property				
number	name				
1					
<u> </u>	PRMX_LOG				
2	POROSITY				
3	SP_S_LOG				
4	RELP_MOD				
5 6	CAP_MOD				
6	PC_MAX				
2					
7	SAT_RBRN				
8	SAT_RGAS				
9	PORE_DIS				•
10	COMP_RCK				
11	SB_MIN				
12	PO_MIN				
13	PCT_A				
14	PCT_EXP			:	,
15	PCT_FLAG				
16	PTINDEX				
17	GRID_ID				
18	REF_PRES				
19	PTHRESH				
20	PRESSURE				,
21	PERM_EXP				
22	PHIMAX			•	
23	DPHIMAX			•	A •
24	PI DELTA				
25	PF_DELTA				
26					
	P_LITHO				• •
27	FRAC_PMX				
28	FRAC_PMY				
29	FRAC_PMZ				
30	KMAXLOG				
31	DELTA_PF				
32					
32	CG_NF				
33	CG_F				
34	CB				
35	SAT_IBRN				
36	SAT_PUD				
37	GRATCORI				
38	GRATCORH				
39	GRATMICI				
40	GRATMICH				
41	STOICOR				
42	STOIMIC		,		
43	DCELLCHW				
44	DCELLRHW				
45	DIRONCHW			,	
46	DIRONRHW				
47	DPLASCHW				•
48	DPLASRHW				•
49	DRUBBCHW				
50	DRUBBRHW				
51	DIRNCCHW				
52	DIRNCRHW				
53	DPLSCCHW				
54	VOLCHW				
54					
55	VOLRHW				
56	DRH_METL				
57	DCH_METL				
58	DRH_BIO				
5 9	DCH_BIO				
60	FRAC_PR				
C1					
61	PRMX				
62	STORAGEC				
63	DIAMMOD				
64	DNSFLUID				
65	WTF				
66	COMPRES				
00					
67	RSOD0				
68	RSOD1				

69	BSLOPE
70	RSLOPE
71	VSLOPE
72	VISCO
73	REF_TEMP
74	INVCHD
75	INVRHD
76	HALFLIFE
77	ATWEIGHT
78	DIS_ENGY
79	LOGSOLM
80	PERM_X
81	PERM_Y
82	PERM_Z
83	SOLUB
84	PAVE
85	COMP

Explanatory Comment: The data below represent CUTTINGS_S's search (through BRAGFLO's output CDB file) for the time of intrusion. End of Comment.

element blog TIME-years	ck 37 3.16599	9995E-01
TIME-years	5.34747819E-08	
TIME-years	2.22228164E-06	
TIME-years	1.09104358E-05	
TIME-years	7.41431795E-05	
TIME-years	2.42588285E-04	
TIME-years	2.23672390E-03	
TIME-years	2.98848525E-02	
TIME-years	6.96178913E-01 3.18654060E+00	
TIME-years TIME-years	1.78978786E+01	
TIME-years	2.58163357E+01	
TIME-years	3.57182579E+01	
TIME-years	5.97380142E+01	
TIME-years	6.10059128E+01	
TIME-years	6.26197319E+01	
TIME-years	7.07265167E+01	
TIME-years	7.27253571E+01	
TIME-years	7.28040161E+01	
TIME-years	7.83843460E+01 8.09984283E+01	
TIME-years TIME-years	8.67234039E+01	
TIME-years	1.00000221E+02	
IIII Jouro		

Interpolation xval 9.99983311E-01 for TINT(KINT) 1.0000000E+02

Explanatory Comment: The listing below represents all the data CUTTINGS_S extracted from the BRAGFLO output CDB file. End of Comment.

.

RESULTS from BRAGFLOW file: CUSP_INP\$BRAGCDB

Unaltered output follows here:

Diameter of borehole 3.16599995E-01

Initial porosity 8.81034851E-01

Intrusion	element	average
time (yrs)	properity	value

Multiple hits_0 0.00 POROS 8.81034851E-01

Multiple hits_1	PRESGAS PRESBRIN SATGAS	6.80242750E+06 1.01325008E+05 9.95905936E-01
0.00	POROS PRESGAS PRESBRIN SATGAS	8.81034970E-01 6.80243050E+06 1.01325000E+05 9.95906115E-01
Multiple hits_2 0.00	POROS PRESGAS PRESBRIN SATGAS	8.81034911E-01 6.80242850E+06 1.01324992E+05 9.95905936E-01
Multiple hits_3 0.00	POROS PRESGAS PRESBRIN SATGAS	8.81034911E-01 6.80242850E+06 1.01324992E+05 9.95905936E-01
Multiple hits_4 0.00	POROS PRESGAS PRESBRIN SATGAS	8.81034911E-01 6.80242850E+06 1.01324992E+05 9.95905936E-01
Multiple hits_0 100.00	POROS PRESGAS PRESBRIN SATGAS	3.00751239E-01 1.65183310E+07 9.81722700E+06 9.83161747E-01
Multiple hits_1 100.00	POROS PRESGAS PRESBRIN SATGAS	3.00751239E-01 1.65183330E+07 9.81722900E+06 9.83178914E-01
Multiple hits_2 100.00	POROS PRESGAS PRESBRIN SATGAS	3.00751209E-01 1.65185040E+07 9.81740000E+06 9.77457225E-01
Multiple hits_3 100.00	POROS PRESGAS PRESBRIN SATGAS	3.00751209E-01 1.65183470E+07 9.81724400E+06 9.89357471E-01
Multiple hits_4 100.00	POROS PRESGAS PRESBRIN SATGAS	3.00751209E-01 1.65181900E+07 9.81708600E+06 9.82479930E-01

From these two relationships:

Code will calculate porosity as: Pf=Pb/(1.0-Pi+Pb) Hf=(1.0-Pi)/(1.0-Pf)*Hi

& output these transformed values to the output cuttings CDB

after max, min

repo pres (pa) 1.5000E+07 permeability 1.0000E-12 ***** (DRILL) Explanatory Comment: The data below are a summary of the inputs and results of CUTTINGS_S's drilling calculations. End of Comment. 7.8000E+00 DOMEGA = ROUTER = 1.5830E-011.0160E-01 RINNER = DIADRIL = 3.1660E-01 UPHOLEV = 6.7942E-011.8340E-02 ETAO = SIGMA1 = SIGMA2 = 1.0821E-06 5.4103E-07 DENSITY = 1.2100E+03 8.1120E+00 TAUFAIL = ABSOLUTE BOREHOLE ROUGHNESS = 2.5000E-02PLASTIC VISCOSITY = 9.1700004E-03 INTRUSION NUMBER (OR HIT) = 1 INTRUSION MINISTRESS = 1.4000001E+00 1.000000E+02 INTRUSION_TIME (UNITS=Years) = FOR REYNOLDS # \leq 2100 ASSUMES LAMINAR HELICAL FLOW FOR REYNOLDS # >2100 ASSUMES TURBULENT AXIAL FLOW INITIAL REYNOLDS NUMBER = 8301., FLOW IS TURBULENT. CRITICAL RADIUS = 9.2572E-01 VBAR MULTIPLIER (CC) = 1.4987E+00 REYNOLDS INTRUSION COLLAR DRILL FINAL CRITICAL VBAR FINAL SOLUTION RADIUS RADIUS RADIUS MULT. AREA NUMBER TTME RADTUS ALGORITHM _____ ____ -- --------1.0000E+02 1.0160E-01 1.5830E-01 1.8928E-01 9.2572E-01 1.4987E+00 1.1255E-01 7.3961E+03 TURBULANT-AXIAL ******** WIPP TRU CONTACT HANDLED WASTE ******* PG PARTDIA AREA VOLUME 1.5000E+07 6.0000E-03 1.206 2.005 VOLUME_S <SPALL> ~~~~ 0.5684 >*<>*<>*<>*<>*<>*<>*<>*<>*<>*<>*<> \!/\!/\!/\!/ HIT NUMBER 1 >*<>*<>* *<>*<>*<> /^\/^\/^\/^\ /^\/^\/^\/^\ >*<>*<>*<>*<>*<>*<>*<>*<>*<>*<> diameter (m) 0.3166 intrusion time (yrs) 100.0 1.5000E+07 repository pre (pa)

Information Only

CUTTINGS_S, Version: 5.10

User's Manual

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porosity, initial (unitless) 0.8810 porosity (bragflow)
porosity (unitless)
permeability (m**2) 0.3008 0.7166 1.0000E-12 6.0000E-03 particle dia (m) FINAL RESULTS: Type: SOLID_BLOWOUT AREA (spall) 1.206 0.1125 AREA (erosion) AREA (total) 1.319 (WRIOUT)

Explanatory Comment: The data below are the results of dividing the initial inventory (CMS-controlled data) by INV.AREA, whose numerical value is 1.11520e5 m². End of Comment.

INITIAL radioisotope LOADING (CI) with an inventory
area of 1.11520E+05 M**2 FOR WIPP TRU CH WASTE)

AM241	2.4000E+05
AM243	0.0000E+00
CF252	2.9600E+03
CM243	0.0000E+00
CM244	0.0000E+00
CM245	0.0000E+00
CM248	0.0000E+00
CS137	2.0700E+04
NP237	6.6700E+01
PA231	0.0000E+00
PB210	0.0000E+00
PM147	5.2500E+02
PU238	4.2400E+06
PU239	3.9200E+05
PU240	6.9300E+04
PU241	1.9300E+06
PU242	4.9100E+04
PU244	0.0000E+00
RA226	0.0000E+00
RA228	0.0000E+00
SR90	9.8500E+03
TH229	0.0000E+00
TH230	0.0000E+00
TH232	6.1100E-01
U233	1.3100E+03
U234	0.0000E+00
U235	1.1500E+00
U236	0.0000E+00
U238	2.0100E-01

Explanatory Comment: The data below represent the principal results of CUTTINGS_S's calculations, namely, releases to the surface (in Curies) at the time of intrusion. These data are also printed in the (binary) output CDB file. End of Comment.

Information Only

radioisotope RELEASE (CI) for 1 hit
 *** CUTTINGS & CAVING ONLY ***

AM241	2.6296E-01	
AM243	0.0000E+00	
CF252	1.1598E-14	
CM243	0.0000E+00	
CM244	0.0000E+00	
CM245	0.0000E+00	

CM248 CS137 PA231 PB210 PM147 PU238 PU239 PU240 PU241 PU242 PU244 RA226 RA228 SR90 TH229 TH230 TH229 TH230 TH232 U233 U234	2.3237E-08 2.0726E-03 7.6122E-05 2.4936E-09 3.0864E-09 1.7776E-15 1.9421E+00 3.9448E-01 6.9201E-02 1.5811E-02 4.9544E-02 1.8757E-14 6.4693E-09 6.1663E-07 9.1989E-04 1.2425E-05 4.2647E-07 6.1663E-07 1.3215E-03 8.3853E-04		
U235 U236	1.1995E-06 2.0594E-07		
U238	2.0362E-07		
radioiso	tope RELEASE (CI) for 1 h: *** SPALL ONLY ***	it	
AM241 AM243 CF252 CM243 CM244 CM245 CM248 CS137 NP237 PA231 PB210 PM147	2.8188E+00 0.0000E+00 1.2432E-13 0.0000E+00 0.0000E+00 2.4910E-07 2.2217E-02 8.1600E-04 2.6731E-08 3.3085E-08 1.9055E-14		
PU238 PU239 PU240 PU241 PU242 PU244 RA226 RA228 SR90 TH229 TH230 TH230 TH232 U233 U234 U235 U236	2.0819E+01 4.2286E+00 7.4181E-01 1.6949E-01 5.3109E-01 2.0107E-13 6.9349E-08 6.6101E-06 9.8609E-03 1.3320E-04 4.5716E-06 6.6101E-06 1.4166E-02 8.9887E-03 1.2858E-05 2.2076E-06		
U238 radioiso *** TO rati to	2.1828E-06 tope RELEASE (CI) for 1 h TAL CUTTINGS & SPALL *** o of spall area cuttings area 10.72 blow out: SOLID_BLOWOUT 	it 	

OF APPENDIX D END

1

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6

6 2

1 0

0

Number of attributes = = Number of node sets Number of side sets = CUSP_PA96 CPU time is 0:06 (minute:second)

WIPPSLT X-4.00VV 02/24/95 10:37:00 MATSET C-8.07ZO 03/22/95 14:37:07 MATSET POSTLHS C-4.04VV 03/22/95 14:38:39 06/04/96 09:52:46 CUSP_PA9 5.03 Number of coordinates per node = = Number of nodes Number of element blocks = Number of elements = Number of nodes per element = = 14 Number of properties

GENNET C-2.03ZO 03/22/95 14:03:42

MAR 95

*** END OF CUSP_PA96 ***

CUSP_PA96 5.03 PROD PA96 05/23/96

CAMDAT Version 1

QA Records (6):

J.W.	BERGLUND:	SPM2
J.W.	BERGLUND:	SPM2

J.W.	BERGLUND:	SPM2

.T.	TAT .	BERGLIND	CDM2	M

· • • • · · · · · · · · · · · · · · · ·	

(CUSP_MAIN)

