

**Keywords: DWPF, glass,  
IDMS, durability, Product  
Consistency Test, vitrification,  
process control**

**WSRC-RP-91-371**

**WSRC-RP--91-371**

**DE93 002154**

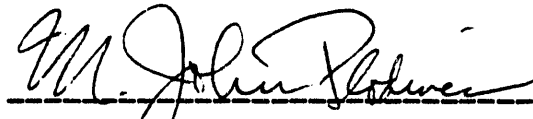
**DURABILITY OF GLASSES FROM THE HG-DOPED  
INTEGRATED DWPF MELTER SYSTEM (IDMS)  
CAMPAIGN (U)**

**CAROL M. JANTZEN**

**Approved by**

**E.W. Holtzscheiter, Research Manager  
Defense Waste Processing Technology**

**Publication Date: August 30, 1992**


  
-----  
**Authorized Derivative Classifier**

---

**Westinghouse Savannah River Co.  
Savannah River Site  
Aiken, SC 29808**

**PREPARED FOR THE U.S. DEPARTMENT OF ENERGY UNDER  
CONTRACT DE-AC09-89SR18035**

**MASTER**

**DISTRIBUTION OF THIS DOCUMENT IS UNLIMITED** 

**This page intentionally left blank.**

## **ABSTRACT**

---

The Integrated DWPF Melter System (IDMS) is designed and constructed to be a 1/9th scale prototype of the full scale Defense Waste Processing Facility (DWPF) melter. The IDMS facility is the first engineering scale melter system capable of processing mercury, and flowsheet levels of halides and noble metals. In order to determine the effects of mercury on the feed preparation process, the off-gas chemistry, glass melting behavior, and glass durability, a three-run mercury (Hg) campaign was conducted. The glasses produced during the Hg campaign were composed of Batch 1 sludge, simulated precipitate hydrolysis aqueous product (PHA) from the Precipitate Hydrolysis Experimental Facility (PHEF), and Frit 202. The glasses were produced using the DWPF process/product models for glass durability, viscosity, and liquidus. The durability model indicated that the glasses would all be more durable than the glass qualified in the DWPF Environmental Assessment (EA). The glass quality was verified by performing the Product Consistency Test (PCT) which was designed for glass durability testing in the DWPF.

The durability of three glasses from each of the three runs was examined with the PCT. There was very little variation in the glass durability during run HG-1 until excess NaOH was added to adjust the glass viscosity to be in the correct processing range. The addition of the excess NaOH caused the glasses in runs HG- 2 and HG-3 to be less durable than the glasses in run HG-1. The nine IDMS Hg glasses were of comparable durability to previously examined sludge-only 165 glasses and to glasses fabricated from Frit 202 with excess alkali. All the IDMS Hg glasses were over 20 times more durable than the glass that was qualified in the DWPF Environmental Assessment (EA). The IDMS Hg campaign glasses would, therefore, meet current Waste Acceptance Preliminary Specification (WAPS) criteria for product consistency.

**This page intentionally left blank.**

## **Table of Contents**

---

	<b><u>Page</u></b>
Abstract . . . . .	3
Introduction . . . . .	7
Experimental . . . . .	9
Glass Homogeneity . . . . .	9
Glass Composition . . . . .	9
Glass Durability . . . . .	9
Quality Assurance . . . . .	10
Glass Sample Identification . . . . .	10
Results and Discussion . . . . .	11
Glass Homogeneity . . . . .	11
Glass Composition . . . . .	11
Glass Durability . . . . .	13
Conclusions . . . . .	17
References . . . . .	19
Appendix I : Measured Glass Compositions . . . . .	21
Appendix II: Leachate Multielement Standard Analyses . . . . .	39
Appendix III: Leachate Data for the IDMS Hg Campaigns 1, 2, 3	41

**This page intentionally left blank.**

# **DURABILITY OF GLASSES FROM THE HG-DOPED INTEGRATED DWPF MELTER SYSTEM (IDMS) CAMPAIGN (U)**

## **INTRODUCTION**

Liquid high-level nuclear waste will be immobilized at the Savannah River Site (SRS) by vitrification into borosilicate glass. The production process to be used in the Defense Waste Processing Facility (DWPF) is designed to reliably produce durable borosilicate nuclear waste glasses.

The Integrated DWPF Melter System (IDMS) was designed and constructed to be a 1/9th scale prototypic pilot plant of the full scale DWPF melter. The IDMS facility is the first engineering scale melter system operated at SRS that is capable of processing mercury, and flowsheet levels of halides and noble metals. The mercury, present in the waste due to the use of mercuric nitrate as a catalyst during nitric acid dissolution of spent fuel rods, decomposes at temperatures much less than the borosilicate glass melt temperature. Any mercury present in the waste at the time of vitrification will likely exit the melter in the off-gas. Therefore, the DWPF has planned to remove most of the mercury prior to vitrification by reduction and steam stripping.<sup>1</sup>

In order to determine the effects of mercury on the feed preparation process, the off-gas chemistry, glass melting behavior, and glass durability, a mercury (Hg) campaign was conducted in the IDMS. The IDMS was operated batch-style for three SRAT/SME (Slurry Receipt Adjustment Tank/Slurry Mix Evaporator) cycles. The facility was operated close to DWPF flowsheet conditions. The process/product models<sup>2</sup> developed for DWPF were used in spreadsheet format since the DWPF process control system, the Product Composition Control System (PCCS)<sup>3</sup> had not yet been fully implemented. Mercury containing Batch 1 sludge<sup>4</sup> without noble metals was blended with precipitate hydrolysis aqueous (PHA) product and Frit 202. This PHA glass was designed to have similar process and product characteristics to SRS sludge-only glasses which were formulated without PHA.<sup>5</sup>

A durability test, designated the Product Consistency Test (PCT), has been developed to measure the consistency of the glass product produced in the Defense Waste Processing Facility (DWPF) during routine production.<sup>6-8</sup> The test was designed to meet the requirements of the Waste Acceptance Preliminary Specifications (WAPS) 1.3.<sup>9</sup> Currently this specification states that based on comparative PCT analyses, a DWPF glass must have a release which is better than the glass that was qualified in the DWPF Environmental Assessment (EA).<sup>10</sup> The PCT measurement of glass durability will take a minimum of 7 days.<sup>6-8</sup> Therefore, a product quality model based on hydration free energy was developed.<sup>11-17</sup> The use of the hydration thermodynamic model allows the glass quality to be predicted from the composition of vitrified melter feed, and then verified after production.

This study documents the final glass chemistry and the measured glass durability of glasses produced during the IDMS-Hg campaign. Over 15,000 pounds of glass was poured during the entire campaign. This represented about 5000 pounds of glass for each of the three runs (Hg-1, Hg-2 and Hg-3). Three samples of IDMS glass were taken from the beginning of each of the three runs (Hg 1-1, Hg 2-1, Hg 3-1), at the middle of each run (Hg 1-2, Hg 2-2 and Hg 3-2), and at the end of each run (Hg 1-3, Hg 2-3, Hg 3-3). This represented glass pouring intervals of ~0, 2500, and 5000 pounds of glass. The process/product properties calculated from the vitrified slurry are compared to the properties calculated from the final glass product as a function of run sequence. The measured glass durability is reported and compared to the following:

- the predicted glass durability
- previously tested SRS sludge-only and PHA glasses
- the sludge-only glass qualified in the DWPF Environmental Assessment.



## **EXPERIMENTAL**

### **Glass Homogeneity**

X-ray Diffraction (XRD) and Scanning Electron Microscopy (SEM) analyses coupled with Energy Dispersive Analysis by X-ray (EDAX) were performed on the nine IDMS-Hg campaign glasses in order to determine the crystallinity and/or homogeneity of the glass.

### **Glass Composition**

The elemental composition of the nine glasses was originally measured by the TNX laboratory. The glasses were reanalyzed by the Analytic Development Section of the Savannah River Technology Center because the original analyses were inconsistent. The nine glasses were analyzed for whole element chemistry and redox. The glass samples were analyzed by the following techniques:

- Dissolution by  $\text{Na}_2\text{O}_2$  with an HCl uptake
  - ICP for Al, Ca, Fe, Mg, Mn, Li, Si, Cr, B, Sr, Ti, P, Ba, Pb, Mo, Zn, Cu, Ni, La, Cd
- Dissolution by HCl/HF microwave
  - ICP for Na, Zr
  - AA for Na, K
- Dissolution by  $\text{H}_2\text{SO}_4/\text{HF}$  in the presence of  $\text{NH}_4\text{VO}_3$ <sup>18</sup>
  - Colorimetric for  $\text{Fe}^{2+}$  and  $\text{Fe}(\text{Total})$

where ICP is Inductively Coupled Plasma Spectroscopy and AA is Atomic Absorption analysis.

### **Glass Durability**

The durability of the nine glasses was studied using Version 3.0 of the PCT.<sup>8</sup> In the PCT, crushed glass of 100-200 mesh is immersed in ASTM Type I water for 7 days at 90°C. The volume of solution (V) used was the recommended 10 mL per gram of glass. Leachates were filtered to remove colloids and/or particulates. At the end of triplicate analyses, both the pH and the leachate concentrations were measured for the glass species of interest. A standard glass was used as a control to eliminate bias in the experimental analysis and in the analytical data. Triplicate analyses and triplicate glass standards were run simultaneously. A multielement solution standard was used to detect any significant biases in the analysis of the leachates.

The leachate concentrations are reported as normalized elemental losses,  $\text{NC}_i$ , released from the glass in grams of glass per L of leachant. This has the advantage that the release concentrations in parts per million are normalized by the weight fraction of that element present in the glass. The

normalized release,  $NC_i$ , is calculated as:

$$NC_i = \frac{C_i}{F_i} \quad (1)$$

where  $NC_i$  = normalized release ( $g_{\text{glass}}/L_{\text{leachant}}$ )  
 $C_i$  = mass of element "i" in the solution ( $g_i/m^3$ )  
 $F_i$  = fraction of element "i" in the glass ( $g_i/g_{\text{glass}}$ )

## **QUALITY ASSURANCE**

All the vitrification activities and glass analyses were performed in accordance with DWPT Task Plan-IDMS Mercury Studies, DWPTQA-89-0041. All tasks were controlled in accordance with the task QA Plan.

The PCT Version 3.0 is a Glass Technology Category 1 Procedure requiring experimenter data input and signoff at every step (GTOP-3-025 in The Glass Technology Procedures Manual, DPSTM-88-700-5, L 13-1). All the ovens, balances, and water purification systems used for the PCT are M&TE Category 1.

Analytical Development Section (ADS) procedures were followed for all chemical and x-ray diffraction data so that the data is readily retrievable.

All the PCT data for this study are recorded in DPSTN-4789 (E-56079) and WSRC-NB-90-271. The glass composition data is recorded in DPSTN-4771 (E-56053).