CS137	2.4290E-02
NP237	8.9212E-04
PA231	2.9224E-08
PB210	3.6172E-08
PM147	2.0833E-14
PU238	2.2761E+01
PU239	4.6231E+00
PU240	8.1101E-01
PU241	1.8530E-01
PU242	5.8064E-01
PU244	2.1983E-13
RA226	7.5818E-08
D 2220	7 22678-06

7.2267E-06

1.0781E-02

4.9981E-06 7.2267E-06

1.4562E-04

1.5488E-02 9.8272E-03

1.4058E-05 2.4136E-06

2.3864E-06

RA228

TH230

TH232 U233

U234 U235

U236 U238

SR90 TH229

06/04/96 09:52:46

DIS_01 LOCAL 00000000

2 4

APPENDIX E: TEMPLATE INPUT DATA FILE FOR PRE_CUTTINGS

This file either (1) defines or (2) calls from the WIPP controlled database all data required to exercise CUTTINGS_S.

```
11
    Template file for PRE_CUSP WIPP property database
11
11
   MODEL_DATA
11
11
    =wipp::my_database:calc_name
11
=wipp::wipp_db:WIPP_CCA
1
1
     The following input variables are hard wired and are related to the
! repository-model paper by J. W. Berglund (see Appendix A)
 PR_MAX
          <>The maximum pressure allowed by model
Ţ
  PR_MAX
                    15.0E6
! PR_MIN
          <>The minimum pressure allowed by model
                     0.0E6
  PR_MIN
! PE_MAX
          <>The maximum permeability allowed by model
  PE_MAX
                   1.0E-12
          <>The minimum permeability allowed by model
! PE_MIN
  PE_MIN
                   1.0E-17
 PCUT
          <> Pressure (Pa) which defines the border between the
!
                       errosion phase and the bowout phase
1
  PCUT
                     8.0E6
          <> Pressure (pa) which defines the border between the gas-
! PGAS
                       erosion phase and the stuck-pipe phase
1
  PGAS
                    10.0E6
          <> Permeability which defines the phase border the blowout
I.
 KCRIT
                      phase and the gas-erosion/stuck pipe phase
1
  KCRIT
                   1.0E-16
1
 Properties relative to the model (repository) <<<<<<<<<<<<<<<<<
1
1
             Delta radius for derivative in stress calculation
 DELTR
          <>
Ţ
                   0.00001
  DELTR
! ITER
          <> Number of iterations to solve for reynolds number
               1
  ITER
              Gravity effectiveness factor
! FGE
          <>
  BLOWOUT : FGE
! FSE
          <> Strength effectiveness factor
  BLOWOUT:FSE
              Cementation effectiveness factor
! FCE
          <>
  BLOWOUT:FCE
! CEMCEM <> Cementation stress
  BLOWOUT : CEMENT
! SUFTEN <> Surface tension (N/N)
```

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BLOWOUT: SUFTEN

!	Properties of the gas (Hydrogen) in the repository <<<<<<<<	
!	KGAS <> Ratio of Specific Heats BLOWOUT:KGAS	5
i	RGAS <> Gas Constant for Hydrogen BLOWOUT:RGAS	
!	VISC <> Hydrogen viscosity BLOWOUT:VISC	
!	PSUF <> Surface presure (Pa) BLOWOUT:PSUF	
!	Properties to do with the repository <<<<<<<<<<<<<<<<<	Υ.
!	TREPO <> Temperature of repository (K) BLOWOUT:TREPO	en Les
	DEPTH <> Distance from repository depth to depth where casing	
!	is set (m) BLOWOUT 716.0	,
	FLWCNST <> Percent volume of material that is carried by Drilling	
!	mud (unitless) FLWCNST 0.05	
!	HREPO <> Height of repository at burial time (4m) BLOWOUT:HREPO	
!	RPANEL <> The equivalent radius of 1 panel (910.0m**2) BLOWOUT:RPANEL	
ł	ROOM <> The equivalent radius of 1 room (11,640m**2) BLOWOUT:ROOM	• •
ł	RHOS <> Waste particle density (kg/m**3) BLOWOUT:RHOS	
,	Variables to do with drilling <<<<<<<<<<<<<	
	L1 <> Collar Length (m)	
:	BOREHOLE:L1	
!	L2 <> Drill pipe length (m) BOREHOLE:L2	
!	COLDIA <> Collar diameter (m) BOREHOLE:COLDIA	
ţ	PIPED <> Drill pipe diameter (m) BOREHOLE:PIPED	
!	ROUGHP <> Friction factor (unitless) BOREHOLE:ROUGHP	
! ! !	APORO <> A constant in equation to determine permeability as a function of porosity BLOWOUT:APORO	
! !	NPORO <> N constant in equation to determine permeability as a function of porosity	
	Information Only	y

BLOWOUT

0.0

! END_OF_MODEL_INPUT

RADIONUCLIDE_DATA

! GEOMETRY

!

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BOREHOLE: INV_AR

BOREHOLE:RHW_AR

BOREHOLE:WUF

!MATERIAL OUT_MAT BOREHOLE

!REPOSITORY_TYPE

REP_NAME WIPP REP_GEOLogy HALITE

RADWASte_type CONtact_handled

!RADIOISOTOPE_chains

```
! chain1/chain2 from U234 & down are the same:
! (Both chains are required input)
```

!	\/	$\backslash /$	\/	17	\/	\/	\/
CHAIN1	PU242 RN222	U238 PO218	TH234 PB214	PA234M BI214		TH230 PB210	RA226 <
CHAIN2	PU238 PB214	U234 BI214	TH230 PO214	RA226 PB210	RN222 <	PO218	

! chain3/chain4 from PU239 & down are the same:

CHAIN3	3 AM243 TH227	NP239 RA223	PU239 RN219	U235 PO215	TH231 PB211	PA231 BI211	AC227 TL207	<
CHAIN4	1 CM243 RA223	PU239 RN219	U235 PO215	TH231 PB211	PA231 BI211	AC227 TL207	TH227 PB209	<
! chai	in5/chain6	from U236	& down	are the	same:			
CHAINS	5 CF252 AC228 PO212	CM248 TH228 <	PU244 RA224	PU240 RN220	U236 PO216	TH232 PB212	RA228 BI212	
CHAIN	5 CM244 RA224	PU240 RN220	U236 PO216	TH232 PB212	RA228 BI212	AC228 PO212	TH228 <	
CHAINT	7 CM245 RA225	PU241 AC225	AM241 FR221	NP237 AT217	PA233 BI213	U233 PO213	TH229 <	
CHAIN	B CS137	BA137M	<					•
CHAINS	9 PM147	SM147	ND143	<				
CHAIN	10 SR90	¥90	ZR90	<				
1	^	^	^	^	^	^	^	

SAVE AM241 AM243 CF252 CM243 CM244 CM245 CM248 CS137 NP237 PA231 PB210 PM147 PU238 PU239 PU242 PU244 RA226 RA228 SR90 TH229 PU240 PU241 TH229 TH230 TH232 11233 U234 U235 U236 U238 <

TABULAR_DATA

! END_OF_RADIOISOTOPE_INPUT

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! Example of how radioisotope data are input: !...1st Line: Radionuclide (an asterisk in column 1 follow by radionuclide name, ex; *AC225) !...2nd & 3rd line [REAL] (3(11x, 1pe14.6)) !...Field#1 Atomic Weight (Kg/Mole) AWT HALFY !...Field#2 Half-Life (Years) [REAL] !...Field#3 Activity Conversion (Ci/Kg) AWTCNV [REAL] !...Field#4 EPA Release Limit (Ci) EPAREL [REAL] !...Field#5 Inventory (Ci) INVCHD [REAL] !...Field#6 Inventory (Ci) INVRHD [REAL] !*PU241 2.410000E-01 HALFY 1.439900E+01 ACTCNV 1.030000E+05 AWT 1 INVRHD 0.000000E+00 EPAREL 1.000000E+07 INVCHD 1.930000E+06 t <TABLE_INPUTS <GENERATE_RADIO> END_TABLES>

END OF APPENDIX E

APPENDIX F: OUTPUT ASCII FILE FROM PRE_CUTTINGS

This appendix lists the ASCII data file that results from exercising PRE_CUTTINGS. It is controlled under CMS and must be used exactly as it appears. Other output files are possible, but they too are recorded and controlled under the CMS and it is clear which results derive from which output ASCII files.

MODEL_DATA

! ! J		following input variables are hard wired and related to the Repository model glund paper: Appendix A, CUTTING_S User's Manual, Ver 5.03
!	PR_MAX PR_MAX	<>The maximum pressure allowed by model 15.0E6
!	PR_MIN PR_MIN	<>The minimum pressure allowed by model 0.0E0
!	PE_MAX PE_MAX	<>The maximum permeability allowed by model 1.0E-12
!	PE_MIN PE_MIN	<>The minimum permeability allowed by model 1.0E-17
!	PCUT PCUT	<> Pressure (Pa) which defines the border between the erosion phase and the blowout phase 8.0E6
! !	PGAS PGAS	<> Pressure (Pa) which defines the border between the gas erosion phase and the stuck pipe phase 10.0E6
! !	KCRIT KCRIT	<> Permeability which defines the phase border between the blowout phase and the gas erosion/stuck pipe phase 1.0E-16
! 1 !	Propert DELTR	ies relative to the model (repository) <<<<<<<<<<<<<<<<>>> Delta radius for derivative in stress calculation
!	DELTR DELTR	0.00001

Information Only

! ITER <> Number of iterations to solve for Reynolds number ITER 1

- ! FGE <> Gravity effectiveness factor FGE 18.4
- ! FSE <> Strength effectiveness factor FSE 0.0

! FCE <> Cementation effectiveness factor FCE 0.0

! CEMENT <> Cementation stress CEMENT 0.0

! SUFTEN <> Surface tension (N/N) SUFTEN 8.0E-2 ! Properties of the gas (Hydrogen) in the repository <<<<<<<<<<> <> Ratio of Specific Heats ! KGAS 1.41E0 KGAS ! RGAS <> Gas Constant for Hydrogen (Nm/kg-K) RGAS 4116.0 ! VISC <> Hydrogen viscosity (Pa-sec) VISC 0.92E-5 ! PSUF <> Surface presure (Pa) 0.089465E6 PSUF ! Properties to do with the repository <<<<<<<<<<<<<<<< <> Temperature of repository (K) Ţ TREPO TREPO 300.0 ! DEPTH <> Distance from repository to where casing is set and gas erosion ceases (m) 1 DEPTH 716.0 ! FLWCNST <> Percent volume of material that is carried by Drilling mud (unitless) 1 FLWCNST 0.05 ! HREPO <> Height of repository at burial time (4m)HREPO 3.96 ! RPANEL <> The equivalent radius of 1 panel (116400m**2) RPANEL 60.87 <> The equivalent radius of 1 room (910.0m**2) ! ROOM 17.1 ROOM <> Waste particle density (kg/m**3) ! RHOS RHOS 2650.0 ! Variables to do with drilling <<<<<<<<<<< 1 ! L1 <> Collar Length (m) 182.88 L1 <> Drill pipe length (m) ! L2 472.12 L2 <> Collar diameter (m) ! COLDIA 0.2032004 COLDIA ! PIPED <> Drill pipe diameter (m) PIPED 0.1143002 <> Friction factor (unitless) ! ROUGHP ROUGHP 0.08 ļ APORO <> A constant in equation to determine permeability as a ı function of porosity APORO 1.0E-11 <> N constant in equation to determine permeability as a ! NPORO function of porosity Ţ NPORO 4.6

! ! END_OF_MODEL_INPUT

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RADIONUCLIDE_DATA ! GEOMETRY INV_AR 1.11520E+05 RHW_AR 1.11520E+05 WUF 5.25 **!MATERIAL** BOREHOLE OUT_MAT !REPOSITORY_TYPE REP_NAME WIPP REP_GEOLogy HALITE RADWASte_type CONtact_handled !RADIOISOTOPE_chains chain1/chain2 from U234 & down are the same: (Both chains are input) $\backslash /$ $\backslash /$ $\backslash /$ \/ ! $\backslash /$ \mathbf{X} \mathbf{V} CHAIN1 PU242 U238 TH234 PA234M U234 TH230 RA226 RN222 PO218 PB214 BI214 PO214 PB210 < U234 TH230 RA226 RN222 PO218 CHAIN2 PU238 PB214 BI214 PO214 PB210 < ! chain3/chain4 from PU239 & down are the same: AC227 AM243 NP239 PU239 U235 TH231 PA231 CHAIN3 PB211 BI211 TL207 TH227 RA223 RN219 PO215 < TH227 CHAIN4 CM243 PU239 U235 TH231 PA231 AC227 PB211 TL207 RA223 RN219 PO215 BI211 < ! chain5/chain6 from U236 & down are the same: TH232 RA228 CHAIN5 CF252 CM248 PU244 PU240 U236 TH228 RA224 RN220 PO216 PB212 BI212 AC228 PO212 < PU240 TH232 RA228 AC228 TH228 CHAIN6 CM244 U236 RA224 RN220 PO216 PB212 BI212 PO212 < PU241 AM241 NP237 PA233 U233 TH229 CHAIN7 CM245 BI213 PO213 PB209 < RA225 AC225 FR221 AT217 BA137M CHAIN8 CS137 < PM147 SM147 ND143 CHAIN9 < CHAIN10 SR90 ¥90 zr90 < ~ ^ ^ ^

SAVE AM241 AM243 CF252 CM243 CM244 CM245 CM248 CS137 PA231 PB210 PM147 PU238 PU239 PU240 PU241 NP237 PU244 SR90 TH230 RA228 TH229 TH232 PU242 RA226 U234 U235 U236 U238 U233 < TABULAR_DATA ļ Example of how radioisotope data are entered: ļ ł ...1st Line: Radionuclide (an asterisk in column 1 follow - 1 by radionuclide name, ex; *AC225) ...2nd & 3rd line ... Field#1 Atomic Weight (Kg/Mole) AWT [REAL] (3(11x,1pe14.6)) [REAL] HALFY ...Field#2 Half-Life (Years) ... Field#3 Activity Conversion (Ci/Kg) AWTCNV [REAL] 11 (Ci) EPAREL [REAL] ... Field#4 EPA Release Limit ... Field#5 Inventory (Ci) INVCHD [REAL] (Ci) INVRHD [REAL] ... Field#6 Inventory !*PU241 1.439900E+01 ACTCNV 1.030000E+05 2.410000E-01 HALFY AWT 0.00000E+00 INVRHD EPAREL 1.00000E+07 INVCHD 1.930000E+06 <TABLE_INPUTS *AC225 2.737909E-02 5.802700E+07 AWT 2.250230E-01 HALFY ACTCNV EPAREL 0.00000E+00 INVCHD 0.00000E+00 *AC227 2.177335E+01 ACTCNV 7.232300E+04 AWT 2.270280E-01 HALFY 0.000000E+00 1.000000E+02 INVCHD EPAREL *AC228 2.241700E+09 2.280310E-01 HALFY 6.993710E-04 ACTCNV AWT 0.00000E+00 EPAREL 0.00000E+00 INVCHD *AM241 4.322347E+02 ACTCNV 3.431200E+03 2.410570E-01 HALFY AWT 1.00000E+02 INVCHD 2.400000E+05 EPAREL *AM243 1.992900E+02 ACTCNV 2.430610E-01 HALFY 7.380313E+03 AWT 0.00000E+00 EPAREL 1.000000E+02 INVCHD *AT217 1.609500E+15 AWT 2.170050E-01 HALFY 1.023547E-09 ACTCNV 0.00000E+00 INVCHD 0.00000E+00 EPAREL *BA137M 5.382400E+11 4.851550E-06 ACTCNV AWT 1.369070E-01 HALFY 0.00000E+00 0.00000E+00 INVCHD EPAREL *BI211 4.049824E-06 ACTCNV 4.184000E+11 2.109870E-01 HALFY AWT 0.00000E+00 INVCHD EPAREL 0.00000E+00 *BI212 1.464800E+10 HALFY 1.151253E-04 ACTCNV 2.119910E-01 AWT 0.00000E+00 INVCHD 0.00000E+00 EPAREL *BT213 1.933800E+10 8.679553E-05 ACTCNV HALFY AWT 2.129940E-01 INVCHD 0.00000E+00 EPAREL 0.00000E+00 *BI214 4.415300E+10 3.783638E-05 ACTCNV 2.139990E-01 HALFY AWT 0.00000E+00 EPAREL 0.00000E+00 INVCHD *CF252 5.375899E+05 HALFY 2.638090E+00 ACTCNV AWT 2.520820E-01 0.00000E+00 INVCHD 2.960000E+03 EPAREL *CM243 2.850088E+01 ACTCNV 5.160700E+04 HALFY ጉመል 2.430610E-01 0.00000E+00 EPAREL 1.000000E+02 INVCHD

*CM244

-

AWT EPAREL	2.440630E-01 0.000000E+00	HALFY INVCHD	1.811013E+01 0.000000E+00	ACTCNV	8.088299E+04
*CM245 AWT EPAREL	2.450650E-01 1.000000E+02	HALFY INVCHD	8.498927E+03 0.000000E+00	ACTCNV	1.716500E+02
*CM248 AWT EPAREL	2.480720E-01 1.000000E+02	HALFY INVCHD	3.390698E+05 0.000000E+00	ACTCNV	4.250200E+00
*CS137 AWT EPAREL	1.369070E-01 1.000000E+03	HALFY INVCHD	2.999975E+01 2.070000E+04	ACTCNV	8.704300E+04
*FR221 AWT EPAREL	2.210140E-01 0.000000E+00	HALFY INVCHD	9.126364E-06 0.000000E+00	ACTCNV	1.772400E+11
*ND143 AWT EPAREL	1.430000E-01 1.000000E+03	HALFY INVCHD	3.168876E+30 0.000000E+00	ACTCNV	0.000000E+00
*NP237 AWT EPAREL	2.370480E-01 1.000000E+02	HALFY INVCHD	2.139942E+06 6.670000E+01	ACTCNV	7.047600E-01
*NP239 AWT EPAREL	2.390530E-01 0.000000E+00	HALFY INVCHD	6.448664E-03 0.000000E+00	ACTCNV	2.319100E+08
*PA231 AWT EPAREL	2.310360E-01 1.000000E+02	HALFY INVCHD	3.276618E+04 0.000000E+00	ACTCNV	4.722500E+01
*PA233 AWT EPAREL	2.330400E-01 0.000000E+00	HALFY INVCHD	7.392988E-02 0.000000E+00	ACTCNV	2.075100E+07
*PA234M AWT EPAREL *PB209	2.340430E-01 0.000000E+00	HALFY INVCHD	2.224551E-06 0.000000E+00	ACTCNV	6.866600E+11
AWT EPAREL *PB210	2.089810E-01 0.000000E+00	HALFY INVCHD	3.764625E-04 0.000000E+00	ACTCNV	4.544100E+09
AWT EPAREL *PB211	2.099840E-01 1.000000E+02	HALFY INVCHD	2.229938E+01 0.000000E+00	ACTCNV	7.634800E+04
AWT EPAREL * PB212	2.109890E-01 0.000000E+00	HALFY INVCHD	6.863786E-05 0.000000E+00	ACTCNV	2.468600E+10
AWT EPAREL *PB214	2.119920E-01 0.000000E+00	HALFY INVCHD	1.213680E-03 0.000000E+00	ACTCNV	1.389500E+09
AWT EPAREL *PM147	2.140000E-01 0.000000E+00	HALFY INVCHD	5.095553E-05 0.000000E+00	ACTCNV	3.278500E+10
AWT EPAREL *PO212	1.469150E-01 0.000000E+00	HALFY INVCHD	2.623513E+00 5.250000E+02	ACTCNV	9.275300E+05
AWT EPAREL *PO213	2.119890E-01 0.000000E+00	HALFY INVCHD	9.506629E-15 0.000000E+00	ACTCNV	1.773900E+20
AWT EPAREL *PO214	2.129930E-01 0.000000E+00	HALFY INVCHD	1.330928E-13 0.000000E+00	ACTCNV	1.261100E+19
AWT EPAREL *PO215	2.139950E-01 0.000000E+00	HALFY INVCHD	5.206464E-12 0.000000E+00	ACTCNV	3.208700E+17
AWT EPAREL *PO216	2.149990E-01 0.000000E+00	HALFY INVCHD	5.640600E-11 0.000000E+00	ACTCNV	2.947900E+16
*PO216 AWT EPAREL *PO218	2.160020E-01 0.000000E+00	HALFY INVCHD	4.753314E-09 0.000000E+00	ACTCNV	3.482000E+14
*PO218 AWT EPAREL *PU238	2.180090E-01 0.000000E+00	HALFY INVCHD	5.799044E-06 0.000000E+00	ACTCNV	2.827800E+11
AWT	2.380500E-01	HALFY	8.774619E+01	ACTCNV	1.711500E+04

EPAREL	1.000000E+02	INVCHD	4.240000E+06		
*PU239 AWT	2.390520E-01	HALFY	2.406445E+04	ACTCNV	6.214600E+01
EPAREL *PU240	1.000000E+02	INVCHD	3.920000E+05		۹.
AWT	2.400540E-01	HALFY	6.537392E+03	ACTCNV	2.278100E+02
EPAREL *PU241	1.000000E+02	INVCHD	6.930000E+04		
AWT EPAREL	2.410570E-01 0.000000E+00	HALFY INVCHD	1.439937E+01 1.930000E+06	ACTCNV	1.030000E+05
* PU242	2.420590E-01	HALFY	3.869198E+05		3.817100E+00
AWT EPAREL	1.000000E+02	INVCHD	4.910000E+04	ACTCNV	3.81/1002+00
* PU244 AWT	2.440640E-01	HALFY	8.261261E+07	ACTCNV	1.773100E-02
EPAREL *RA223	1.000000E+02	INVCHD	0.00000E+00		
AWT	2.230190E-01	HALFY	3.130533E-02	ACTCNV	5.120600E+07
EPAREL *RA224	0.000000E+00	INVCHD	0.000000E+00		
AWT EPAREL	2.240200E-01 0.000000E+00	HALFY INVCHD	1.001999E-02 0.000000E+00	ACTCNV	1.592700E+08
*RA225			4.052993E-02		3.919900E+07
AWT EPAREL	2.250240E-01 0.000000E+00	HALFY INVCHD	4.052993E-02 0.000000E+00	ACTCNV	3.919900E+07
*RA226 AWT	2.260250E-01	HALFY	1.599966E+03	ACTCNV	9.885801E+02
EPAREL *RA228	1.000000E+02	INVCHD	0.000000E+00		
AWT	2.280310E-01	HALFY	6.699955E+00	ACTCNV	2.340000E+05
EPAREL *RN219	0.00000E+00	INVCHD	0.000000E+00		
AWT EPAREL	2.190090E-01 0.000000E+00	HALFY INVCHD	1.254875E-07 0.000000E+00	ACTCNV	1.300800E+13
*RN220				20002007	9.222600E+11
AWT EPAREL	2.200110E-01 0.000000E+00	HALFY INVCHD	1.761895E-06 0.000000E+00	ACTCNV	9.2226002+11
*RN222 AWT	2.220180E-01	HALFY	1.046997E-02	ACTCNV	1.538000E+08
EPAREL *SM147	0.000000E+00	INVCHD	0.000000E+00		.*
AWT	1.470000E-01	HALFY	1.059355E+11 0.000000E+00	ACTCNV	2.295400E-05
EPAREL *SR90	1.000000E+02	INVCHD			
AWT EPAREL	8.990801E-02 1.000000E+03	HALFY INVCHD	2.912197E+01 9.850002E+03	ACTCNV	1.365400E+05
*TH227 AWT	2.270280E-01	HALFY	5.124073E-02	ACTCNV	3.073200E+07
EPAREL *TH228	0.000000E+00	INVCHD	0.000000E+00		
AWT	2.280290E-01	HALFY	1.913051E+00	ACTCNV	8.195301E+05
EPAREL *TH229	0.000000E+00	INVCHD	0.000000E+00		
AWT EPAREL	2.290320E-01 1.000000E+02	HALFY INVCHD	7.339118E+03 0.000000E+00	ACTCNV	2.126900E+02
*TH230		HALFY	7.700370E+04	ACTCNV	2.018300E+01
AWT EPAREL	2.300330E-01 1.000000E+01	INVCHD	0.000000E+00	ACICINV	2.0105002+01
*TH231 AWT	2.310360E-01	HALFY	2.911247E-03	ACTCNV	5.315200E+08
EPAREL *TH232	0.000000E+00	INVCHD	0.00000E+00		· .
AWT	2.320380E-01	HALFY	1.405080E+10	ACTCNV	1.096500E-04
EPAREL *TH234	1.000000E+01	INVCHD	6.110000E-01		
AWT EPAREL	2.340440E-01 0.000000E+00	HALFY INVCHD	6.597601E-02 0.000000E+00	ACTCNV	2.315200E+07
*TL207	2.069770E-01	HALFY	9.069325E-06	ACTCNV	1.904500E+11
AWT EPAREL	0.000000E+00	INVCHD	0.000000E+00	110 1 0144	T12030000111

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*U233					
AWT	2.330400E-01	HALFY	1.585072E+05	ACTCNV	9.678300E+00
EPAREL	1.000000E+02	INVCHD	1.310000E+03		
*U234					1
AWT	2.340410E-01	HALFY	2.445105E+05	ACTCNV	6.247300E+00
EPAREL	1.000000E+02	INVCHD	0.00000E+00		· .
*U235					
AWT	2.350440E-01	HALFY	7.038074E+08	ACTCNV	2.161100E-03
EPAREL	1.000000E+02	INVCHD	1.150000E+00		
*U236					
AWT	2.360460E-01	HALFY	2.341483E+07	ACTCNV	6.468300E-02
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*U238					
AWT	2.380510E-01	HALFY	4.468115E+09	ACTCNV	3.361100E-04
EPAREL	1.000000E+02	INVCHD	2.010000E-01		
*Y90					
AWT	8.990701E-02	HALFY	7.301091E-03	ACTCNV	5.446300E+08
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
*ZR90					
AWT	9.000000E-02	HALFY	3.168876E+30	ACTCNV	0.00000E+00
EPAREL	1.000000E+03	INVCHD	0.00000E+00		
END_TABLES>					
1					

C NAGE.

I END_OF_RADIOISOTOPE_INPUT

!