## **GLASS SAMPLE IDENTIFICATION**

The glass nomenclature Hg 1-1, 1-2, 1-3 was used to facilitate the understanding of the chemical variation occurring in the first (Hg-1) run of the IDMS Hg campaign as a function of the sequence in which the glasses were sampled.<sup>19</sup> Glasses Hg 1-1, 1-2, and 1-3 appear in the IDMS production records as GLAS 2340, 2427, and 2518, respectively. Glasses Hg 2-1, 2-2, and 2-3 are GLAS 2630, 2838, and 2877, respectively. Glass Hg 3-1 is a combined sample of production run GLAS 2124 and 2125. Glass Hg 3-2 is a combined sample of production run GLAS 3236 and 3237. Glass Hg 3-3 is production run GLAS 3357.

## RESULTS AND DISCUSSION

### Glass Homogeneity

The x-ray diffraction analysis of the crushed and sieved glass indicated that there was no crystallization present in any of the IDMS-Hg glasses. During SEM/EDAX analysis of the crushed glass, small amounts of metallic Cu<sup>0</sup> were observed. The Cu<sup>0</sup> appeared to be on the glass surface, indicating that it may have been contamination from the brass sieves. No evidence of glass crystallization was observed during SEM/EDAX analysis.

### Glass Composition

The chemical analyses indicated that there had been an excursion in Na<sub>2</sub>O content in the glass vitrified in IDMS between runs Hg-1 and Hg-2. This excursion occurred when the vitrified slurry analyses indicated that the resulting melt would be too viscous to pour, e.g., over the viscosity process control limit of 100 poise.<sup>19</sup> NaOH was added to run Hg-2 to adjust the feed so that it would form a glass of < 100 poise viscosity. Additional PHA, which is Na<sub>2</sub>O rich, was added to the run Hg-3 feed instead of NaOH, so that the Hg-3 run glass would conform to the viscosity process limit of <100 poise.

Analysis of the glass used in the durability studies indicated that the glasses run during run Hg-1 (Hg 1-1, Hg 1-2, and Hg 1-3) and the very first sample taken at the beginning of run Hg-2 (Hg 2-1) were consistent in composition while the sequence of glasses after the NaOH (Hg 2-2, Hg 2-3) and the PHA (Hg 3-1, Hg 3-2, Hg 3-3) additions were consistent. The glass compositions were measured in duplicate and were determined to be biased low compared to the original analyses of these glasses at the TNX laboratory. Therefore, the reanalyzed glass compositions were bias-corrected to the original IDMS Hg-1 glass data (Appendix I). The measured redox ratios (Fe<sup>2+</sup>/ΣFe) of the IDMS glasses were <0.05 during runs Hg-1 and Hg-2 and between 0.07 and 0.1 during run Hg-3 (Appendix I). The measured redox for Hg-1 and Hg-2 glasses were more oxidizing than those measured from the vitrified feed for Hg-1 and Hg-2. The measured redox for Hg-3 glasses was slightly more reduced than the redox measured from the vitrified feed for Hg-3.<sup>19</sup>

The effect of the glass composition variability calculated for each of the compositionally dependent process (viscosity and liquidus) and product (durability) constraints is shown in Figure 1a-c as a function of production run sequence. The process and product parameters calculated from the average glass compositions given in Appendix I followed the same trends as the parameters calculated<sup>19</sup> from the vitrified melter feed for all runs. The chemical analysis of vitrified feed from run Hg 2-1 indicated that the melt would have exceeded the upper viscosity limit of 100 poise (Figure 1a) and the upper liquidus temperature of 1050°C (Figure 1b) had the glass composition not been remediated to the value indicated by the Hg 2-2 glass analysis. Glass Hg 2-1 had been sampled early in run 2 before the melt volume had completely turned over after the addition of the NaOH. Moreover, Figure 1c indicates that the remediation of the glass composition within the process constraints did not adversely affect the glass durability, e.g. the calculated glass durability, expressed as the glass free energy of hydration, remained a more positive value than -7 kcal/mole.

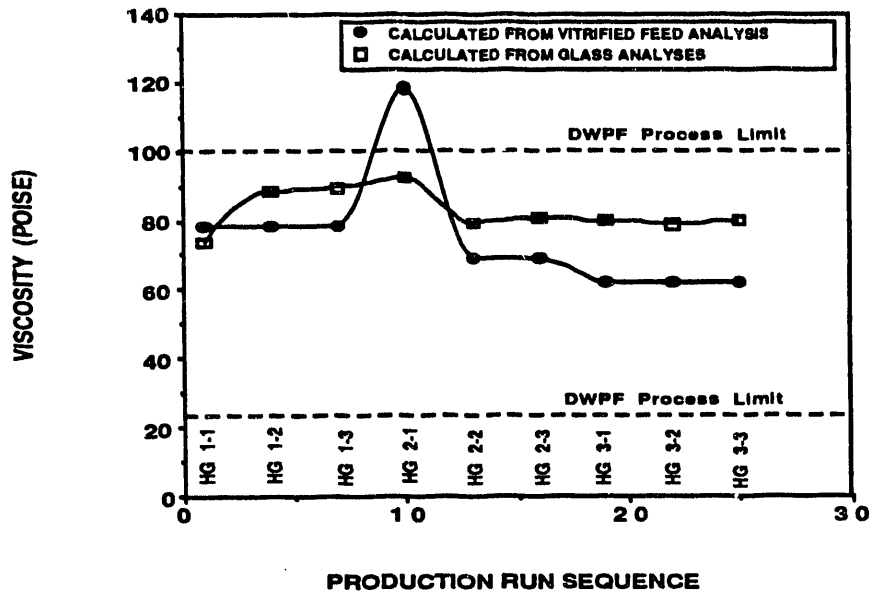


Figure 1a. Comparison of predicted glass viscosity based on vitrified feed analyses and final glass composition analysis.

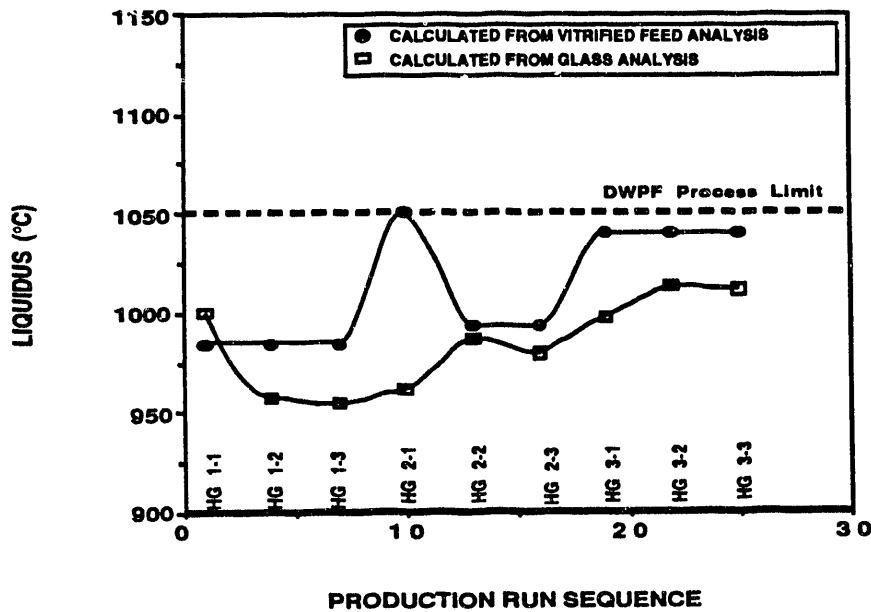


Figure 1b. Comparison of predicted glass liquidus based on vitrified feed analyses and final glass product analysis.

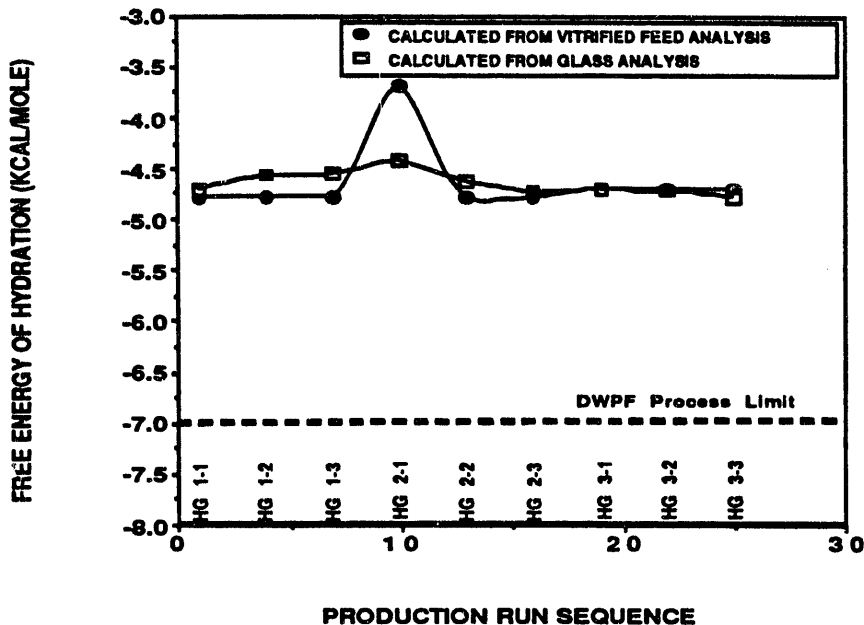
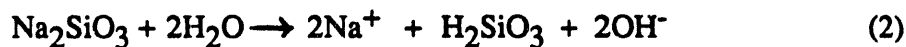


Figure 1c. Comparison of predicted glass liquidus based on vitrified feed analyses and final glass product analysis.

### Glass Durability

The triplicate glass durability analysis pH values shown in Figure 2 also reflect the changing alkali content of the Hg campaign glass as a function of production run sequence. The leachate pH changed significantly between campaign Hg 2-1 and Hg 2-2 when the excess NaOH addition to the sludge became a significant contribution in the resulting glass. This is not surprising since excess Na<sub>2</sub>O in the glass would cause excess OH<sup>-</sup> in the leachate which causes the leachate pH to become more basic via the following reaction:



The pH corrected free energy of hydration was calculated from the glass composition and the leachate pH (Appendix I).<sup>12</sup> The pH corrected free energy of hydration was plotted as a function of glass production run sequence for the IDMS-Hg glasses (Figure 3). A significant change in the free energy of hydration,  $\Delta G_{\text{hyd}}$ , is noted between runs Hg 2-1 where the calculated value is about -7 kcal/mole and Hg 2-2 where the calculated value is about -8 kcal/mole. The change in the pH corrected  $\Delta G_{\text{hyd}}$  coincided with the the point at which the excess Na<sub>2</sub>O became a significant contribution to the glass.