END OF APPENDIX F

APPENDIX G: SAMPLE COMMAND FILE FOR PRE_CUTTINGS

This command file exercises PRE_CUTTINGS.

```
$!@CUSP_RUN_PRE_CUTTINGS.COM
Ś!
Ś
  ON control_y
                 THEN goto TERMINATE
                 THEN goto TERMINATE
  ON control_c
Ś
$
  ON error
                 THEN goto TERMINATE
$!
$!
$! The production executable
$!
  DEFINE CUSP$EXE WP$PRODROOT: [CUSP.EXE]CUSP_PA96.EXE
$
$1
$ !
$! This sets CUSP into the pre_cuttings mode
$!
  DEFINE CUSP$PRE_CUTTINGS TRUE
$
$!
$!
   Input files required by PRE_CUTTINGS
$!
$!
   This file is a template to reference the WIPP CMS-controlled Data Base
$!
    It will have the chains and some hardwired numbers which
$!
     can be found in the theoretical manual.
$!
   The file will be placed into CMS CUSP
<u>$!</u>
$!
                 CUSP_INP$CDB
                                  CUSP_TEST_1.CDB
$ DEFINE/NOLOG
$!
$ DEFINE/NOLOG
                 CUSP_INP$TXT1
                                  CUSP_TEST_1.INP
$!
$ DEFINE/NOLOG
                 CUSP_INP$SDB
                                  CUSP_CCA_TEMPLATE.INP
$!
$ DEFINE/NOLOG
                 CUSP_INP$DESC
                                  CUSP_CCA.SDB
$!
$!
$! A summary of the results of the pre_cuttings
51
$ DEFINE/NOLOG
                 CUSP_OUT$DBG
                                       PRE_CUSP_OUTPUT.DBG
$!
$!
$!
   The output ASCII file that is read by all production runs
$!
 DEFINE/NOLOG
                 CUSP_OUT$DB_ASC
                                       CUSP_CCA_SDB.ASC
Ś
$1
$!
   The output BINARY file that could be read by all the production runs
$!
                                 ~~~~
$1
$!
S DEFINE/NOLOG
                 CUSP_OUT$DB_BIN
          CUSP_INPUT.BINARY
$!
$!
   CUSP
         := $CUSP$EXE
$!
$!
   CUSP
$
$!
$!
$ TERMINATE:
$ !
$! Turn off pre_cuttings mode
$!
$ DEASSIGN
                 CUSP$PRE_CUTTINGS
<u>$</u>!
$ DEASSIGN
                 CUSP_INP$CDB
```

:

CUSP_INP\$TXT1 \$ DEASSIGN CUSP_INP\$SDB CUSP_OUT\$DB_ASC CUSP_OUT\$DB_BIN \$ DEASSIGN \$ DEASSIGN \$ DEASSIGN \$! \$TERMINATE:

.

\$ EXIT

END OF APPENDIX G

APPENDIX H: SAMPLE CHECK_OUT PROBLEMS

This file includes a command file that exercises CHECK_OUT and a sequence of actual input control files associated with selected applications of interest.

Information Only

General Purpose ".COM" file to run CHECK_OUT modes

```
$!@CUSP_CHECK_OUT.COM
                          "p1"
$!
$!
    'p1' > Check_Out Type
$!
$!
   Where 'p1': CALIBRATE
                 EXPERIMENT
$!
$!
                 VALIDATE
$!
$!
    ON control_y THEN goto TERMINATE ON control_c THEN goto TERMINATE
$
$
$
                  THEN goto TERMINATE
    ON error
$!
$!
$
$!
    LIBCUSP
                .
$! Clea up any prexisting logicals
$!
                                              $
   IF F$SEARCH("CUSP_DEASSIGN.COM") .eqs.
      THEN CFE CUSP_DEASSIGN.COM
$!
   @CUSP_DEASSIGN
$
$!
$!
$
$
   IF F$SEARCH("CUSP$EXE") .eqs. ""
   THEN
$!
     IF F$SEARCH("CUSP_TEST_SETUP_CUSP.COM") .eqs.
                                                        0.0
$
        THEN CFE CUSP_TEST_SETUP_CUSP.COM
$!
$
     @CUSP_TEST_SETUP_CUSP
$!
$
   ENDIF
$!
$!
$
   DEFINE
               CUSP$CHECK_OUT TRUE
$!
$!
$!
    Input files required by CUSP$CHECK_OUT
$!
                  CUSP_INP$TXT0
                                      CUSP_CHECK_OUT.DAT
$ DEFINE/NOLOG
$ DEFINE/NOLOG
                  CUSP_INP$TXT1
                                     CUSP_'p1'.INP
$!
$!
   Output files required by CUSP$CHECK_OUT
$!
                  CUSP_OUT$NVERFY
                                      CUSP_'p1'.VER
$ DEFINE/NOLOG
                                      CUSP_'p1'.PLT
$ DEFINE/NOLOG
                  CUSP OUT$PLT1
$!
$!
   CUSP
          := $CUSP$EXE
$!
   CUSP
Ŝ
$!
$ DEASSIGN CUSP$CHECK_OUT
$!
$ DEASSIGN CUSP_INP$TXT0
$ DEASSIGN CUSP_INP$TXT1
```

ζ.

\$ DEASSIGN CUSP_OUT\$NVERFY \$ DEASSIGN CUSP_OUT\$PLT1 \$! \$! **\$TERMINATE:** \$! \$ EXIT

Input file for: CALIBRATE

!		
1	TEST_FUNCT	
•	CALIBRATE	
	FSEMIN	10.0
	DELFSE	5.0
	IFSE	3
	FGEMIN	10.0
	DELFGE	10.0
	JFGE	3
	FCEMIN	0.0
	DELFCE	0.0
	KFCE	0
	ITMAX	200
	DELERR	0.01
	POROST	0.40

START_SEARCH

!

.000151	.01021	2.133	394379.	.0007221	.261	295.	85151.	#28
.000151	.01021	2.133	290958.	.0005760	.261	295.	85151.	#28
.000151	.01021	2.133	290958.	.0005057	.262	295.	84805.	#29
.000151	.01021	2.133	290958.	.0006104	.279	296.	84048.	#30
.000151	.01021	2.133	222010.	.0004148	.262	295.	84805.	#29
.000151	.01021	2.133	222010.	.0004817	.261	295.	85151.	#28
.000151	.01021	2.133	222010.	.0004800	.279	296.	84048.	#30
.000151	.01021	2.133	153064.	.0003005	.262	295.	84805.	#29
.000151	.01021	2.133	153064.	.0003642	.261	295.	85151.	#28
.000151	.01021	2.133	153064.	.0003444	.279	296.	84048.	#30
.000151	.01021	2.133	118590.	.0001261	.262	295.	84805.	#29
.000151	.01021	2.133	118590.	.0000663	.261	295.	85151.	#28
.000151	.01021	2.133	118590.	.0001740	.279	296.	84048.	#30

Input file for: EXPERIMENT

!	TEST_FUNCT										
:	EXPERIMENT										
	DIA 0.311	0									
	PARTMIN	0.0002	PARTMAX	0.2	PARTMUL	1.5					
	SAT	0.3									

PG	15.0E6
POROF	0.4
FSE	0.010
FGE	1.0

Input file for: VALIDATE

!	TEST_FUN	СТ					
!	VALIDAT	Е					
:	DIA	0.3110					
: :	PARTMIN	0.00002	PARTMAX	0.2	PARTMU	L	2.0
:	SATMIN	0.20	DELSAT	0.05	ISA	9	
:	PGMIN	8.0E6	DELPG	1.0E6	IPG	8	
!	PORMIN	0.2	DELPOR	0.10	IPOR	7	

.

Input file: CUSP_CHECK_OUT.DAT

MODEL_DATA

Ţ

```
I.
 The following input variables are related to the Repository and
!
  are hard wired or related to the model by J. W. Berglund (see Appendix A)
1
1
! Properties relative to the model (experiments) <<<<<<<<<<<<<<
! DELTR
         <> Delta radius for derivative in stress calculation
              0.0001
 DELTR
! ITER
             Number of iterations to solve for Reynolds number
         <>
 ITER
              7
! SUFTEN <> Surface tension (N/m)
 SUFTEN
             7.3E-2
! Properties of Air (experiments) <<<<<<<<<
         <> Ratio of Specific Heats
! KGAS
 KGAS
              1.4E0
! RGAS
         <> Gas Constant for Air
              286.5
 RGAS
! VISC
         <> Air viscosity
 VISC
              1.77E-5
! Properties to do with the experiments <<<<<<<<
         <> Height of repository at burial time
! HREPO
 HREPO
              0.0762
```

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!

<> The equivalent radius of 1 room ! ROOM ROOM 0.254 ! RHOS <> Waste particle density (kg/m**3) 2650.0 RHOS ! Variables to do with drilling <<<<<<<< <> Drill pipe length (m) ! L2 L2 1.0 ! COLDIA <> Collar diameter (m) COLDIA 0.0 ! PIPED <> Drill pipe diameter (m) 0.0 PIPED ! END_OF_MODEL_INPUT

d'

END OF APPENDIX H

1

APPENDIX I: EXAMPLE OF THE INPUT FILE: CUSP_INP\$TXT0 FOR TEST RUNS

This input file provides various data relevant to the position, size, and nature of the repository and to physical properties of the materials in and around the borehole. The file itself is located in the CMS portion of the WIPP Alpha cluster as : CUSP_TEST_REPOSITORY.DAT. The file is discussed in Section 6.1 as file number 2.

To extract and edit this file from the CMS system, type the following sequence of commands, each followed by a carriage return.

\$LIBCUSP \$CFE CUSP_TEST_REPOSITORY.DAT EDIT CUSP_TEST_REPOSITORY.DAT _____ OF BEGINNING CUSP_TEST_REPOSITORY.DAT _____ MODEL DATA ! 1 The following input variables are hard wired and related to the 1 ! Repository model J. W. Berglund paper: Appendix A, CUTTING_S User's Manual, Ver 5.03 1 ! PR_MAX <>The maximum pressure allowed by model 15.0E6 PR_MAX ! PR_MIN <>The minimum pressure allowed by model 0.0E0 PR_MIN ! PE_MAX <>The maximum permeability allowed by model PE_MAX 1.0E-12 ! PE_MIN <>The minimum permeability allowed by model PE_MIN 1.0E-17 ! PCUT <> Pressure (Pa) which defines the border between the 1 erosion phase and the blowout phase PCUT 8.0E6 <> Pressure (Pa) which defines the border between the gas ! PGAS erosion phase and the stuck pipe phase 1 PGAS 10.0E6 <> Permeability which defines the phase border between the blowout ! KCRIT phase and the gas erosion/stuck pipe phase Ţ 1.0E-16 KCRIT

! Properties relative to the model (repository) <<<<<<<<<<<<<<

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! !	DELTR DELTR	<>	Delta radius for derivative in stress calculation 0.00001
!	ITER ITER	<>	Number of iterations to solve for reynolds number 1
!	FGE FGE	<>	Gravity effectiveness factor 18.4
!	FSE FSE	<>	Strength effectiveness factor 0.0
!	FCE FCE	<>	Cementation effectiveness factor 0.0
!	CEMENT CEMENT	<>	Cementation stress 0.0
!	SUFTEN SUFTEN	<>	Surface tension (N/N) 8.0E-2
!	Propert:	ies	of the gas (Hydrogen) in the repository <<<<<<<<<<<
!	KGAS KGAS	<>	Ratio of Specific Heats 1.41E0
!	RGAS RGAS	<>	Gas Constant for Hydrogen (Nm/kg-K) 4116.0
!	VISC VISC	<>	Hydrogen viscosity (Pa-sec) 0.92E-5
!	PSUF PSUF	<>	Surface presure (Pa) 0.089465E6
!	Propert	ies	to do with the repository <<<<<<<<<<>>>>>>>>>>>>>>>>>>>>>>>>>>>
!	TREPO TREPO	<>	Temperature of repository (K) 300.0
! !	DEPTH	<>	Distance from repository to where casing is set and gas erosion ceases (m)
·	DEPTH		716.0 ·
	FLWCNST	<>	Percent volume of material that is carried by Drilling
!	FLWCNST		mud (unitless) 0.05
!	HREPO HREPO	<>	Height of repository at burial time (4m) 3.96
!	RPANEL RPANEL	<>	The equivalent radius of 1 panel (116400m**2) 60.87
!	ROOM	<>	The equivalent radius of 1 room (910.0m**2)

```
ROOM
                      17.1
          <> Waste particle density (kg/m**3)
! RHOS
 RHOS
                    2650.0
! Variables to do with drilling <<<<<<<<<
1
! L1
          <> Collar Length (m)
 L1
                    182.88
! L2
          <> Drill pipe length (m)
 L2
                    472.12
! COLDIA <> Collar diameter (m)
                 0.2032004
 COLDIA
! PIPED
         <> Drill pipe diameter (m)
                 0.1143002
  PIPED
! ROUGHP <> Friction factor (unitless)
                      0.08
 ROUGHP
1
! APORO
          <> A constant in equation to determine permeability as a
              function of porosity
1
                   1.0E-11
 APORO
! NPORO
          <> N constant in equation to determine permeability as a
              function of porosity
1
 NPORO
                       4.6
1
! END_OF_MODEL_INPUT
Į.
```

RADIONUCLIDE_DATA

! GEOMETRY

INV_AR 1.11520E+05

RHW_AR 1.11520E+05

WUF 5.25

!MATERIAL OUT_MAT BOREHOLE

!REPOSITORY_TYPE

REP_NAME WIPP REP_GEOLogy HALITE

RADWASte_type CONtact_handled

!RADIOISOTOPE_chains

!

! chain1/chain2 from U234 & down are the same: ! (Both chains are required input)								
!	\/	\/	\/ \	\/ \	\/	\/	\/	
CHAIN1		U238 PO218					RA226 <	
CHAIN2		U234 BI214				PO218	1	
! chain3/chain4 from PU239 & down are the same:								
CHAIN3	AM243 TH227	NP239 RA223	PU239 RN219		TH231 PB211			<
CHAIN4		PU239 RN219					TH227 <	
! chain!	5/chain6	from U236	& down	are the	same:			
CHAIN5		TH228	PU244 RA224					
CHAIN6	CM244 RA224	PU240 RN220		TH232 PB212	RA228 BI212			
CHAIN7	CM245 RA225	5 PU241 5 AC225	AM241 FR221					<
CHAIN8	CS137	BA137M	<					ı
CHAIN9	PM147	SM147	ND143	<				
CHAIN10	SR90	¥90	ZR90	<				
!	^	^	^	^	^	^	^	
SAVE	NP237 PU242	AM243 CF2 PA231 PB2 PU244 RA2 U234 U23	10 PM14 26 RA22	7 PU238 8 SR90	PU239 TH229			
TABULAR_DATA								
! ! Example of how the radioisotope data are input: !								
<pre>!1st Line: Radionuclide (an asterisk in column 1 follow ! by radionuclide name, ex; *AC225)</pre>								

...2nd & 3rd line ! ! (3(11x, 1pe14.6))...Field#1 Atomic Weight (Kg/Mole) AWT [REAL] 1 ...Field#2 Half-Life (Years) HALFY [REAL] ...Field#3 Activity Conversion (Ci/Kg) AWTCNV [REAL] 11 ! = ! -...Field#4 EPA Release Limit (Ci) EPAREL [REAL] !

...Field#5 Inventory (Ci) INVCHD [REAL] 1 " (Ci) 1 ... Field#6 Inventory INVRHD [REAL] 1 !*PU241 AWT 2.410000E-01 HALFY 1.439900E+01 ACTCNV 1.030000E+05 1.00000E+07 INVCHD 1.930000E+06 INVRHD 0.00000E+00 EPAREL <TABLE_INPUTS *AC225 2.250230E-01 HALFY 2.737909E-02 ACTCNV 5.802700E+07 AWT EPAREL 0.00000E+00 INVCHD 0.00000E+00 *AC227 2.270280E-01 HALFY 2.177335E+01 ACTCNV 7.232300E+04 AWT 1.00000E+02 INVCHD 0.00000E+00 EPAREL, *AC228 AWT 2.280310E-01 HALFY 6.993710E-04 ACTCNV 2.241700E+09 0.00000E+00 INVCHD 0.00000E+00 EPAREL *AM241 4.322347E+02 ACTCNV 3.431200E+03 2.410570E-01 HALFY AWT 1.000000E+02 INVCHD 2.400000E+05 EPAREL *AM243 AWT 2.430610E-01 HALFY 7.380313E+03 ACTCNV 1.992900E+02 1.000000E+02 0.00000E+00 EPAREL. INVCHD *AT217 HALFY 1.023547E-09 1.609500E+15 2.170050E-01 ACTCNV AWT 0.00000E+00 0.00000E+00 INVCHD EPAREL *BA137M 1.369070E-01 HALFY 4.851550E-06 5.382400E+11 ACTCNV AWT 0.00000E+00 0.00000E+00 INVCHD EPAREL *BI211 4.049824E-06 4.184000E+11 2.109870E-01 HALFY ACTCNV AWT 0.00000E+00 INVCHD 0.00000E+00 EPAREL *BI212 2.119910E-01 HALFY 1.151253E-04 ACTCNV 1.464800E+10 AWT 0.00000E+00 0.00000E+00 INVCHD EPAREL *BI213 HALFY 8.679553E-05 ACTCNV 1.933800E+10 2.129940E-01 AWT 0.00000E+00 INVCHD 0.00000E+00 EPAREL *BI214 2.139990E-01 HALFY 3.783638E-05 ACTCNV 4.415300E+10 AWT 0.00000E+00 INVCHD 0.00000E+00 EPAREL *CF252 AWT 2.520820E-01 HALFY 2.638090E+00 ACTCNV 5.375899E+05 0.00000E+00 INVCHD 2.960000E+03 EPAREL *CM243 2.430610E-01 2.850088E+01 ACTCNV 5.160700E+04 HALFY AWT 0.00000E+00 1.00000E+02INVCHD EPAREL *CM244 ACTCNV 8.088299E+04 AWT 2.440630E-01 HALFY 1.811013E+01 0.00000E+00 0.00000E+00 INVCHD EPAREL *CM245 8.498927E+03 ACTCNV 1.716500E+02 2.450650E-01 HALFY AWT 1.00000E+02 INVCHD 0.00000E+00 EPAREL *CM248 4.250200E+00 ACTCNV 2.480720E-01 HALFY 3.390698E+05 AWT 1.00000E+02 INVCHD 0.00000E+00 EPAREL *CS137

Information Only

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AWT EPAREL	1.369070E-01 1.000000E+03	HALFY INVCHD	2.999975E+01 2.070000E+04	ACTCNV	8.704300E+04
*FR221					
AWT EPAREL	2.210140E-01 0.000000E+00	HALFY INVCHD	9.126364E-06 0.000000E+00	ACTCNV	1.772400E+11
*ND143	0.00000100	INVEID	0.0000005+00		
AWT	1.430000E-01	HALFY	3.168876E+30	ACTCNV	0.00000E+00
EPAREL	1.000000E+03	INVCHD	0.00000E+00		1
*NP237 AWT	2.370480E-01	HALFY	2.139942E+06	ACTCNV	7.047600E-01
EPAREL	1.000000E+02	INVCHD	6.670000E+01		
*NP239					
AWT	2.390530E-01	HALFY	6.448664E-03 0.000000E+00	ACTCNV	2.319100E+08
EPAREL *PA231	0.000000E+00	INVCHD	0.00000E+00		
AWT	2.310360E-01	HALFY	3.276618E+04	ACTCNV	4.722500E+01
EPAREL	1.000000E+02	INVCHD	0.00000E+00		
*PA233 AWT	2.330400E-01	HALFY	7.392988E-02	ACTCNV	2.075100E+07
EPAREL	0.000000E+00	INVCHD	0.000000E+00	ACICNV	2.0/31008+0/
*PA234M			•		
AWT	2.340430E-01	HALFY	2.224551E-06	ACTCNV	6.866600E+11
EPAREL *PB209	0.00000E+00	INVCHD	0.00000E+00		,
AWT	2.089810E-01	HALFY	3.764625E-04	ACTCNV	4.544100E+09
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
* PB210	2.099840E-01	HALFY	2.229938E+01	ACTCNV	7.634800E+04
AWT EPAREL	1.000000E+02	INVCHD	0.000000E+00	ACICNV	7.034000£+04
*PB211	2.0000002.02				
AWT	2.109890E-01	HALFY	6.863786E-05	ACTCNV	2.468600E+10
EPAREL *PB212	0.00000E+00	INVCHD	0.00000E+00		
AWT	2.119920E-01	HALFY	1.213680E-03	ACTCNV	1.389500E+09
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
* PB214	2.140000E-01	HALFY	5.095553E-05	ACTCNV	3.278500E+10
AWT EPAREL	2.140000E-01 0.000000E+00	INVCHD	0.000000E+00	ACICINV	2.2/0200E+10
*PM147					
AWT	1.469150E-01	HALFY	2.623513E+00	ACTCNV	9.275300E+05
EPAREL *PO212	0.00000E+00	INVCHD	5.250000E+02		
AWT	2.119890E-01	HALFY	9.506629E-15	ACTCNV	1.773900E+20
EPAREL	0.00000E+00	INVCHD	0.000000E+00		
* PO213 AWT	2.129930E-01	HALFY	1.330928E-13	ACTCNV	1.261100E+19
EPAREL	0.000000E+00	INVCHD	0.000000E+00	ACTOW	1.2011001119
*PO214					
AWT	2.139950E-01	HALFY	5.206464E-12	ACTCNV	3.208700E+17
EPAREL * PO215	0.00000E+00	INVCHD	0.00000E+00		
AWT	2.149990E-01	HALFY	5.640600E-11	ACTCNV	2.947900E+16
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
* PO216	2.160020E-01	HALFY	4.753314E-09	ACTCNV	3.482000E+14
AWT ÉPAREL	2.160020E-01 0.000000E+00	INVCHD	4.753514E-09 0.000000E+00		2.402000E114
*PO218					
AWT	2.180090E-01	HALFY	5.799044E-06	ACTCNV	2.827800E+11
EPAREL *PU238	0.00000E+00	INVCHD	0.00000E+00		

*PU238

AWT EPAREL	2.380500E-01 1.000000E+02	HALFY INVCHD	8.774619E+01 4.240000E+06	ACTCNV	1.711500E+04
*PU239					
AWT	2.390520E-01	HALFY	2.406445E+04	ACTCNV	6.214600E+01
EPAREL	1.000000E+02	INVCHD	3.920000E+05		
*PU240	1.000000102	INVCIID	3.920000103		
	2 4005400 01	113 T 132	C E272020-02		2 2781005.02
AWT	2.400540E-01	HALFY	6.537392E+03	ACTCNV	2.278100E+02
EPAREL	1.000000E+02	INVCHD	6.930000E+04		ł
*PU241					
AWT	2.410570E-01	HALFY	1.439937E+01	ACTCNV	1.030000E+05
EPAREL	0.000000E+00	INVCHD	1.930000E+06		
*PU242					
AWT	2.420590E-01	HALFY	3.869198E+05	ACTCNV	3.817100E+00
EPAREL	1.000000E+02	INVCHD	4.910000E+04		
*PU244	1.000000102	INVCIID	4.9100000104		
	2 4406405 01		0.0010010.07		1 7721005 02
AWT	2.440640E-01	HALFY	8.261261E+07	ACTCNV	1.773100E-02
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*RA223					
AWT	2.230190E-01	HALFY	3.130533E-02	ACTCNV	5.120600E+07
EPAREL	0.00000E+00	INVCHD	0.00000E+00		
*RA224					
AWT	2.240200E-01	HALFY	1.001999E-02	ACTCNV	1.592700E+08
EPAREL	0.000000E+00	INVCHD	0.000000E+00		ı
*RA225					
AWT	2.250240E-01	HALFY	4.052993E-02	ACTCNV	3.919900E+07
EPAREL	0.000000E+00	INVCHD	0.000000E+00	11010111	515255002.07
*RA226	0.0000000000000000000000000000000000000	INVCID	0.000000100		
	2 2602505 01		1.599966E+03	ACTCNV	9.885801E+02
AWT	2.260250E-01	HALFY		ACTUNV	9.885801E+02
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*RA228					
AWT	2.280310E-01	HALFY	6.699955E+00	ACTCNV	2.340000E+05
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*RN219					
AWT	2.190090E-01	HALFY	1.254875E-07	ACTCNV	1.300800E+13
EPAREL	0.000000E+00	INVCHD	0.00000E+00		· ·
*RN220					
AWT	2.200110E-01	HALFY	1.761895E-06	ACTCNV	9.222600E+11
EPAREL	0.000000E+00	INVCHD	0.000000E+00		
*RN222	0.0000002.00				
AWT	2.220180E-01	HALFY	1.046997E-02	ACTCNV	1.538000E+08
EPAREL	0.000000E+00	INVCHD	0.000000E+00	11010111	1.0000001 00
	0.0000000000000000000000000000000000000	INVCID	0.000000100		
*SM147	1.470000E-01		1.059355E+11	ACTCNV	2.295400E-05
AWT		HALFY		ACICINV	Z.Z95400E-05
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*SR90					1 2654005.05
AWT	8.990801E-02	HALFY	2.912197E+01	ACTCNV	1.365400E+05
EPAREL	1.000000E+03	INVCHD	9.850002E+03		
*TH227					
AWT	2.270280E-01	HALFY	5.124073E-02	ACTCNV	3.073200E+07
EPAREL	0.000000E+00	INVCHD	0.00000E+00		· ·
*TH228					
AWT	2.280290E-01	HALFY	1.913051E+00	ACTCNV	8.195301E+05
EPAREL	0.000000E+00	INVCHD	0.00000E+00		
*TH229					
AWT	2.290320E-01	HALFY	7.339118E+03	ACTCNV	2.126900E+02
EPAREL	1.000000E+02	INVCHD	0.000000E+00		
*TH230	1.0000000000	TIAACUD	0.0000000000000000000000000000000000000		
	2 2002205 01	HALFY	7.700370E+04	ACTCNV	2.018300E+01
AWT	2.300330E-01			ACICINV	7.010J00E+01
EPAREL	1.000000E+01	INVCHD	0.000000E+00		
*TH231					