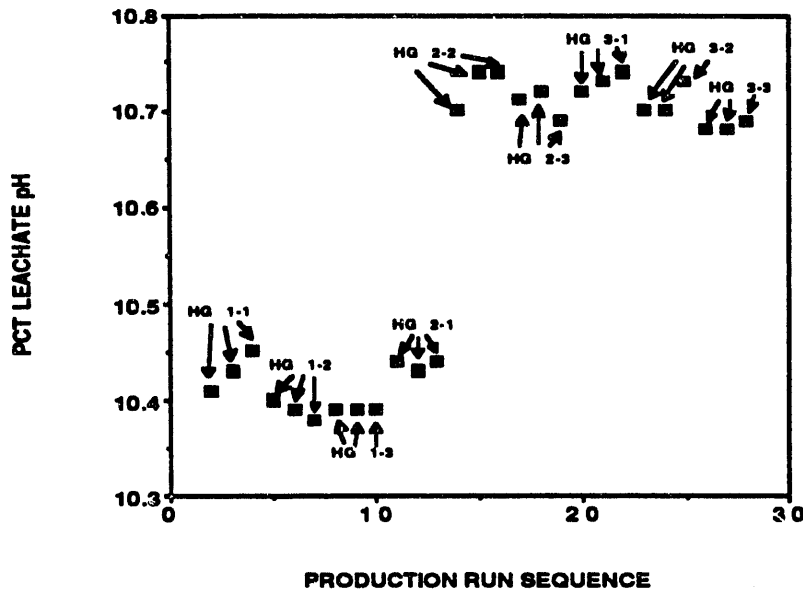


Figure 2. Variation of leachate pH with IDMS Hg glasses in the sequence in which the glasses were produced and poured.

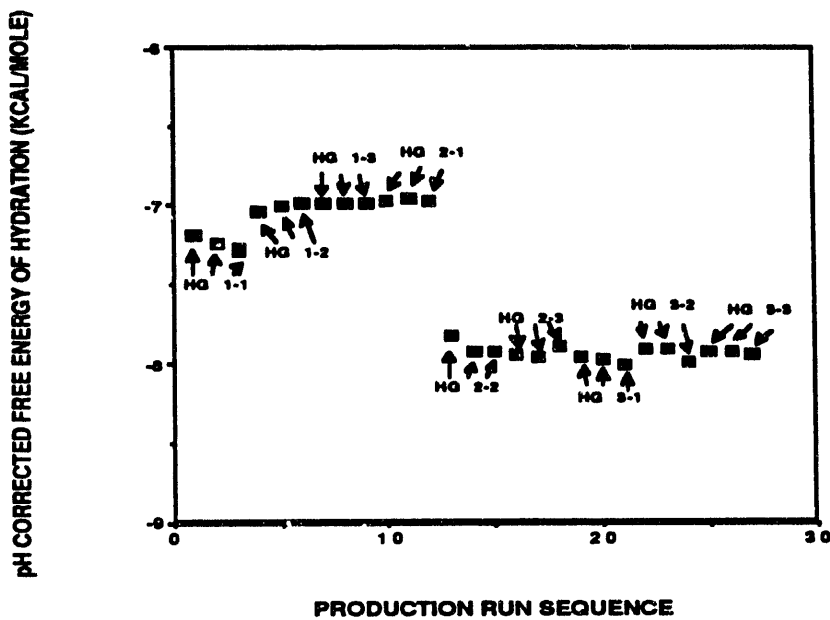


Figure 3. Variation of the calculated free energy of hydration (calculated from the glass composition and the solution pH) for IDMS Hg glasses shown in the sequence in which the glasses were produced and poured.

A multielement solution standard was run after every 5 leachate samples. Analysis of the solution standard data indicated that there was less than 0.5 ppm bias in the ICP analyses for Na, B, Li, Al, and Fe and in the Atomic Absorption (AA) analyses for K. The data also indicated that there was about a 1.4 ppm bias in the Na determinations made by AA and a -1.47 bias in the Si determinations by ICP analysis (Appendix II).

The raw data used to convert the leachate concentrations to  $NC_i$  is given in Appendix III. Boron and lithium releases from glass are considered to be the most accurate indicators of glass durability since these elements are leached from glass faster than any other elements. Boron is probably the best indicator of glass durability because it does not saturate in the leachate and does not participate in precipitation reactions caused by solution supersaturation. The elemental releases measured in the leachates were checked for internal consistency by plotting the normalized releases for Li, Si, Na, and K against the normalized releases of B (Figure 4). Normalized units are plotted since the glasses before run Hg 2-2 are considerably different in composition than those after run Hg 2-2. The plots of Li, Si, Na and K versus B were all consistent, e.g. if a given glass released a great deal of B it, likewise, released a great deal of Li and/or Si, Na, or K. Figure 4 also demonstrates a consistently high bias in the Na analyses by AA over the Na analyses by ICP.

The normalized elemental release of B was plotted against the pH corrected  $\Delta G_{hyd}$ . The elemental release of B from the IDMS Hg glasses were compared with PCT results from previous experimentation (Table 1 and Figure 5). The previous experimental results had been collected prior to May 1989. Historical control charting of the ARM-1 standard glass indicated a shift in the elemental releases for all elements in May 1989. This shift correlated with a change from PCT 2.0 to PCT 3.0 when a glass powder washing procedure was introduced to remove adhering fine particles from the sieved glass. Since removal of the glass fines prior to conducting the durability test biases the PCT 3.0 leachate data lower than the historic data, the historic data was bias corrected to be consistent with the data derived from PCT 3.0.

The bias corrected normalized B release for the historic data and the PCT 3.0 leachate data for the Hg glasses was plotted against the  $\Delta G_{hyd}$  calculated from the glass composition and the solution pH. This comparison indicated that the glasses from the IDMS Hg-1 campaign (Hg 1-1, Hg 1-2 and Hg 1-3) and the glass from the first sample of the Hg-2 campaign (Hg 2-1) were very similar in durability to DWPF startup glass, 165 sludge-only glasses, and Frit 202 glass with excess alkali (Figure 5). The remaining glasses in the Hg-2 campaign (Hg 2-2 and Hg 2-3) and in the Hg-3 campaign (Hg 3-1, Hg 3-2, and Hg 3-3) were higher in  $Na_2O$  content due to NaOH addition during run Hg-2 and the excess PHA addition during run Hg-3 in order to adjust the glass viscosity. These glasses were less durable than the Frit 202 glass with excess PHA but more durable than 131 sludge-only glasses. All the IDMS Hg glasses are more durable than the glass that was qualified in the DWPF Environmental Assessment (EA).<sup>10</sup> Comparisons of the data in Table 1 for the average EA glass leachate concentrations and the data in Appendix I for the Hg Campaign glasses indicates that the Hg Campaign glasses are over 20 times more durable than the EA glass. The IDMS Hg campaign glasses would, therefore, meet current Waste Acceptance Preliminary Specification (WAPS) criteria for product consistency.

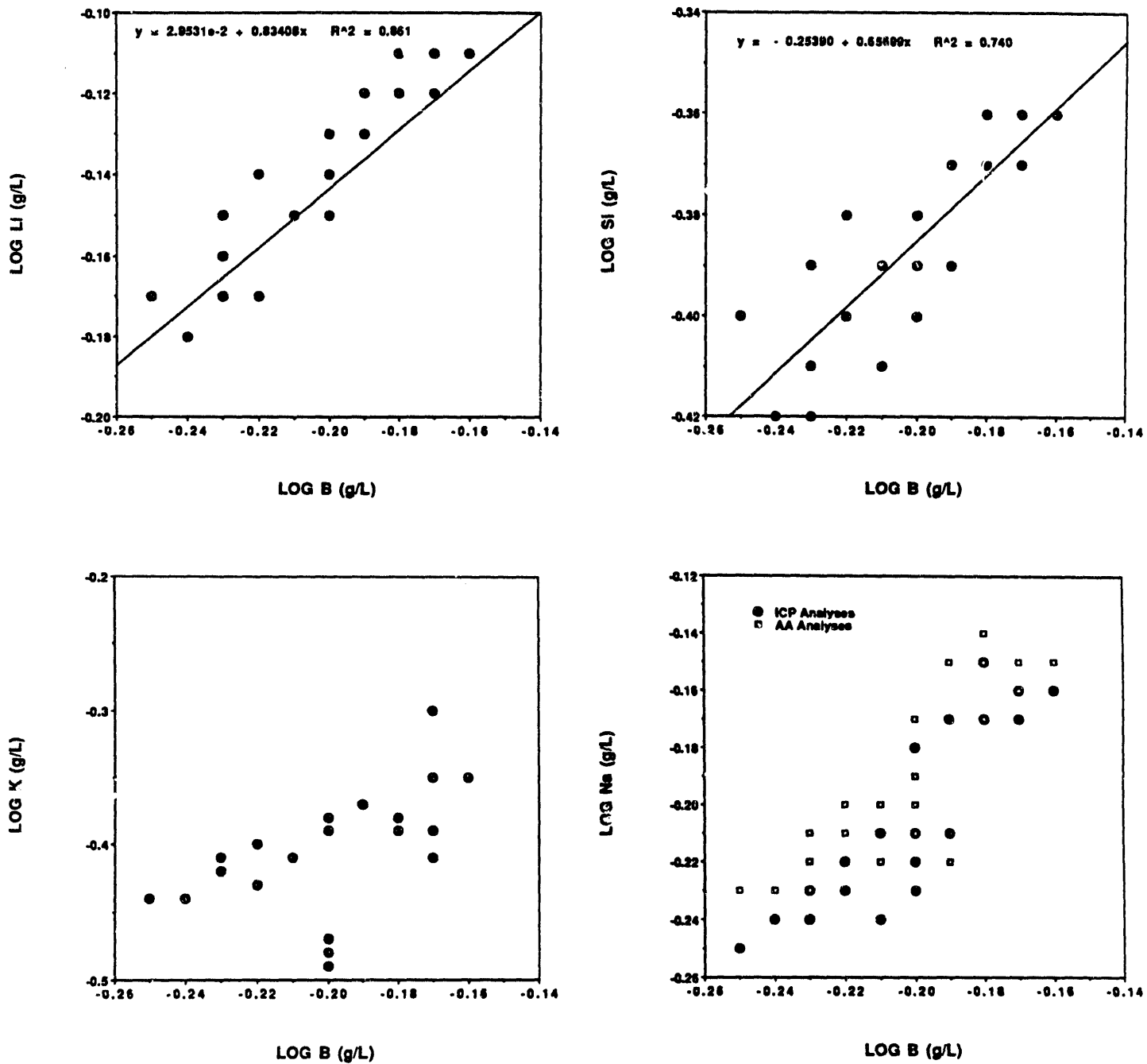
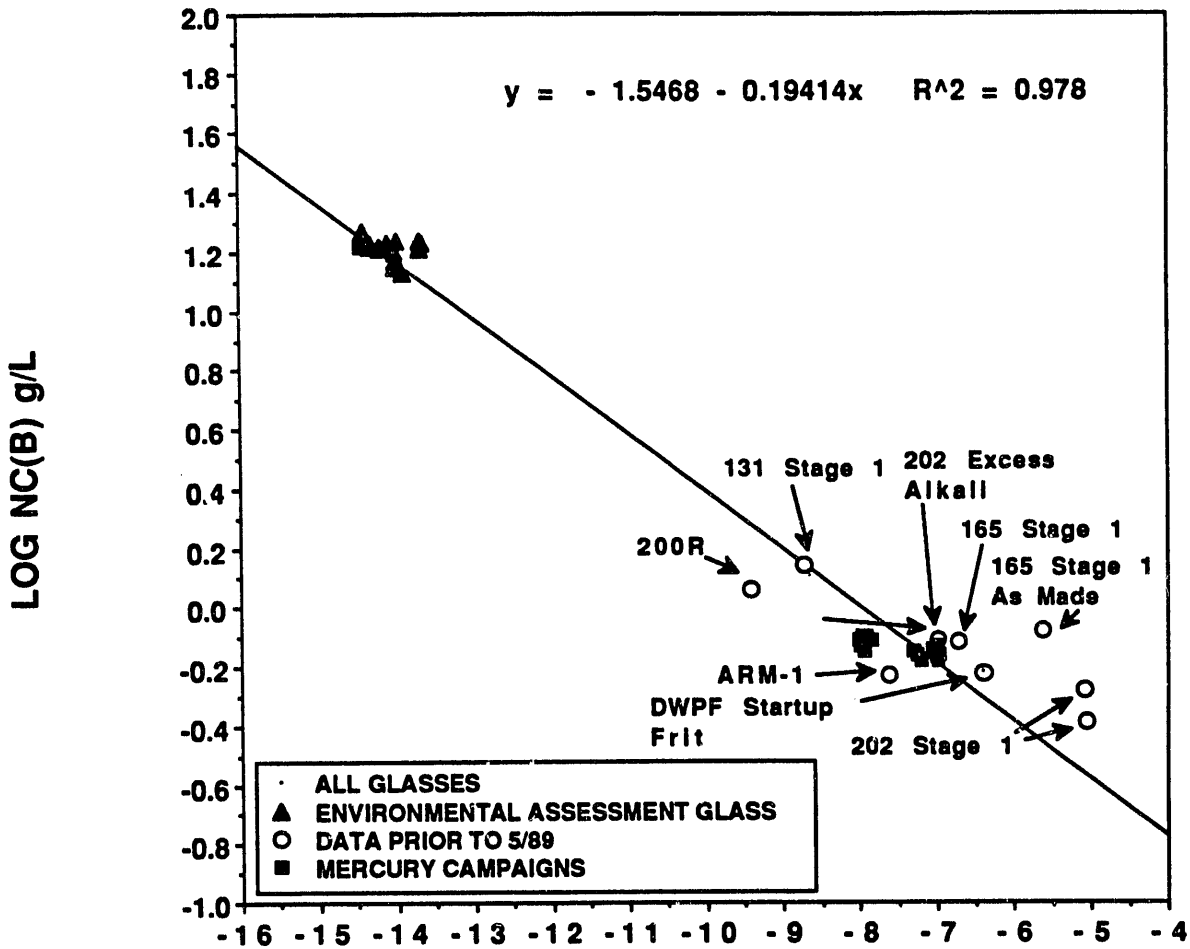