		T_INP_CH_REPOSITO		
DIOISOTOPE_INPU	D.I.			
1.000000E+03	INVCHD	0.000000E+00		
9.000000E-02	HALFY	3.168876E+30	ACTCNV	0.000000E+00
0.0000000000000000000000000000000000000	INVCHD	0.0000000000000000000000000000000000000		
			ACTCNV	5.446300E+08
				5 4462000 00
1.000000E+02	INVCHD	2.010000E-01		
2.380510E-01	HALFY	4.468115E+09	ACTCNV	3.361100E-04
1.0000000102	TIMETID	0.0000000000000000000000000000000000000		
			VCICIAA	0.4000000-02
2 2604600 01	UNTEV	2 2/1/02 5 . 07	<u>λ</u> .Ο.Φ.Ο.Ν.Τ.Ζ	6.468300E-02
1.000000E+02	INVCHD	1.150000E+00		
2.350440E-01	HALFY	7.038074E+08	ACTCNV	2.161100E-03
1.000000E+02	INVCHD	0.00000E+00		
2.340410E-01	HALFY	2.445105E+05	ACTCNV	6.247300E+00
1.0000001.02	1100 CIID	2.0200002/00		
			110 1 014 4	5.0705008+00
2 330400E-01	HALFY	1 585072E+05	ACTCNN	9.678300E+00
0.00000E+00	INVCHD	0.0000000000000000000000000000000000000		
			ACTCNV	1.904500E+11
2.340440E-01 0.000000E+00	HALFY INVCHD	6.597601E-02 0.000000E+00	ACTCNV	2.315200E+07
1.0000001.01	INVEID			
			ACTCNV	1.096500E-04
0.00000E+00	INVCHD	0.000000E+00		
	2.320380E-01 1.00000E+01 2.340440E-01 0.000000E+00 2.069770E-01 0.000000E+00 2.330400E-01 1.000000E+02 2.340410E-01 1.000000E+02 2.350440E-01 1.000000E+02 2.360460E-01 1.000000E+02 2.380510E-01 1.000000E+02 8.990701E-02 0.000000E+03	2.320380E-01 HALFY 1.000000E+01 INVCHD 2.340440E-01 HALFY 0.00000E+00 INVCHD 2.069770E-01 HALFY 0.00000E+00 INVCHD 2.330400E-01 HALFY 1.000000E+02 INVCHD 2.330400E-01 HALFY 1.000000E+02 INVCHD 2.340410E-01 HALFY 1.000000E+02 INVCHD 2.350440E-01 HALFY 1.000000E+02 INVCHD 2.360460E-01 HALFY 1.000000E+02 INVCHD 2.380510E-01 HALFY 1.000000E+02 INVCHD 8.990701E-02 HALFY 0.00000E+00 INVCHD 9.000000E-02 HALFY	2.320380E-01 1.000000E+01HALFY INVCHD1.405080E+10 6.110000E-012.340440E-01 0.00000E+00HALFY INVCHD6.597601E-02 0.000000E+002.069770E-01 0.00000E+00HALFY INVCHD0.00000E+002.330400E-01 1.00000E+02HALFY INVCHD1.585072E+05 1.310000E+032.340410E-01 1.00000E+02HALFY INVCHD1.310000E+032.350440E-01 1.00000E+02HALFY INVCHD2.445105E+05 0.00000E+002.350440E-01 1.00000E+02HALFY INVCHD7.038074E+08 1.50000E+002.360460E-01 1.00000E+02HALFY INVCHD2.341483E+07 0.00000E+002.380510E-01 1.00000E+02HALFY INVCHD2.010000E+002.380510E-01 1.00000E+02HALFY INVCHD3.168876E+30 0.00000E+009.000000E-02 1.00000E+03HALFY INVCHD3.168876E+30 0.00000E+00	2.320380E-01 HALFY 1.405080E+10 ACTCNV 1.000000E+01 INVCHD 6.110000E-01 ACTCNV 2.340440E-01 HALFY 6.597601E-02 ACTCNV 0.000000E+00 INVCHD 0.000000E+00 ACTCNV 2.069770E-01 HALFY 9.069325E-06 ACTCNV 0.000000E+00 INVCHD 0.000000E+00 ACTCNV 2.330400E-01 HALFY 1.585072E+05 ACTCNV 1.000000E+02 INVCHD 1.310000E+03 ACTCNV 2.340410E-01 HALFY 2.445105E+05 ACTCNV 1.000000E+02 INVCHD 0.000000E+00 ACTCNV 1.000000E+02 INVCHD 1.150000E+00 ACTCNV 2.350440E-01 HALFY 7.038074E+08 ACTCNV 1.000000E+02 INVCHD 1.000000E+00 ACTCNV 2.360460E-01 HALFY 2.341483E+07 ACTCNV 1.000000E+02 INVCHD 0.000000E+00 ACTCNV 2.380510E-01 HALFY 4.468115E+09 ACTCNV 1.000000E+02 INVCHD 0.000000E+00 ACTCNV

The input file listed above was designed specifically to apply to WIPP-site calculations. In principle, CUTTINGS_S could be applied to other sites, although no such applications are foreseen for regulatory calculations. As a convenience to users, the list below specifies those parameters that could be changed to reflect sites other than the WIPP. It is not to be regarded as part of the code's QA documentation.

!Repository Geometry

VARIable_name <default value=""> Only capitalized portion of name is or required VARI MATERIAL_NAME:PROPERITY_NAME fetch input from CDB file</default>
!Repository Data/Input
INV_ARea <no default=""> - Area of repository according to the WIPP TRU waste type (CH or RH)</no>
[REAL]. INV_AR MATERIAL_NAME:PROPERITY_NAME fetch input from CDB file
OUT_MAT <no default=""> - The CAMDAT material region to which CUTTINGS_S output variables are written (CHARACTER*8).</no>
RADWAS <contact_handled> - Type of radioisotope waste form (CHARACTER). Valid reposes are:</contact_handled>
REMote_handled; WIPP Remote-Handled (RH) TRU waste, CONtact_handled; WIPP Contact-Handled (CH) TRU waste,
or INEL_winco; INEL/WINCO HLW/SNF waste(s).
CHAIN <no default=""> rad_name_1, rad_name_2, rad_name_3, <</no>
A list of all the radioisotope names in a decay chain. The order of the list is from parent to daughter,, terminating at a stable element. Repeat the CHAIN keyword and list of radioisotopes for as many chains as desired. The half-lives, atomic weights, activity conversions, EPA release limits, and waste-unit information will be retrieved from the generic radioisotope data base. The inventory (in Curies) of each radioisotope will be extracted from the secondary database (CHARACTER).
SAVE <all radioisotopes="" unique=""> rad_name_1, rad_name_2, rad_name_3, <</all>

A list of all unique radioisotope names for which values are to be written to the output CAMDAT and to the debug/diagnostics file. As a default, all unique radioisotope values will be stored/saved to the output CAMDAT and the debug/diagnostics file.

- Name of the repository (CHARACTER). Other valid REP_NAME <WIPP> responses are: INEL/WINCO. REP_GEOLogy <SALT> - Repository geologic medium (CHARACTER). Other valid responses are: HALITE, or GRANITE. - Waste Disposal Group number (INTEGER). Valid DISPOSal_group <1> waste-disposal group numbers are from 1 to 5. If NAME=WIPP, and GEOLOGY=SALT, the waste-disposal group number is not used. By specifying a WDG number, radioisotope inventory data is accessed from the Secondary Database according to the WDG number.

Do not specify PACKage_id !!

PACKage_id <1> Waste Package ID (INTEGER). Valid waste package (WP) IDs are from 1 to 26. If NAME=WIPP, and GEOLOGY=SALT, the waste package ID is

not used. By specifying a WP ID, Radioisotope inventory data is accessed from the Secondary Database according to the WP ID.

Do not specify DISPOSal_group !!

END OF APPENDIX I

APPENDIX J: EXAMPLE OF THE INPUT FILE: CUSP_INP\$TXT1

This input file provides various data relevant to the drilling equations and the intrusion time. Because it is designed for a regulatory-type calculation, it does not list specific numerical values for the various parameters. Rather, it identifies them by name (see Appendix C, item 4, for an explanation of input-file formats). Numerical values are then copied over from the input CDB files. This file is discussed in Section 6.1 above as file number 3.

Definitions of the various parameters used in this input file are given at the end of Appendix J.

BEGINNING OF A PRODUCTION - LIKE INPUT FILE ______ 1 Intrusion 100.0 TINTR PARTDIA BLOWOUT: PARTDIA ! Properities TAUFAIL BOREHOLE: TAUFAIL DIAMMOD BOREHOLE: DIAMMOD DOMEGA BOREHOLE: DOMEGA DNSFLUID DRILLMUD: DNSFLUID VISCO DRILLMUD:VISCO YLDSTRSS DRILLMUD: YLDSTRSS ABSROUGH WAS_AREA: ABSROUGH ! Bragflow Multiple hits (max of 10, 0 thru 9) MHIT_0 is associated with the hit that 1 CUTTINGS used for BRAG flow properities t ţ INTR_0 CAVITY 1 < 625 < INTR_1 619 622 INTR_2 618 621 624 < 617 620 623 < < INTR_3 INTR_4 1010 < INTR_5 1023 < INTR_6 471 < INTR_7 606 ! Spallings MODEL3 ! Properties for Model 3 PTHRESH WAS_AREA: PTHRESH VOLSPALL WAS_AREA:VOLSPALL REPPRES CAVITY_1 < OR (MODEL 4) MODEL4

CUTTINGS_S, Version: 5.10 User's Manual		ERMS# 5323 October 20 Page 1	03
REPPR_NM PRESGAS REPPRES CAVITY_1 < RNDSPALL DUM5:RNDSPAL PTHRESH DUM5:PTH	L	· · · · · · · · · · · · · · · · · · ·	
a,			
END	OF	PRODUCTION - LIKE INPUT FILE	

END OF APPENDIX J

APPENDIX K: EXAMPLE OF THE INPUT FILE: CUSP_INP\$TXT1 FOR A TEST RUN

This input file provides various data relevant to the drilling equations and the intrusion time. Because it is designed for a test run, specific numerical values are defined for each of the various parameters. The file is stored in the CMS portion of the WIPP Alpha cluster under the name: CUSP_TEST_2.INP. Corresponding input files for the other test problems may be found by replacing the number 2 in the file's name by the problem number of interest. This file is discussed in Section 6.1 above as file number 3.

Definitions of the various parameters used in this input file are given at the end of this appendix.

			OF		 mp.cm		INPUT	 FTI P	-
		BEGINNING		A	TEST -		INPUT	FILE	_
! ! !		2, STUCK PIPE							
!	INTRUSION								
	TINTR DIAMMOD PRESSURE TCLOUT	2000.0 0.3556 12.5E6 6.40E4							
! ! !	Initial poros	sity is required,	hence	two en	tries				
	POROSITY	0.88 0.07	0405						
!	PROPERTIES								
	DOMEGA 0 VISCO 0	.1210E+04 .7800E+01 .9170E-02 .4400E+01						•	:
	ABSROUGH 0	.2500E-01 .1120E+00							
_									-
		END OF		A T	EST - L:	ike ini	PUT FILI	8	

Explanation of Input Parameters Used in Appendices J and K

In the list that follows, two entries are made for each listing. The first applies to a test-like input file such as the one given in Appendix E. The second applies to a regulatory-calculation-like input file such as the one given in Appendix D. Both pertain to drilling and intrusion parameters.

VARIable or VARI	_name <default values=""> Only capitalized portion of name is required MATERIAL_NAME:PROPERITY_NAME fetch input from CDB file</default>
Drilling Par	ameters
DENSity	<1.210E+03> - Density of the drilling fluid (mud) in kg/m^3 .
DENS	MATERIAL_NAME: PROPERITY_NAME
DOMEga	<0.7800E+01> - Angular velocity of the drill string in rad/s.
or DOMÉ	MATERIAL_NAME: PROPERITY_NAME
ROUGH or ROUGH	<.0250> - Absolute borehole roughness (m) MATERIAL_NAME:PROPERITY_NAME
TAUFail or	<0.1000E+03> - Effective shear resistance to erosion (Pa).
TAUF	MATERIAL_NAME: PROPERITY_NAME
VISCo or	<0.9170E-02> - Plastic viscosity of drilling mud associated with an Oldroyd type fluid model (Pa-s)
VISC	(zero shear-rate viscosity). MATERIAL_NAME:PROPERITY_NAME
YLDSTrss or	<0.4400E+01> - Yield Stress of the drilling fluid (mud) (Pa).
YLDS	MATERIAL_NAME: PROPERITY_NAME

Intrusion Parameters VARIable_name <default value> Only capitalized portion of name is required or MATERIAL_NAME: PROPERITY_NAME fetch input from CDB file VARI if nhits .gt. 1 (testing mode), then or VARI variable(1) variable(2) ... variable(nhits) UNITS <SEC> - Type of time units that the times of intrusions UNITS are specified [CHARACTER]. Valid responses are: YEARS or SECONDS. <1> - Number of intrusions [REAL]*. NHITS NHITS

^{* [}REAL] signifies a real, floating-point number.