Figure 4. Internal consistency of IDMS leachate analyses. Note that the IDMS leachate analyses are significantly lower than the average values of Li, Na, B and Si for the EA glass which are  $189 \pm 17$ ,  $1650 \pm 131$ , and  $576 \pm 46$ ,  $882 \pm 105$  ppm, respectively. See Table 1.





### pH CORRECTED FREE ENERGY OF HYDRATION (Kcal/mole)

Figure 5. Correlation of the free energy of hydration,  $\Delta G_{hyd}$ , calculated from the glass composition and the leachate pH, versus the logarithm of the normalized boron release from glasses investigated for the Defense Waste Processing Facility (DWPF). The measured durability values for the IDMS Hg-1, Hg-2, and Hg-3 campaigns with Batch 1 sludge, PHA, and 202 frit are overlain for reference.

### CONCLUSIONS

Glasses produced during the three IDMS Hg run campaigns were produced using the DWPF process/product models. The durability model indicated that the glasses would all be more durable than the glass qualified in the DWPF Environmental Assessment (EA). The measured glass durability demonstrated that the Hg campaign glasses were as durable as previously examined Frit 202 PHA glasses and sludge-only 165 glasses and more durable than the EA glass. The IDMS Hg glasses, therefore, meet the current Waste Acceptance Preliminary Specification (WAPS) 1.3 criteria for control of product consistency.

**Table 1. Calculated Free Energies of Hydration and PCT Durability of DWPF Borosilicate Waste Glasses**

<b>Glass/Reference</b>	<b>pH Corrected Free Energy of Hydration (Kcal/mole)</b>	<b>Log NC(Si)</b>	<b>Log NC(B)</b>	<b>Log NC(Na)</b>	<b>Log NC(Li)</b>	<b>pH</b>	<b>PCT</b>
		<b>(g<sub>glass</sub>/L<sub>leachate</sub>)</b>					
EA Glass-Average of 24 PCT Tests <sup>20</sup>	-14.05	0.58	1.21	1.12	0.98	11.92	3.0
202 Stage 1 (AH-10) DPSTN-4724	-5.050	-0.53*	-0.39*	-0.51*	-0.35	9.59	2.1
165 Stage 1 (AH-165) DPSTN-4630	-5.62	-0.32	-0.08	-0.14	-0.061	9.73	2.1W
165 Standard (Corning) DPSTN-4575	-6.72	-0.29*	-0.12*	-0.07*	-0.14*	10.31	2.0
DWPF Startup Frit-Remelted DPSTN-4630	-6.39	-0.90*	-0.22*	-0.23*	-0.33*	10.24	2.1
202-Stage 1 Excess Alkali (Corning 202P) DPSTN-4630	-6.99	-0.41*	-0.11*	-0.08*	-0.16*	10.18	2.1
200 Average Radioactive DPSTN-4570	-9.41	-0.33	0.06	0.00	-0.01	10.55	2.1W
131 Stage 1 (AH 131) DPSTN-4630	-8.71	-0.51	0.14	0.10	0.02	10.43	2.1W
202-Stage 1 (Corning 202G) DPSTN-4631	-5.08	-0.46*	-0.28*	-0.20	-0.24	9.79	2.1

\* Solution data corrected for Si bias of -13.31 ppm, B bias of -4.10 ppm, Na bias of -8.4 ppm, and an Li bias of -3.34 ppm; higher solution values were noted when glass fines were not removed from the meshed glass prior to durability testing (PCT 2.0 and 2.1). PCT 2.1W indicates PCT procedure 2.1 with fines washing and these data are directly comparable to data from PCT 3.0 without bias correction.

## **REFERENCES**

1. R. E. Eibling and J.R. Fowler, "Mercury Reduction and Removal During High-Level Radioactive Waste Processing and Vitrification," Scientific Basis for Nuclear Waste Management, IV, S.V. Topp (Ed.), Elsevier Publishers, New York, 617-622 (1982).
2. C.M. Jantzen, "Relationship of Glass Composition to Glass Viscosity, Resistivity, Liquidus Temperature, and Durability: First Principles Process-Product Models for Vitrification of Nuclear Waste," Proceedings of the 5th International Symposium on Ceramics in Nuclear Waste Management, G.G. Wicks, D.F. Bickford, and R. Bunnell (Eds.), American Ceramic Society, Westerville, OH, 37-51 (1991).
3. K.G. Brown, R.L. Postles, and R.E. Edwards, "The Product Composition Control System (PCCS)," Proceedings of the 5th International Symposium on Ceramics in Nuclear Waste Management, G.G. Wicks, D.F. Bickford, and R. Bunnell (Eds.), American Ceramic Society, Westerville, OH, 37-51 559-568 (1991).
4. Waste Form Compliance Plan for the Defense Waste Processing Facility," USDOE Report WSRC-SW4-6, Rev. 0, Westinghouse Savannah River Co., Savannah River Laboratory, Aiken, SC (March, 1990).
5. C. M. Jantzen, "Glass Compositions and Frit Formulations Developed for DPWF," USDOE Report DPST-88-952, E.I. duPont deNemours & Co., Savannah River Laboratory, Aiken, SC (November, 1988).
6. C. M. Jantzen and N. E. Bibler, "Product Consistency Test (PCT) and Test Protocol," USDOE Report DPST-87-575, E.I. duPont deNemours & Co., Savannah River Laboratory, Aiken, SC (1987).
7. C.M. Jantzen and N.E. Bibler, "The Product Consistency Test for the DWPF Wasteform," Requirements for Waste Acceptance and Quality Control, Proceedings of the 2nd International Seminar on Radioactive Waste Products, E. Warnecke, R. Odoj, and R. Simon (Eds.), Kernforschunganlanger (KFA) Research Centre Julich, Federal Republic of Germany, 609-622 (1990).
8. C.M. Jantzen and N.E. Bibler, "Nuclear Waste Glass Product Consistency Test (PCT)-Version 3.0," USDOE Report WSRC-TR-90-539, Rev. 1, Westinghouse Savannah River Co., Aiken, SC (November, 1990).
9. Waste Acceptance Preliminary Specifications for the Defense Waste Processing Facility High-Level Waste Form, U. S. Department of Energy, Office of Civilian Radioactive Waste Management, USDOE Document OGR/B-8, Washington, DC (1988).
10. U.S. Department of Energy, "Environmental Assessment Waste Form Selection for SRP High-Level Waste," U.S. DOE Report DOE-EA-0179 (July, 1982).

11. M.J. Plodinec, C.M. Jantzen, and G.G. Wicks, "Thermodynamic Approach to Prediction of the Stability of Proposed Radwaste Glasses," Advances in Ceramics, 8, 491-495 (1984).
12. C.M. Jantzen and M.J. Plodinec, "Thermodynamic Model of Natural, Medieval, and Nuclear Waste Glass Durability," J. Non-Cryst. Solids, 67, 207-233 (1984).
13. M.J. Plodinec, C.M. Jantzen, and G.G. Wicks, "Stability of Radioactive Waste Glasses Assessed from Hydration Thermodynamics," Sci. Basis for Nuclear Waste Management, VII, G.L. McVay (ed). Elsevier Publ., New York 755-762 (1984).
14. C.M. Jantzen, "Prediction of Nuclear Waste Glass Durability from Natural Analogs" Advances in Ceramics, 20, D.E. Clark, W.B. White and A.J. Machiels (Eds.), American Ceramic Society, Westerville, OH, 703-712 (1986).
15. C.M. Jantzen, "Prediction of Glass Durability as a Function of Environmental Conditions", Proceedings of the Symposium on Materials Stability and Environmental Degredation, A. Barkatt et. al. (Eds.), Materials Research Society, Pittsburgh, PA, 143-159 (1988).
16. C.M. Jantzen, "Prediction of Glass Durability as a Function of Glass Composition and Test Conditions: Thermodynamics and Kinetics", Proceedings of the First Intl. Conference on Advances in the Fusion of Glass, American Ceramic Society, Westerville, OH , p. 24.1-24.17 (1988).
17. C.M. Jantzen, "Thermodynamic Approach to Glass Corrosion," Corrosion of Glass, Ceramics, and Ceramic Superconductors, Noyes Publications, Park Ridge, NJ, 153-215 (1992).
18. E.W. Baumann, "Colorimetric Determination of Iron (II) and Iron (III) in Glass," Analyst, v. 117, 913-916 (1992).
19. N.D. Hutson, J.R. Zamecnik, M.E. Smith, D.H. Miller and J.A. Ritter, "Integrated DWPF Melter System (IDMS) Campaign Report: Mercury Operation," USDOE Report WSRC-TR-91-0363, Westinghouse Savannah River Co., Savannah River Laboratory, Aiken, SC (May, 1991).
20. C.M. Jantzen, "Characterization of the Defense Waste Processing Facility (DWPF) Environmental Assessment (EA) Glass Standard Reference Material," USDOE Report WSRC-TR-92-346, Rev. 0, Westinghouse Savannah River Co., Savannah River Laboratory, Aiken, SC (September, 1992).

# APPENDIX I Measured Glass Compositions

GLASS ID= HG1-1A (SPREADSHEET REVISION 5.0, OCTOBER 23, 1990)

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	MOLE FRACTION	MOLE %	NORMOXIDE MOLE %	COMPONENT FREEENERGY	FREEENERGY HYDRATION	NBD	KCAL/KG
Al2O3	2.3410	1.8895	4.4233	1.21	5.35	101.9600	0.0524	3.3227	3.0400	3.0400	0.1010	-0.0665	1.5943
CaO	0.7770	1.3992	1.0872	1.11	1.21	56.0794	0.0215	1.3646	-16.1160	-16.1160	-0.2199	0.0273	-3.4711
Fe2O3	6.1820	1.4297	11.6512	1.04	12.09	159.6922	0.0757	4.7959	15.5000	0.7434	0.7434	-0.0959	11.7326
Fo	N/A	1.2865	0.0419	1.00	0.04	71.8464	0.0006	0.0370	-14.6090	-0.0054	-0.0054	0.0007	-0.0853
MgO	0.7980	1.6583	1.3233	1.12	1.49	40.3114	0.0369	2.3388	-13.8680	-0.3248	-0.3248	0.0466	-5.1266
MnO	1.4320	1.2912	1.8490	1.31	2.42	70.9374	0.0341	2.1617	-14.8710	-0.3215	-0.3215	0.0432	-5.0737
Na2O	6.5850	1.3480	8.8766	1.03	9.14	61.9790	0.1474	9.5034	-22.8150	-2.8910	-2.8910	0.1868	-42.4721
Li2O	1.9680	2.1525	4.2318	1.06	4.48	29.8774	0.1500	9.5034	-22.7400	-2.1611	-2.1611	0.1901	-34.1084
NO	0.5720	1.2726	0.7279	1.18	0.86	74.7094	0.0115	0.7302	-14.3470	-0.1048	-0.1048	0.0146	-1.6534
SiO2	23.0310	2.1393	49.2702	1.10	54.17	60.0848	0.9016	57.1238	5.5900	1.8398	0.0587	-0.0001	25.8763
Cr2O3	0.1620	1.4816	0.2368	1.59	0.38	151.9974	0.0025	0.1571	37.3600	0.0587	0.0587	0.0000	0.9264
B2O3	2.1240	3.2201	6.8395	1.02	6.86	69.6204	0.0999	6.3317	-9.9300	-0.6287	-0.6287	0.0000	-9.9235
Li2O	1.1344	0.0000	0.0000	1.00	0.00	270.0398	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000	0.0000
TiO2	1.1379	0.0000	0.0000	1.00	0.00	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000	0.0000
SO	0.0220	1.1826	0.0260	1.00	0.03	103.6194	0.0003	0.0159	-24.4000	-0.0039	-0.0039	0.0003	-0.0613
ZrO2	0.1180	1.3508	0.1594	1.00	0.16	123.2188	0.0013	0.0818	45.1000	0.0369	0.0369	0.0033	0.5620
TiO2	0.2170	1.6680	0.3620	0.98	0.35	79.8988	0.0044	0.2801	15.9900	0.0448	0.0448	0.0112	0.7069
K2O	2.4100	1.2046	2.9031	0.92	2.67	94.2034	0.0283	1.7952	-41.7350	-0.7492	-0.7492	0.0359	-11.8250
CaO	0.0850	1.0602	0.0688	1.00	0.07	281.8094	0.0002	0.0155	-46.8200	-0.0073	-0.0073	0.0003	-0.1145
Si2O3	1.1970	0.0000	0.0000	1.00	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0275	0.0275	1.00	0.03	141.9370	0.0002	0.0123	-53.8200	-0.0066	-0.0066	0.0000	-0.1039
Na2O3	1.1660	0.0000	0.0000	1.00	0.00	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	0.0000	1.00	0.00	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	0.0000	1.00	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000	0.0000
B2O	1.1165	0.0692	0.0692	1.14	0.08	153.3394	0.0005	0.0326	-30.5700	-0.0100	-0.0100	0.0007	-0.1571
RO	0.0940	1.0772	0.1013	2.11	0.21	223.1900	0.0010	0.0608	-8.8100	-0.0053	-0.0053	0.0012	-0.0843
CaO2	1.2284	0.0000	0.0000	1.00	0.00	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000	0.0000
Mn2O3	1.5003	0.0000	0.0000	1.00	0.00	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000	0.0000
Zn	1.2447	0.2029	0.2029	2.00	0.41	81.3900	0.0050	0.3168	-2.4000	-0.0076	-0.0076	0.0063	-0.1199
ClO	1.2518	0.1953	0.1953	1.18	0.23	79.5454	0.0029	0.1838	1.7900	0.0033	0.0033	0.0037	0.0519
Fe2+/Fe3+													
SUMS	51.2890	94.6743	102.8107	100.0000	100.0000	1.5763	100.0000	SUM + pHSI	1.7900	-4.8193	-4.8193	0.4069	-72.9076
								SUM + pHB		-5.4068	-5.4068		
								pH=		-7.1494	-7.1494		
										10.43	10.43		

ALK/SIO2 = 0.362  
 ALK+ B2O3/SIO2 = 0.472  
 B2O3/SIO2 = 0.111  
 VISCOSITY @TEMP(C) = 85.319  
 LIQUIDUS RATIO = 0.084  
 LIQUIDUS (C) = 967

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

(SPREADSHEET REVISION 5.8, OCTOBER 23, 1990)

GLASS ID= HG1-1B

ELEMENT	WT%	GRAV FACTOR	OXIDE WT%	BIAS CORRECT	BIAS CORRECT	OXIDE WT%	NORMOXIDE	M.W.	MOLE FRACTION	NORMOXIDE MOLE %	COMPONENT FREEENERGY	FREEENERGY HYDRATION	NBO	KCAL/KG
Al2O3	2.4220	1.8895	4.5764	1.21	5.5321	5.4931	101.9600	0.0543	3.5547	3.0400	0.1081	0.1081	-0.0711	1.8494
CaO	0.9160	1.3992	1.2817	1.11	1.4239	1.4139	56.0794	0.0254	1.8935	-16.1160	-0.2661	-0.2661	0.0333	-4.0920
Fe2O3	9.3400	1.4297	13.3002	1.04	13.7986	13.7013	151.6922	0.0864	5.6611	15.5000	0.0864	0.0864	-0.1132	13.3931
FeO	N/A	1.2665	0.0479	1.00	0.0479	0.0475	71.8464	0.0007	0.0075	-14.6090	-0.0064	-0.0064	0.0009	-0.0973
MgO	0.8500	1.6583	1.4096	1.12	1.5950	1.5738	40.3114	0.0393	2.5760	-13.8880	-0.3578	-0.3578	0.0515	-5.4607
MnO	1.1310	1.2912	1.4603	1.31	1.9115	1.8960	70.9374	0.0269	1.7654	-14.8710	-0.2625	-0.2625	0.0353	-4.0072
Na2O	6.6510	1.3480	8.9655	1.03	9.2270	9.1618	61.9790	0.1489	9.7536	-28.8150	-2.8105	-2.8105	0.1951	-42.8978
Li2O	1.9120	2.1525	4.1156	1.06	4.3593	4.3278	29.8774	0.1459	9.5570	-22.7400	-2.1793	-2.1793	0.1911	-33.1716
NO	0.5720	1.2726	0.7279	1.18	0.8610	0.8549	74.7094	0.0115	0.7550	-14.3470	-0.1083	-0.1083	0.0151	-1.8534
SiO2	21.3870	2.1393	45.7532	1.10	50.3051	49.9503	60.0648	0.8372	54.8522	5.9900	1.4837	1.4837	0.0000	22.6470
Cr2O3	0.1360	1.4616	0.1988	1.59	0.3164	0.3142	151.9974	0.0021	0.1364	37.3600	0.0510	0.0510	-0.0027	0.7777
B2O3	2.1650	3.2201	6.9715	1.02	7.0918	7.0418	69.6204	0.1019	6.6737	-9.9300	-0.6627	-0.6627	0.0000	-10.1151
Li2O	1.1344	1.1344	0.0000	1.00	0.0000	0.0000	270.0398	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000	0.0000
ThO2	1.1379	1.1379	0.0000	1.00	0.0000	0.0000	284.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000	0.0000
SiO	0.0030	1.1826	0.0035	1.00	0.0035	0.0035	103.6194	0.0000	0.0000	-24.4000	-0.0005	-0.0005	0.0000	-0.0084
ZrO2	0.1180	1.3508	0.1594	1.00	0.1594	0.1579	123.2188	0.0013	0.0845	45.1000	0.0381	0.0381	0.0034	0.5820
TiO2	0.3540	1.6880	0.5905	0.98	0.5763	0.5722	79.8988	0.0072	0.4725	15.9900	0.0758	0.0758	0.0189	1.1533
K2O	2.4080	1.2046	2.9007	0.92	2.6689	2.6481	94.2034	0.0283	1.8548	-41.7350	-0.7741	-0.7741	0.0371	-11.8152
Ca2O	1.0802	0.6899	0.0899	1.00	0.0899	0.0884	291.4992	0.0002	0.0160	-46.8200	-0.0075	-0.0075	0.0003	-0.1145
Si2O3	1.1970	1.1970	0.0000	1.00	0.0000	0.0000	141.9370	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	2.2910	0.0000	1.00	0.0000	0.0000	336.4782	0.0000	0.0000	-53.6200	0.0000	0.0000	0.0000	0.0000
Na2O3	1.1660	1.1660	0.0000	1.00	0.0000	0.0000	325.8100	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000	0.0000
La2O3	1.1728	1.1728	0.0000	1.00	0.0000	0.0000	225.8082	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000	0.0000
Y2O3	1.2699	1.2699	0.0000	1.00	0.0000	0.0000	223.1900	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000	0.0000
B2O	0.1160	1.1165	0.1295	1.14	0.1474	0.1464	153.3394	0.0010	0.0630	-30.5700	-0.0183	-0.0183	0.0013	-0.2939
RbO	1.0772	1.0772	0.0000	2.11	0.0000	0.0000	172.1200	0.0000	0.0000	-8.8100	0.0000	0.0000	0.0000	0.0000
CaO2	1.2284	1.2284	0.0000	1.00	0.0000	0.0000	143.9382	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000	0.0000
Mn2O3	1.5003	1.5003	0.0000	1.00	0.0000	0.0000	81.3800	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000	0.0000
ZnO	0.0870	1.2447	0.1083	2.00	0.2170	0.2155	79.5454	0.0027	0.1747	-2.4000	-0.0042	-0.0042	0.0035	-0.0640
CaO	0.2790	1.2518	0.3493	1.18	0.4127	0.4098	79.5454	0.0052	0.3399	1.7900	0.0081	0.0081	0.0068	0.0929
Fe2+/Fe3+	0.0040													
SUMS	50.9120		93.1186	100.7103	100.0000	100.0000		1.5263	100.0000				0.4065	-73.4957