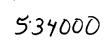
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```
- Drilling bit diameter (m); (outside
                   <0.5080E+00>
       DBDTAM
                                    diameter of drill-string annulus)
                                    corresponding to the first
        or
                                     intrusion time [REAL].
                  MATERIAL_NAME: PROPERITY_NAME
       DBDIAM
        or
       DBDIAM(nhits)
       TINTR
                  <> - First intrusion time (m)
                                                     [REAL].
        or
       TINTR
                  MATERIAL_NAME: PROPERITY_NAME
        or
       TINTR(nhits)
                <> - Clean-out time** (s) [REAL].
       TCLOUT
        or
       TCLOUT
                  MATERIAL_NAME: PROPERITY_NAME
        or
       TCLOUT(nhits)
                 <> - Particle diameter (m) [REAL].
       PARTDIA
        or
       PARTDIA(nhits)
                The following two parameters are NOT required if a BRAGFLO CDB file
Note!!!
is attached.
                  <> - Porosity at time of intrusion (unitless) [REAL].
       POROSITY
        or
                  MATERIAL_NAME: PROPERITY_NAME
       POROSITY
        or
       POROSITY (nhits)
                  <> - Repository pressure at time of intrusion (Pa) [REAL].
       PRESSure
        or
       PRESS
                  MATERIAL_NAME: PROPERITY_NAME
        or
       PRESS(nhits)
```

Note!!! The preceding two parameters are NOT required if a BRAGFLO CDB file is attached.

END OF APPENDIX K

^{**} Clean-out time is the time required to replace all of the borehole drilling fluid at the present pumping rate.



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Jennifer Long SCM Coordinator



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Phone: (505) 234-0106 Fax: (505) 234-0061 Internet: jjlong@sandia.gov

Date: February 25, 2004

To: SNL WIPP Records Center Carlsbad Programs Group

From: Jennifer Long SCM Coordinator

Subject: Record Correction to User's Manual Criteria Form and Cover Sheet for CUTTINGS_S Version 5.10

Please disregard the User's Manual Criteria Form and Cover Page for CUTTINGS_S Version 5.10 and replace with the attached User's Manual Criteria Form and Cover Page. The originals incorrectly listed the Document Version as 1.00. The correct Document Version is 5.10.

Cc: C. Hansen

WIPP:1.5.1:SFT:QA-L:532334

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				Revision 10 Page 25 of 34
	Appendix G			
NUCLEAR WASTE MANAGEMENT Sandia National Laboratories	User's Manual Criteria For	n	Ν	m Number: IP 19-1-6 age 1 of 1
Does the user's man 1. Software Name: 2. Software Versio 3. Document Versi	n: <u>5.10</u>			
4. ERMS #: Prior to sign-off of the l	532340	code spo	onsor s	o that "Yes"
or " N/A " may be check 5. A statement(s) o	ed. Include this form as part of the User's Manual. f functional requirements (consistent RD) and system limitations?		Yes	
6. An explanation o	f the mathematical model and s, where applicable as based on code functionality?	\boxtimes	Yes	□ N/A
7. Physical and ma based on code fu	thematical assumptions, where applicable as unctionality?	\boxtimes	Yes	🗋 N/A
8. The capabilities	and limitations inherent in the software?	\boxtimes	Yes	
Instructions that software?	describe the user's interaction with the	\boxtimes	'Yes	
10. The identification	n of input parameters, formats, and valid ranges?	\boxtimes	Yes	,
11. Messages initiate the user can resp	ed as a result of improper input and how bond?	\boxtimes	Yes	
12. The identification	n and description of output specifications and formats?	\boxtimes	Yes	
13. A description of	any required training necessary to use the software?	\boxtimes	Yes	
14. The identification	n of components of the code that were not tested?	\boxtimes	Yes	
	Sponsor (print) Signature	vr		10/22/03 Date
	eviewer (print) Signature			<u>10/22/03</u> Date
	Manager (print) Signature	~		<u>10/22/03</u> Date
18. Jennifer Lona SCM Coo	rdinator (print) Signature	>		<u>10/22/03</u> Date

NP 19-1

Key for check boxes above:

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Software Requirements

Check Yes for each item reviewed and found acceptable

Check N/A for items not applicable, where applicable as based on code functionality

WIPP PA

User's Manual

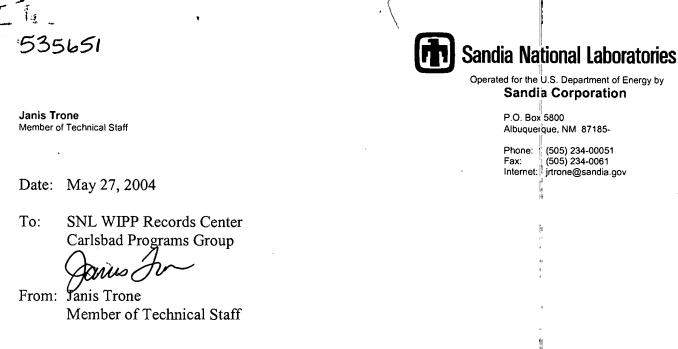
for

CUTTINGS_S, Version 5.10

DOCUMENT VERSION 5.10 .

ERMS # 532340

OCTOBER 2003



Subject: Record Correction to User's Manual for CUTTINGS_S Version 5.10

Please replace pages 8-10, 14-15, 17 - 19 and 152-153 of the User's Manual for CUTTINGS_S Version 5.10 with the attached pages. The attached pages contain change bars beside text which was revised, rather than the hand edits in red ink found in the original pages.

This page change correction is in response to Finding 1 of the February 24-26, 2004 EPA Audit of SNL.

Cc: C. Hansen J. Long

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WIPP:1.5.1:SFT:QA-L:532334

Exceptional Service In the National Interest

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4.0 DESCRIPTION OF THE MODEL AND METHODS

The discussion of the model and methods given below is but a brief overview c¹f the theory and methods employed by CUTTINGS_S. It is intended for users who have only a cursory interest in the mathematical and fluid dynamical details. A thorough, detailed treatment of the theory and numerics applied in CUTTINGS_S is given in Appendix A of this manual.

4.1 Description of the Model

<u>CUTTINGS</u> S is a multi-faceted computational procedure to assess the effects of *direct removal* of wastes buried in a WIPP-like nuclear-waste repository. Direct removal is hypothesized to occur as the result of inadvertent penetration of the repository by a borehole drilled in connection with oil or gas exploration at some unknown time in the future (Berglund, 1993). The word "direct" refers to the fact that CUTTINGS S removals from the repository to the surface occur *at the time* of drilling. Since drilling operations are normally completed on the time scale of weeks, removal of drilled wastes from the repository to the surface (i.e., the accessible environment) takes place entirely within a single timestep of the much-longer-term groundwater transport codes. Hence, the word "direct" should be taken to mean "from the repository to the surface within a single timestep of a typical WIPP groundwater code."

<u>Removal is subdivided according to process, which results in three principal categories, cuttings, cavings, and spallings.</u>

First, the borehole is assumed to penetrate directly through waste containers stored within the repository. The drill bit cuts directly through the waste containers, allowing those repository wastes located on a collision course with the drill bit (called cuttings) to be transported directly and immediately to the surface via the cooling fluid (muddy brine) that is circulated through the borehole during drilling. The volume of repository material removed as cuttings is the portion of the deformed repository cut by the borehole.

Wastes that originally bordered the borehole on its exterior are permitted to erode into the cooling fluid as a result of the fluid's frictional effects on the outer cylindrical surface of the borehole. If shear stresses in the cooling fluid at the wall exceed the shear strength of the matrixed materials located at the wall, erosion will occur and wall material (cavings) will be removed*. As a result, the borehole diameter may increase locally at the level of the repository. Because of that action, the volume of wastes removed from the repository can actually be larger than the volume of the borehole originally cut through the wastes. As the diameter of the borehole increases, the fluid shear stresses at the outer boundary decrease. If they decrease enough that the failure stress of the wall material equals or exceeds the fluid shear stress at the outer boundary of the flow region, then the caving process will terminate. CUT TINGS_S allows for both laminar and turbulent shear flow in the circulating fluids, which, in turn, leads to two

^{*} In tightly matrixed or crystaline media such as the WIPP-site halite formation, exploratory boreholes are usually unlined. Rather, the drilling fluid is changed to saturated salt solution to minimize dissolution at the walls, and the borehole is permitted to support itself until the drill bit reaches granular strata below the halite horizon.

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different models to assess the effects of material removal due to erosion. If the cooling-fluid flow is laminar, CUTTINGS_S employs an analytical solution of the modeling equations to evaluate wall stresses and other flow parameters. The analytical solution takes the form of nonlinear integral equations, which CUTTINGS_S evaluates via an iterative numerical procedure. If the flow is turbulent, CUTTINGS_S turns to a simpler empirical model to evaluate the turbulent shear stresses. Transition to turbulence is assumed to occur at a Reynolds number greater than 2100. In nature, the transition to turbulence actually occurs over a range of Reynolds numbers and is not an abrupt process. Normally there is a transitional regime during which the flow is partly laminar and partly turbulent with active and chaotic transitions between the two states. CUTTINGS_S does not attempt to model the transition regime. Rather, it assumes the flow becomes fully and abruptly turbulent when the Reynolds exceeds 2100. CUTTINGS_S employs a corrective procedure to compensate for the effects of that assumption (see Appendix A for details). To date, all WIPP-like runs have required turbulent cooling-fluid flow.

The cuttings/cavings model described thus far is functionally similar to the CUTTINGS model used in previous WIPP PAs (WIPP PA Dept, 1992, Vol 1). However, an important new element has been added to the newer CUTTINGS_S code, namely, the inclusion of a third materialremoval mechanism, removal in the form of spallings, which is discussed below.

If the repository is pressurized at the time of penetration, and if the pressure within the repository is high enough (i.e., above the hydrostatic pressure in the cooling-fluid column), repository materials originally surrounding the borehole, but interior to the repository-fluid interface, can be forced to deform inward toward the borehole so as to replace the materials already removed by action of the cutting bit and the erosive brine. The solid fraction of these materials are called spallings^{*}. Depending on the parameters of the problem, spalling deformations^{*} can be rapid, or they can be a sluggish, more like a large-amplitude, viscoelastic flow. If repository materials are relatively impermeable, the spalled materials can deform slowly inward so as to press against the rotating drill string. If repository pressures are high enough, frictional forces against the drill string may be high enough to stop the drill from turning. In that event, drillers hormally withdraw the drill (vertically) and redrill downward through the spalled materials that clog the borehole. On occasion, that procedure must be repeated more than once. If repository materials are more permeable, the pressurized gas may simply seep through the matrix carrying loose particles with it. Alternatively, if repository pressures are too low, the spallings part of CUTTINGS S may do nothing at all. CUTTINGS S may have to be applied more than once during an assessment run, at more than one intrusion time, since multiple borehole intrusions are permissible in WIPP PAs.

The total volume of repository material removed by CUTTINGS S is the sum over all intrusions of cuttings, cavings, and spallings. Options cuttings and cavings are assumed to operate

^{*} If the repository is wet with brine at time of penetration, which possibility is allowed in WIPP PAs, three phases of material could, in principle, flow toward the borehole, namely: pressurized gases, pressurized brine, and solid wastes. CUTTINGS S estimates only the amount of solids that are moved into the borehole. Pressurized brine that may be released into the borehole is treated by an independent WIPP code named DBR_BRACFLO. The gas flow, although crucial to the overall deformational process, transports no radioisotopes and is therfore not assessed.

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continuously, whereas spallings operates only when repository pressure conditions permit it (the cut off being 8MPa.). The amount of material removed by each process, during each hypothesized drilling intrusion is reported in CUTTINGS S's output in terms of an equivalent area. Downstream WIPP codes, namely CCDFGF, use BRAGFLO's repository-height data to convert the equivalent areas to volumes of repository material, and then assign various sampled aged repository contents to convert these material volumes to releases. Thus, CCDFGF evaluates the release effects of striking contact-handled versus remote-handled wastes. The reader is referred to Appendix A for details.

The fluid-flow portion of CUTTINGS S is based on a non-linear analytical integral solution for laminar helical flow of a non-Newtonian (Oldroyd-type) fluid in an annular pipe, and on empirical equations for turbulent flow in the same geometrical configuration. CUTTINGS S acquires its parameter data from (i) the input CDB files provided by the WIPP codes MATSET and BRAGFLO, (ii) a user-supplied input text file, and (iii) the carefully controlled and QAed WIPP property database. Discrete values of sampled data are provided by the WIPP sampling code POSTLHS in cases where parameter uncertainty is judged to impact assessments of the repository's performance. CUTTINGS S input includes the rotational speed of the drill string, cutting-bit diameter, the shear-rate-dependent viscosity and density parameters for the drilling mud, borehole roughness, the compacted repository material. A few required parameters, namely drilling-mud flow rate and drill-collar diameter, are calculated internally by the code, based on present-day oil-drilling industry guidelines.

4.2 Description of the Modeling Methods