TEST= TIME= SA/IV=

SUM + pH SI SUM + pH B pH=

-4.8151 -5.6024 -7.3452 10.43

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

ELEMENT	GRAV	OXIDE	FACTORS FOR	NORMOXIDE	M.W.	MOLE	NORMOXIDE	COMPONENT	FREE ENERGY	NEO	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
Al2O3	2.2690	4.2873	1.21	5.18	101.9600	0.0508	3.1800	3.0400	0.0967	-0.0636	1.5452
CaO	0.6940	0.9710	1.11	1.08	56.0794	0.0192	1.2035	-16.1180	-0.1940	0.0241	-3.1003
Fe2O3	6.1340	11.5713	1.04	12.00	159.8922	0.0752	4.7032	15.5000	0.7290	-0.0941	11.6522
FeO	1.2665	0.0521	1.00	0.05	71.8464	0.0007	0.0453	0.0000	-0.0086	0.0009	-0.1059
MgO	0.8120	1.3465	1.12	1.51	40.3114	0.0378	2.3499	-13.8880	-0.3264	0.0470	-5.2165
MnO	1.3560	1.7509	1.31	2.29	70.9374	0.0323	2.0212	-14.8710	-0.3008	0.0404	-4.8044
Na2O	6.4160	8.6498	1.03	8.90	61.9790	0.1436	8.9848	-28.8150	-2.5890	0.1797	-41.3821
Ni2O	2.0290	4.3674	1.06	4.63	29.8774	0.1548	9.8848	-22.7400	-2.2023	0.1937	-35.2014
Li2O	2.1525	4.3674	1.06	4.63	29.8774	0.1548	9.8848	-22.7400	-2.2023	0.1937	-35.2014
NiO	0.5310	0.6758	1.18	0.80	74.7094	0.0107	0.6693	-14.3470	-0.0960	0.0134	-1.5349
SiO2	23.5070	50.2885	1.10	55.29	60.0848	0.9202	57.5715	5.5900	1.8677	0.0000	28.8569
Cr2O3	0.1550	0.2265	1.59	0.36	151.9874	0.0024	1.4844	37.3600	0.0555	-0.0030	0.8863
B2O3	3.2201	6.8170	1.02	6.93	69.8204	0.0996	6.2315	-9.9300	-0.8188	0.0000	-9.8908
UO2	1.1344	0.0000	1.00	0.00	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1378	0.0000	1.00	0.00	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SO	0.0220	0.0260	1.00	0.03	0.0250	0.0003	0.0157	-24.4000	-0.0038	0.0003	-0.0613
ZrO2	0.0640	0.0865	1.00	0.09	103.6194	0.0007	0.0438	45.1000	0.0197	0.0018	0.3158
TiO2	1.6680	0.2886	0.98	0.28	79.8988	0.0035	0.2205	15.9900	0.0353	0.0088	0.5638
K2O	3.1660	3.8138	0.92	3.51	94.2034	0.0372	2.3287	-41.7350	-0.9719	0.0468	-15.5344
CaO	0.0650	0.0689	1.00	0.07	281.8094	0.0002	0.0153	-46.8200	-0.0072	0.0003	-0.1145
SrO	1.1970	0.0000	1.00	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0684	1.00	0.07	141.9370	0.0005	0.0293	-53.8200	-0.0157	0.0000	-0.2510
Nb2O5	1.1660	0.0000	1.00	0.00	325.8100	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	336.4782	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
B2O	1.1165	0.0681	1.14	0.08	153.3394	0.0005	0.0316	-30.5700	-0.0097	0.0008	-0.1548
FeO	1.0772	0.0851	2.11	0.18	223.1900	0.0008	0.0503	-8.8100	-0.0044	0.0010	-0.0709
CaO	1.2284	0.0000	1.00	0.00	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MnO	1.5003	0.0000	1.00	0.00	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.1904	2.00	0.38	81.3800	0.0047	0.2934	-2.4000	-0.0070	0.0059	-0.1128
ClO	0.1530	0.1915	1.18	0.23	79.5454	0.0028	0.1780	1.7900	0.0032	0.0036	0.0509
Fe2+/Fe3+	0.0050										
SUMS	51.9850	95.8854	103.8371	103.0000	1.5984	100.0000	SUM + pH SI		-4.7463	0.4074	-75.8646
							SUM + pH B		-5.4929		
							pH =		-7.1839		
									10.39		

TEST=

TIME=

SA/F=

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

(SPREADSHEET REVISION 5.0, OCTOBER 23, 1990)

GLASS ID= HG1-28

ELEMENT	GRAV	OXIDE	FACTORS FOR	NORMOXIDE	M.W.	FRACTION	NORMOXIDE	COMPONENT	FREEENERGY	NED	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	WT%		FRACTION	WT%	FREEENERGY	HYDRATION		
			BIAS CORRECT								
Al2O3	2.2480	4.2478	1.21	4.9933	101.9600	0.0504	3.1749	3.0400	0.0965	-0.0635	1.5309
CaO	0.7010	0.9808	1.11	1.0597	56.0794	0.0194	1.2250	-16.1160	-0.1974	0.0245	-3.1315
Fe2O3	8.1240	11.5571	1.04	11.6602	159.6922	0.0751	4.7335	15.5000	0.7337	-0.0947	11.6379
FeO	N/A	0.0520	1.00	0.0506	71.8464	0.0007	0.3456	-14.6090	-0.0067	0.0009	-0.1057
MgO	0.8160	1.3532	1.12	1.4797	40.3114	0.0377	2.3797	-13.9880	-0.3305	0.0476	-5.2422
MnO	1.3580	1.7534	1.31	2.2320	70.9374	0.0324	2.0398	-14.8710	-0.3033	0.0408	-4.9115
Na2O	6.2400	8.4115	1.03	8.6188	61.9790	0.1397	8.8058	-28.8150	-2.5373	0.1761	-40.2469
Li2O	2.0300	2.3696	1.06	4.4899	29.8774	0.1549	9.7640	-22.7400	-2.2203	0.1953	-35.2188
NO	0.5200	0.6618	1.18	0.7812	74.7094	0.0095	0.6605	-14.3470	-0.0948	0.0132	-1.5031
SiO2	23.5330	50.3441	1.10	53.8294	60.0848	0.9212	58.0789	5.5900	1.7285	0.0000	27.4171
Cr2O3	0.1420	0.2075	1.59	0.3213	151.9974	0.0022	0.1370	37.3600	0.0512	-0.0027	0.8120
B2O3	2.1270	6.8492	1.02	6.7758	69.8204	0.1001	6.3092	-9.9300	-0.6265	0.0000	-9.9375
UO2	1.1344	0.0000	1.00	0.0000	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.0000	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SnO	0.0210	0.0248	1.00	0.0242	103.8194	0.0002	0.0151	-24.4000	-0.0037	0.0003	-0.0585
ZrO2	0.0640	0.0865	1.00	0.0839	123.2188	0.0007	0.0441	45.1000	0.0199	0.0018	0.3158
TiO2	0.1710	0.2852	0.98	0.2707	79.8988	0.0035	0.2196	15.9900	0.0351	0.0088	0.5571
K2O	2.3720	2.8573	0.92	2.5547	94.2034	0.0279	1.7581	-41.7350	-0.7337	0.0352	-11.6385
Ca2O	0.0650	0.0689	1.00	0.0670	281.8094	0.0002	0.0154	-46.8200	-0.0072	0.0003	-0.1145
Sn2O3	1.1970	0.0000	1.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0321	1.00	0.0312	141.9370	0.0002	0.0142	-53.6200	-0.0076	0.0000	-0.1212
Na2O3	1.1860	0.0000	1.00	0.0000	338.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.0000	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.0000	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	0.0580	0.0648	1.14	0.0717	153.3394	0.0005	0.0303	-30.5700	-0.0093	0.0008	-0.1470
PbO	0.0930	1.0772	2.11	0.2055	223.1900	0.0009	0.0597	-8.8100	-0.0053	0.0012	-0.0834
Ca2O2	1.2284	0.0000	1.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MgO3	1.5003	0.0000	1.00	0.0000	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	0.1630	0.2029	2.00	0.3954	81.3900	0.0050	0.3150	-2.4000	-0.0076	0.0063	-0.1199
UO	0.1490	0.1865	1.18	0.2143	79.5454	0.0028	0.1747	1.7900	0.0031	0.0035	0.0498
Fe2+/Fe3+	0.0050										
SUMS	51.0090	94.6970	102.8901	130.0000		1.5662	100.0000		-4.4232	0.3954	-70.1601
ALK/SiO2 =	0.359			TEST=				SUM + pH SI			
ALK+ B2O3/SiO2 =	0.459			TIME=				SUM + pH B			
B2O3/SiO2 =	0.109			SAV=				pH =			
VISCOSITY @TEMP(C)	92.898										
LIQUIDUS RATIO =	0.089										
LIQUIDUS (C) =	937										

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)



ELEMENT	WT%	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	MOLE FRACTION	MOLE %	NORMOXIDE COMPONENT FREE ENERGY	FREE ENERGY FORMATION	NBD	KCAL/KG
Al2O3	2.2180	1.8895	4.1871	1.21	5.06	4.9279	101.9600	0.0498	0.0193	3.1249	3.0400	0.0950	-0.0625	1.5091
CaO	0.6950	1.3992	0.9724	1.11	1.08	1.0518	56.0794	0.0193	0.0193	1.2127	-16.1160	-0.1954	0.0243	-3.1047
Fe2O3	8.0630	1.4297	11.3573	1.04	11.78	11.4717	159.6922	0.0738	0.0021	4.6446	15.5000	0.7199	-0.0929	11.4367
FeO	N/A	1.2865	0.1533	1.00	0.15	0.1492	71.8464	0.0021	0.0021	0.1343	-14.8090	-0.0196	0.0027	-0.3117
MgO	0.8150	1.6583	1.3515	1.12	1.52	1.4796	40.3114	0.0377	0.0377	2.3731	-13.8880	-0.3296	0.0475	-5.2358
MnO	1.3370	1.2912	1.7263	1.31	2.26	2.2000	70.9374	0.0377	0.0377	2.0052	-14.8710	-0.2982	0.0401	-4.7371
Na2O	6.3980	1.3480	8.6245	1.03	8.88	8.6416	61.9790	0.1432	0.1432	9.0147	-28.8150	-2.5976	0.1803	-41.2860
Li2O	2.0430	2.1525	4.3978	1.06	4.66	4.5339	29.8774	0.1559	0.1559	9.8115	-22.7400	-2.2311	0.1962	-53.4443
NiO	0.4940	1.2728	0.6287	1.18	0.74	0.7240	74.7094	0.0100	0.0100	0.6265	-14.3470	-0.0899	0.0125	-1.4280
SiO2	23.5590	2.1393	50.3998	1.10	55.41	53.9506	60.0848	0.9223	0.0019	58.0543	5.9900	1.7183	0.0000	27.2876
Cr2O3	0.1230	1.4616	0.1798	1.59	0.29	0.2786	151.9974	0.0019	0.0019	0.1185	37.3600	0.0443	-0.0024	0.7033
B2O3	2.1390	3.2201	6.8878	1.02	7.01	6.8216	69.6204	0.1006	0.1006	6.3351	-9.8300	-0.6291	0.0000	-9.8936
UO2	1.1344	0.0000	0.0000	1.00	0.00	0.0000	270.0388	0.0000	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	0.0000	1.00	0.00	0.0000	264.0400	0.0000	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SiO	1.1828	0.0260	0.0260	1.00	0.03	0.0253	103.6194	0.0003	0.0003	0.0158	-24.4000	-0.0039	0.0003	-0.0813
ZrO2	1.3508	0.0027	0.0027	1.00	0.00	0.0026	123.2188	0.0004	0.0004	0.0074	45.1000	0.0008	0.0001	0.0099
TiO2	1.6680	0.2702	0.2702	0.98	0.26	0.2587	79.8988	0.0033	0.0033	0.2078	15.9800	0.0332	0.0083	0.5278
K2O	2.3390	1.2048	2.8178	0.92	2.59	2.5221	94.2034	0.0275	0.0275	1.7310	-41.7950	-0.7224	0.0346	-11.4766
CaO	1.0602	0.0889	0.0889	1.00	0.07	0.0671	281.8094	0.0002	0.0002	0.0154	-48.8200	-0.0072	0.0003	-0.1145
Sn2O3	1.1970	0.0000	0.0000	1.00	0.00	0.0000	291.4982	0.0000	0.0000	0.3070	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0573	0.0573	1.00	0.06	0.0558	141.9370	0.0004	0.0004	0.0254	-53.8200	-0.0138	0.0000	-0.2164
N2O3	1.1660	0.0000	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	0.0000	1.00	0.00	0.0000	325.8100	0.0000	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2698	0.0000	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
B2O	1.1185	0.0659	0.0659	1.14	0.07	0.0730	153.3394	0.0005	0.0005	0.0308	-30.5700	-0.0094	0.0008	-0.1495
PbO	1.0772	0.0928	0.0928	2.11	0.20	0.1903	223.1900	0.0009	0.0009	0.0551	-8.8100	-0.0049	0.0011	-0.0771
CaO2	1.2284	0.0000	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MnCO3	1.5003	0.0000	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.1855	0.1855	2.03	0.37	0.3619	81.3800	0.0046	0.0046	0.2875	-2.4000	-0.0069	0.0058	-0.1098
ClO	1.2518	0.1885	0.1885	1.18	0.22	0.2148	79.5454	0.0028	0.0028	0.1744	1.7900	0.0031	0.0035	0.0496
Fe2+/Fe3+	50.9400		94.6393	102.7124	100.0000	100.0000	1.5886	100.0000	1.5886	100.0000	SUM + pH SI	-4.5443	0.4004	-72.1921
SUMS											SUM + pH B	-5.2909		
											pH =	10.39		

TEST= TIME= SA/Y=

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FeO (OCTOBER 23, 1990)

GLASS ID= HG1-3B (SPREADSHEET REVISION 5.0, OCTOBER 23, 1990)

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	MOLE FRACTION	NORMOXIDE MOLE %	COMPONENT FREEENERGY	FREEENERGY HYDRATION	NED	KCAL/KG
Al2O3	2.1670	1.8895	1.21	4.95	4.8536	101.9600	0.0485	3.0766	3.0400	0.0935	-0.0615	1.4758
CaO	0.7380	1.3992	1.11	1.15	1.1249	56.0794	0.0205	1.2965	-16.1160	-0.2069	0.0259	-3.2968
Fe2O3	8.0190	1.4297	1.04	11.72	11.4912	159.6922	0.0734	4.6507	15.5000	0.7209	-0.0930	11.3743
FeO	N/A	1.2865	1.00	0.15	0.1495	71.8464	0.0021	0.1345	-14.6090	-0.0198	0.0027	-0.3100
MgO	0.8050	1.6583	1.12	1.50	1.4720	40.3114	0.0372	2.3600	-13.8680	-0.3278	0.0472	-5.1716
MnO	1.3290	1.2912	1.31	2.25	2.2028	70.9374	0.0128	0.9377	-14.8710	-0.2984	0.0401	-4.7087
Ni2O	6.3710	1.3480	1.03	8.84	8.6671	61.9790	0.1548	9.8105	-22.7400	-2.2309	0.1868	-41.0918
Li2O	2.0290	1.2525	1.06	4.63	4.5353	29.8774	0.0102	0.6487	-14.3470	-0.0931	0.0130	-1.4684
NiO	0.5080	1.2726	1.18	0.76	0.7498	74.7094	0.0018	0.1183	37.3600	0.0442	-0.0024	0.8976
SiO2	23.3800	2.1393	1.10	54.99	53.9256	60.0849	0.0095	6.3036	-9.9300	-0.8259	0.0000	0.0000
Cr2O3	0.1220	1.4616	1.59	0.28	0.2793	151.9974	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
B2O3	2.1140	3.2201	1.02	6.92	6.7904	69.6204	0.0000	0.0000	-2.5330	-0.0042	0.0003	-0.0668
Li2O	1.1344	0.0000	1.00	0.00	0.0000	270.0368	0.0000	0.0000	45.1000	0.0006	0.0001	0.0099
ThO2	0.0240	1.1379	1.00	0.00	0.0000	103.6194	0.0000	0.0000	-24.4000	-0.0042	0.0000	0.0000
ZrO2	0.0020	1.1826	1.00	0.00	0.0000	284.0400	0.0000	0.0000	15.9900	0.0328	0.0082	0.5147
TiO2	0.1580	1.6680	1.00	0.00	0.0000	123.2188	0.0000	0.0000	-41.7350	-0.0073	0.0000	-11.5502
K2O	2.3540	1.2046	0.98	2.61	2.5565	94.2034	0.0000	0.0000	-46.8200	-0.0073	0.0000	-0.1145
CaO	0.0650	1.0602	1.00	0.07	0.0678	281.8094	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
Si2O3	0.0190	1.1970	1.00	0.00	0.0000	291.4982	0.0000	0.0000	-53.6200	-0.0104	0.0000	-0.1644
P2O5	2.2910	2.2910	1.00	0.04	0.0427	141.9370	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
Nd2O3	1.1680	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	0.0000	325.8100	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
Y2O3	1.2698	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	-30.5700	-0.0092	0.0006	-0.1444
BaO	0.0570	1.1165	1.14	0.07	0.0710	153.3394	0.0000	0.0000	-8.8100	-0.0041	0.0009	-0.0848
PbO	0.0720	1.0772	2.11	0.16	0.1605	223.1900	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
CaO2	1.2284	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
MdO3	1.5003	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	-2.4000	-0.0070	0.0058	-0.1104
ZnO	0.1500	1.2447	2.00	0.37	0.3669	81.3800	0.0046	0.2914	1.7900	0.0031	0.0034	0.0466
CuO	0.1450	1.2518	1.18	0.22	0.2118	79.5454	0.0027	0.1721	0.0000	0.0000	0.0000	0.0000
Fe2+/Fe3+	0.0150											
SUMS	50.6290	93.9837	101.9789	100.0000	100.0000	1.5779	100.0000	SUM + pH S	-6.5891	0.4037	-72.2699	
ALK/SiO2 =	0.355			TEST=				SUM + pH B	-5.3268			
ALK+ B2O3/SiO2 =	0.464			TIME=				pH=	-7.0177			
B2O3/SiO2 =	0.109			SA/IV=					10.39			
VISCOSITY @TEMP(C)	88.602											
LIQUIDUS RATIO =	0.079											
LIQUIDUS (°C) =	954											

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FeO (OCTOBER 23, 1990)

ELEMENT	GRAV	OXIDE	FACTORS FOR	NORMOXIDE	M.W.	MOLE	NORMOXIDE	COMPONENT	FREE ENERGY	NEO	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
Al2O3	1.8895	4.2884	1.21	5.16	101.9600	0.0506	3.1887	3.0400	0.0969	-0.0638	1.5384
CaO	1.3992	0.9766	1.11	1.09	56.0784	0.0193	1.2191	-16.1160	-0.1985	0.0244	-3.1181
Fe2O3	1.4297	11.7287	1.04	12.17	159.8922	0.0762	4.8012	15.5000	0.7442	-0.0960	11.8108
FeO	1.2665	0.0739	1.00	0.07	71.8464	0.0010	0.0648	-14.6090	-0.0095	0.0013	-0.1502
MgO	1.6563	1.3449	1.12	1.51	40.3114	0.0375	2.3638	-13.8880	-0.3283	0.0473	-5.2101
MnO	1.2912	1.7612	1.31	2.31	70.8374	0.0325	2.0477	-14.8710	-0.3045	0.0410	-4.8327
Na2O	1.3480	6.5706	1.03	8.82	61.9790	0.1423	8.9672	-28.6150	-2.5638	0.1793	-41.0080
NiO	2.1525	4.3696	1.06	4.63	29.8774	0.1549	9.7587	-22.7400	-2.2181	0.1952	-35.2188
Ni2O3	1.2728	0.8554	1.18	0.78	74.7094	0.0104	0.6538	-14.3470	-0.0938	0.0131	-1.4887
SiO2	2.1393	50.4511	1.10	55.47	60.0848	0.8232	58.1709	5.5900	1.7384	0.0000	27.5573
Cr2O3	1.4816	0.2163	1.59	0.34	151.9974	0.0023	0.1427	37.3600	0.0533	-0.0029	0.8483
B2O3	3.2201	6.7236	1.02	6.84	69.8204	0.0982	9.1902	-9.9300	-0.6147	0.0000	-9.7553
Li2O	1.1344	0.0000	1.00	0.00	270.0368	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SnO	1.1828	0.0260	1.00	0.03	103.6184	0.0003	0.0158	-24.4000	-0.0039	0.0003	-0.0613
ZnO	1.3508	0.0627	1.00	0.06	123.2188	0.0000	0.0014	45.1000	0.0006	0.0001	0.0099
TiO2	1.6680	0.2635	0.98	0.26	79.8988	0.0032	0.2028	15.9900	0.0324	0.0081	0.5147
K2O	1.2046	2.6588	0.92	2.44	94.2034	0.0259	1.8348	-41.7350	-0.6823	0.0327	-10.8289
Ca2O	1.0602	0.0477	1.00	0.05	281.8084	0.0002	0.0107	-48.8200	-0.0050	0.0002	-0.0793
Sn2O3	1.1870	0.0000	1.00	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0458	1.00	0.05	141.9370	0.0003	0.0203	-53.6200	-0.1009	0.0000	-0.1731
N2O3	1.1860	0.0000	1.00	0.00	338.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2899	0.0000	1.00	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
B2O	1.1165	0.0659	1.14	0.07	153.3394	0.0005	0.0308	-30.5700	-0.0094	0.0000	-0.1495
RO	1.0772	0.0808	2.11	0.17	172.1200	0.0008	0.0481	-8.8100	-0.0042	0.0010	-0.0673
CaO2	1.2284	0.0000	1.00	0.00	223.1900	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
Mn2O3	1.5003	0.0000	1.00	0.00	143.8382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.1878	2.00	0.38	81.3600	0.0048	0.2918	-2.4000	-0.0070	0.0058	-0.1111
LiO	1.2518	0.1865	1.18	0.22	79.5454	0.0028	0.1748	1.7900	0.0031	0.0035	0.0496
Fe2+/Fe3+											
SLIMS	51.0030	94.7056	102.8479	100.0000		1.5871	100.0000	SUM + pH SI	-4.4068	0.3912	-89.8255
ALK/SiO2 =	0.350			TEST=				SUM + pH B	-5.2038		
ALK+ B2O3/SiO2 =	0.457			TIME=					-6.9594		
B2O3/SiO2 =	0.106			SA/VE=					10.44		
VISCOSITY @TEMP(C)	92.986										
LIQUIDUS RATIO =	0.081										
LIQUIDUS (C) =	961										