CUTTINGS_S does not endeavor to integrate conservation equations characterizing the flow on a point-for-point, time-for-time basis in search of a detailed space-time history of the removal process. Rather, removal is treated as instantaneous, that is, within a single timestep of the longer-term groundwater codes. Thus, numerical methods are used primarily to determine the limiting values of the physical parameters that represent the final configurations of the various (sampled) drilled systems, which is all that is required to assess radioactive waste removal due to the cuttings, cavings, and spallings that arise during drilling operations. These methods are fairly straightforward, iterative, recursive approximation techniques that involve systems of algebraic equations. They are discussed thoroughly in Appendix A (Sections 2.2.1.1, 2.2.2.1, and 2.4.1.2.1), which is highly recommended to readers who aim to understand this code.

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where XY:[ABC.DEF.GHI]JKL.MNO represents the specific input file the user wishes to associate with the input logical CUSP_INP\$CDB. The path need be included only if the file resides other then in the local working-level directory.

6.2 CUTTINGS S's Input/Output File Structure

Up to 12 logical symbols may be associated with CUTTINGS S's various input/output files. Of them, 6 are required and 5 are optional in regulatory runs. Before discussing the nature and function of the 12 possible input/output files, we first list them one-by-one, for ready reference, together with their corresponding logical symbols. They are:

File ID No.	Input/Output File Names	Associated Logical Symbol	ls the file Required or Not?
1. Binary in	put CAMDAT file (from MATSET)	CUSP INP\$CDB	Yes
repositor	ntrol file (text) specifying y/model parameters, initial es, generic radioisotope database.	CUSP INP\$TXT0	Yes_
	ntrol file (text) specifying drilling sion parameters	CUSP INP\$TXT1	Yes
4. BRAGFL Regulato Test runs		CUSP INP\$BRAGCDB	<u>Yes</u>
	ntrol file for Spall Model 4 specifing s and their volumes by vector	CUSP SPL4\$DAT	Yes (spall model 4 only)
6. CUTTIN	GS S's binary output .CDB file	CUSP OUT\$CDB	Yes
7. CUTTIN	GS S's output debug text file	CUSP OUT\$DBG	
<u>8. Text file</u>	for verification testing	CUSP OUT\$NVERIFY	<u>No</u>
9. Radioiso output f	tope inventory history (in CURIES)	CUSP OUT\$IHISTO	n No
(EPA 40	ized-release-history output text file CFR191, Subpt B) ASCII format	CUSP OUT\$NHISTO	No
Long	istory plotting text file for: half-life radioisotope decay h half-life radioisotope decay half-life radioisotope decay	CUSP OUT\$PLT1 CUSP OUT\$PLT2 CUSP OUT\$PLT3	<u>No</u> No No
	text file for transfer to PP code CCDFPERM	CUSP OUT\$ICTRN	<u>No</u>

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The first five files of the above list are CUTTINGS S's input data and control files. They are described briefly in Section 7.0 of this user's manual. The text-format input files include specific examples printed out in full in the appendices of this user's manual. Appendix C is included to assist the user in understanding the format used in CUTTINGS S's input control files (file numbers 1 and 2).

The last seven files of the above list are CUTTINGS S's output files. They are described briefly in Section 9.0 of this user's manual. Of the required output files, only the debug file is created in text format. One example of a CUTTINGS S output file is included in Appendix D. It is a debug file. It is annotated with italicized remarks that identify and explain the various groups of data that are reported therein.

6.3 Exercising CUTTINGS_S

Once the logical symbol CUSP has been defined (see Section 6.1), and the necessary input and output files have been properly assigned (see Sections 6.1 and 6.2), CUTTINGS_S may be exercised simply by typing the logical symbol:

\$CUSP

followed by a carriage return. If all goes well, CUTTINGS_S will run and issue a "normal completion" message. If errors are encountered, CUTTINGS_S will abort and issue an error message(s). See Section 8.0 for a listing of the possible error messages CUTTINGS_S can produce.

6.4 Using CUTTINGS_S's PRE_CUTTINGS Mode

Because interactions with the controlled WIPP database require large and unnecessary expenditures of input/output time during production runs, CUTTINGS_S includes a special operational mode designed to expedite data acquisitions during production runs. It is called PRE_CUTTINGS, and in it, the WIPP QAed property database is accessed but once during an entire production run. PRE_CUTTINGS reads a template file that interfaces the WIPP property database, extracting the required values and writing them to an ASCII file that can be read time and again by CUTTINGS_S without subsequent reference to the property database. The ASCII data file thus produced is under strict CMS control.

PRE_CUTTINGS's template file, which is shown in Appendix E, treats three data types. The first type is exemplified by CUTTINGS_S's code parameters, which are discussed in the theoretical Manual (see Appendix A). These values are essential to exercising CUTTINGS_S, but were not entered as part of the WIPP's CCA controlled database. They are discussed in detail in Appendix A. The second type are common property values as taken from the WIPP's CCA controlled database. For example, the datum BOREHOLE:L1 is called. In this example, the colon at mid word identifies the form as a WIPP controlled database property. BOREHOLE is the material identifier and L1 is the parameter identifier. Using these two keywords, the WIPP

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7.0 DESCRIPTION OF CUTTINGS_S'S INPUT FILES

CUTTINGS_S's input files are listed in Section 6.2 together with their corresponding CUTTINGS_S logical symbols. They are file numbers 1 through 4 in that listing.

File 1. CUSP_INP\$CDB is the principal input CDB file and is absolutely required by CUTTINGS_S. It is the file to which CUTTINGS_S's results will be transferred and it is normally provided by an upstream code such as MATSET. In production runs, sampled data from POSTLHS would be fed to this input file. In that case, there would be N input CDB files, one for each of the sampled values where N is the number of samples employed. In test-case mode or debug mode, it is inconvenient to open the editor and hand build an input CDB file for specific input values. However, even in test or debug mode, CUTTINGS_S expects to receive an input CDB file. As a convenience to users, an all-purpose, input CDB file for use with test problems and in debug runs is stored in the configuration-management-system (CMS) portion of the WIPP Alpha cluster. It is named: CUSP_TEST_1.CDB and can be called by typing:

\$LIBCUSP

\$CFE CUSP_TEST_1.CDB

File 2. CUSP_INP\$TXT0 is an ASCII input control file that contains various repository, drill geometry, and material input parameters required by CUTTINGS_S. File 2 also contains the required radioisotope inventory data for the scenario currently under study. In regulatory calculations, these data are taken directly from the controlled WIPP property database. File 2 also provides radioisotope data essential for (1) the Bateman (1910) decay calculation, for (2) calculation of radioactive release values in Curies, and for (3) normalizing calculated releases in terms of the Environmental Protection Agency's (EPA's) 40CFR191 release I mits (WIPP PA Dept [Vol 1], 1992). A sample CUSP_INP\$TXT0 file usable for QA test and/or debug problems is available in the CMS portion of the WIPP Alpha cluster. It is named: CUSP_TEST_REPOSITORY.DAT, and it is printed in full in Appendix I of this user's manual. Most of its parameters are specific to the CUTTINGS_S model and to individual radioisotopic properties.

File 3. CUSP_INP\$TXT1 is an ASCII input control file that contains various drilling and intrusion input parameters required by CUTTINGS_S. A sample CUSP_INP\$TXT1 file usable for test and/or debug runs is available in the CMS portion of the WIPP Alpha cluster. It is named CUSP_TEST_#.INP_where # is the QA test problem number and varies from 1 through 5. In a normal production run, this file would contain no specific input data other than an intrusion time. Instead, only character variables would appear, identifying the element property name on the input CDB file. Values would then be assigned from the secondary database (see item 4, below). Appendix J is sample printout of a typical production input file. For specific testing or debugging, this file would have fixed specific numerical values. Appendix K is a listing of CUSP_TEST_2.INP.

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This file also identifies which spall model will be used. At the end of the file the model is determined by entering MODEL3 or MODEL4, depending on which model will be used. This is followed by the parameters that are needed for the model. If model three is selected then three parameters will be needed. Model four uses three parameters and one dummy parameter. The parameter PTHRESH is not used in model four. An example of how to call each model and the parameters it requires is below.

MODEL3PTHRESHWAS AREA: PTHRESHVOLSPALLWAS AREA: VOLSPALLREPPRESCAVITY 1

MODEL4REPPRNMREPRESCAVITY 1RNDSPALLDUM5:RNDSPALLPTHRESHDUM5:PTH

File 4. CUSP_INP\$BRAGCDB is the output CDB file from BRAGFLO as it is exercised in the undisturbed or disturbed scenarios. CUTTINGS_S extracts from it (1) the repository porosity at time of decommissioning, and (2) the repository porosity and (3) repository pressure at the specified intrusion time. In regulatory calculations, these data would be provided by a regulatory BRAGFLO run. For QA testing and debug purposes, a fixed BRAGFLO output CDB file has been stored in the CMS portion of the WIPP Alpha cluster. It is named CUSP_TEST_INP_BF.CDB, and it is the BRAGFLO output file used for test problem 1.

File 5. CUSP_SPL4\$DAT is a data file that contains information pertinent to spall model 4. The file supplies the code with pressures and their corresponding data to determine the spall volume. The first line of the file lists the number of vectors in pressure data set. The second line is the number of pressure data sets. The third line lists the different pressures in increasing order separated by blank spaces. The following lines list data relating to each pressure. The first data set is for the first pressure given in line three, the second data set is for the second pressure in line three, and so on. The data sets have three columns of data that are separated by blank spaces. The first column is the vector numbers, the second column is the time (not used), and the third column is the volume. Each of the data sets is followed by a blank line. Each data set must contain the same number of vectors and be in the same order. An example data file is shown below.

$ \frac{5}{2} $ 1000000 1 0 2 0 3 0 4 0 5 0 blank)0 2	0000000
1 0	1	
20	1.2	5
3 0	1.5	,
4 0	1.7	5
50	2.0	0
blank		
$\frac{1}{2}$ 0	10	
2 0	12.	5

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3	0	15
4	0	17.5
5	0	20
bla	nk	

8.0 ERROR MESSAGES

Numerous procedural errors can cause CUTTINGS_S to abort. In general, these errors relate to the improper input of data. By noting where the code was when the abort occurred, and what the code was trying to read at the time, these problems can generally be resolved. However, there is not specific advice that can be offered to users in resolving errors beyond the general statement: locate the error and repair it. However, Appendix L, which was extracted primarily from the source code, contains useful operating notes and tips to the user. They are included purely because the code sponsor felt they might be useful to novice users.

To familiarize readers with CUTTINGS_S's error-message patterns, CUTTINGS_S's various error messages are listed below, section for section within the code.

SUBROUTINE ADDMEM(MORMEMR, MORMEMC)

CALL QAABORT(

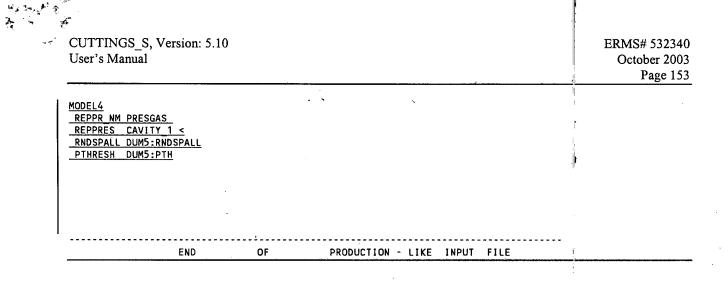
- \$ '(ADDMEM) REAL array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM) REAL array memory errors detected ***') CALL QAABORT(
- \$ '(ADDMEM) CHARACTER array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM) CHARACTER array memory errors detected ***') CALL QAABORT('(ADDMEM1) - REAL array size errors detected ***') CALL QAABORT('(ADDMEM1) - REAL memory errors detected ***') CALL QAABORT(
- \$ '(ADDMEM1) CHARACTER array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM1) CHARACTER array memory errors detected ***') CALL QAABORT('(ADDMEM2) - REAL array size errors detected ***') CALL QAABORT('(ADDMEM2) - REAL memory errors detected ***') CALL QAABORT(
- \$ '(ADDMEM2) CHARACTER array size errors detected ***') CALL QAABORT(
- \$ '(ADDMEM2) CHARACTER array memory errors detected ***') CALL QAABORT('(ADDMEM3) - REAL array size errors detected ***') CALL QAABORT('(ADDMEM3) - REAL memory errors detected ***') CALL QAABORT(

APPENDIX J: EXAMPLE OF THE INPUT FILE: CUSP_INP\$TXT1

This input file provides various data relevant to the drilling equations and the intrusion time. Because it is designed for a regulatory-type calculation, it does not list specific numerical values for the various parameters. Rather, it identifies them by name (see Appendix C, item 4, for an explanation of input-file formats). Numerical values are then copied over from the input CDB files. This file is discussed in Section 6.1 above as file number 3.

Definitions of the various parameters used in this input file are given at the end of Appendix J.

BEGINNING OF A PRODUCTION - LIKE INPUT FILE Intrusion TINTR 100.0 PARTDIA BLOWOUT: PARTDIA ! Properities TAUFAIL BOREHOLE: TAUFAIL DIAMMOD BOREHOLE: DIAMMOD DOMEGA BOREHOLE:DOMEGA DNSFLUID DRILLMUD:DNSFLUID VISCO DRILLMUD:VISCO YLDSTRSS DRILLMUD:YLDSTRSS ABSROUGH WAS_AREA:ABSROUGH Bragflow Multiple hits (max of 10, 0 thru 9) MHIT O is associated with the hit that CUTTINGS used for BRAG flow properities INTR_O CAVITY_1 < INTR 1 619 622 625 < 618 621 624 < INTR 2 623 < 620 INTR 3 617 INTR 4 1010 < INTR 5 1023 < INTR_6 471 < INTR 7 606 < Spallings i MODEL3 Properties for Model 3 PTHRESH WAS AREA: PTHRESH VOLSPALL WAS AREA: VOLSPALL REPPRES CAVITY 1 < OR (MODEL 4) **Information Onl**



END OF APPENDIX J