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	MOLE FRACTION	NORMOXIDE MOLE %	COMPONENT FREE ENERGY	FREE ENERGY HYDRATION	NEO	KCAL/KG
Al2O3	2.2490	1.8895	4.2495	1.21	5.0035	101.9600	0.0504	3.1810	3.0400	0.0967	-0.0636	1.5316
CaO	0.7010	1.3992	0.9808	1.11	1.09	56.0794	0.0194	1.2268	-16.1160	-0.1977	0.0245	-3.1315
Fe2O3	8.2420	1.4297	11.7017	1.04	11.8247	159.6922	0.0760	4.7998	15.5000	0.7440	-0.0960	11.7835
FeO	N/A	1.2865	0.0737	1.00	0.0718	71.8464	0.0010	0.0648	-14.6090	-0.0095	0.0013	-0.1499
MgO	0.8110	1.6583	1.3449	1.12	1.4730	40.3114	0.0375	2.3686	-13.8890	-0.3290	0.0474	-5.2101
MnO	1.3620	1.7562	1.7562	1.31	2.2421	76.9374	0.0325	2.0488	-14.8710	-0.3047	0.0410	-4.8257
MgO	6.4000	1.3480	8.6272	1.03	8.6481	61.9790	0.1433	9.0447	-28.8150	-2.6062	0.1809	-41.2789
N2O	2.0250	2.1525	4.3588	1.06	4.4960	29.8774	0.1545	9.7544	-22.7400	-2.2181	0.1951	-35.1320
Li2O	0.5260	1.2726	0.6694	1.18	0.79	74.7094	0.0106	0.8691	-14.3470	-0.0960	0.0134	-1.5205
SiO2	23.4920	2.1393	50.2350	1.10	55.23	60.0948	0.9192	58.0388	5.5900	1.7202	0.0000	27.2452
Cr2O3	0.1460	1.4616	0.2134	1.59	0.34	151.9874	0.0022	0.1411	37.3600	0.0527	-0.0028	0.8349
B2O3	2.0850	3.2201	6.7139	1.02	6.6523	69.6204	0.0981	6.1937	-9.9300	-0.6150	0.0000	-9.7413
Li2O	1.1344	0.0000	0.0000	1.00	0.0000	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	0.0000	1.00	0.0000	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SiO2	0.0210	1.1626	0.0248	1.00	0.02	103.6194	0.0002	0.0151	-24.4000	-0.0037	0.0003	-0.0585
ZrO2	0.0020	1.3508	0.0027	1.00	0.0028	123.2188	0.0000	0.0014	45.1000	0.0006	0.0001	0.0099
TiO2	0.1560	1.6680	0.2602	0.98	0.2473	79.8988	0.0032	0.2007	15.9900	0.0321	0.0080	0.5082
K2O	2.2330	1.2048	2.6899	0.92	2.4088	94.2034	0.0263	1.6575	-41.7350	-0.6918	0.0332	-10.9565
CaO	0.0450	1.0602	0.0477	1.00	0.0465	281.8094	0.0002	0.0107	-46.9200	-0.0050	0.0002	-0.0793
Si2O3	1.1970	0.0000	0.0000	1.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0321	0.0321	1.00	0.0312	141.9370	0.0002	0.0143	-53.9200	-0.0077	0.0000	-0.1212
N2O3	1.1660	0.0000	0.0000	1.00	0.0000	338.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	0.0000	1.00	0.0000	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	0.0000	1.00	0.0000	225.6382	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0648	0.0648	1.14	0.0718	153.3394	0.0005	0.0304	-30.5700	-0.0083	0.0006	-0.1470
PbO	1.0772	0.0948	0.0948	2.11	0.1948	223.1900	0.0008	0.0566	-8.8100	-0.0050	0.0011	-0.0789
CaO2	1.2284	0.0000	0.0000	1.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MgO3	1.5003	0.0000	0.0000	1.00	0.0000	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.2041	0.2041	2.00	0.3985	81.3600	0.0050	0.3174	-2.4000	-0.0076	0.0063	-0.1206
CaO	1.2518	0.1753	0.1753	1.18	0.2017	78.5454	0.0026	0.1644	1.7900	0.0029	0.0033	0.0466
Fe2+/Fe3+												
SUMS	50.9500	94.5233	102.6676	100.0000	100.0000	1.5639	100.0000	SUM + pH SI	-4.4578	0.3942	-70.5920	
ALK/SiO2 =	0.353			TEST=				SUM + pH B	-5.2548			
ALK+ B2O3/SiO2 =	0.459			TIME=				pH =	-7.0103			
B2O3/SiO2 =	0.167								10.44			
VISCOSITY @TEMP(C)	91.060											
LIQUIDUS RATIO =	0.882											
LIQUIDUS (°C) =	961											

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1999)

GLASS ID= HG2-2A (SPREADSHEET REVISION 8.8, OCTOBER 23, 1990)

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	FRACTION MOLE	NORMOXIDE MOLE %	COMPONENT FREEENERGY	FREEENERGY HYDRATION	NBD	KCAL/KG
Al2O3	2.4620	1.8895	1.21	5.62	5.3862	101.9600	0.0552	3.4490	3.0400	0.1048	-0.0690	1.8787
CaO	0.7420	1.3982	1.11	1.15	1.1048	56.0784	0.0206	1.2862	-16.1160	-0.2073	0.0257	-3.3147
Fe2O3	8.8710	1.4287	1.04	13.08	12.5278	159.6922	0.0819	5.1219	15.5000	0.7939	-0.1024	12.8953
FeO	N/A	0.8681	1.00	0.07	0.0652	71.8464	0.0009	0.0592	-14.6090	-0.0087	0.0012	-0.1384
MgO	0.7510	1.2454	1.12	1.40	1.3413	40.3114	0.0347	2.1724	-13.8880	-0.3017	0.0434	-4.8246
MnO	1.4610	1.8664	1.31	2.47	2.3851	70.9374	0.0348	2.1767	-14.8710	-0.3237	0.0435	-5.1764
Na2O	7.5770	10.2138	1.03	10.51	10.0682	61.9790	0.1696	10.6058	-28.9150	-3.0561	0.2121	-48.8703
Li2O	1.9780	2.1525	1.06	4.51	4.3207	29.8774	0.1510	9.4417	-22.7400	-2.1470	0.1888	-34.3340
NO	0.5210	0.6630	1.18	0.78	0.7511	74.7084	0.0105	0.6564	-14.3470	-0.0942	0.0131	-1.5060
SiO2	23.3810	50.0190	1.10	55.00	52.6749	60.0848	0.9153	57.2369	5.9900	1.8235	0.0000	25.9628
Cr2O3	0.1330	1.4618	1.59	0.31	0.2984	151.9974	0.0020	0.1273	37.3600	0.0476	-0.0025	0.7605
B2O3	1.8840	3.2201	6.0687	6.17	5.9110	69.6204	0.0886	5.5432	-9.9300	-0.5504	0.0000	-8.8022
UO2	1.1344	0.0000	1.00	0.00	0.0000	270.0368	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000
ThO2	1.1378	0.0000	1.00	0.00	0.0000	284.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SO	1.1828	0.0284	1.00	0.03	0.0272	103.6194	0.0003	0.0171	-24.4000	-0.0042	0.0003	-0.0688
ZrO2	0.1550	1.3508	1.00	0.21	0.2000	123.2188	0.0017	0.1060	45.1000	0.0478	0.0042	0.7644
TiO2	0.1420	1.6680	0.98	0.23	0.2214	79.8988	0.0029	0.1809	15.9900	0.0289	0.0072	0.4826
K2O	1.6970	1.2046	0.92	1.88	1.8002	94.2034	0.0200	1.2478	-41.7350	-0.5207	0.0250	-8.3285
Ca2O	0.0350	1.0602	1.00	0.04	0.0355	281.8094	0.0001	0.0082	-46.8200	-0.0039	0.0002	-0.0616
Si2O3	1.1970	0.0000	1.00	0.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	0.0230	2.2910	1.00	0.05	0.0505	141.9370	0.0004	0.0232	-53.6200	-0.0124	0.0000	-0.1991
Na2CO3	1.1860	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	0.0000	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2899	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0758	1.14	0.09	0.0828	153.3384	0.0008	0.0352	-30.5700	-0.0108	0.0007	-0.1723
RbO	0.1040	1.0772	2.11	0.24	0.2284	223.1900	0.0011	0.0662	-8.8100	-0.0058	0.0013	-0.0933
CaO2	1.2284	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MnCO3	1.5003	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	0.1640	1.2447	2.00	0.41	0.3918	61.3900	0.0050	0.3144	-2.4000	-0.0075	0.0063	-0.1208
ClO	1.2518	0.1339	1.18	0.16	0.1516	79.5454	0.0020	0.1244	1.7900	0.0022	0.0025	0.0358
Fe2+/Fe3+												
SLIMS	52.2810	96.0486	104.4050	100.0000	100.0000	1.5991	100.0000	SUM + pH/SI	-4.6956	0.4017	-73.8492	
ALK+/SiO2 =	0.372		TEST=					SUM + pH/B	-5.6949			
ALK+ B2O3/SiO2 =	0.468		TIME=						-7.7940			
B2O3/SiO2 =	0.897		SA/IV=						10.7			
VISCOSITY @TEMP(C)	80.753											
LIQUIDUS RATIO =	0.998											
LIQUIDUS (°C) =	986											

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

(SPREADSHEET REVISION 5.0, OCTOBER 23, 1990)

GLASS ID= HG2-28

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORM OXIDE WT%	M.W.	FRACTION MOLE	NORM OXIDE MOLE %	COMPONENT FREE ENERGY	FREE ENERGY HYDRATION	NEO	KCAL/KG
Al2O3	2.4350	1.6895	1.21	4.6009	5.56	101.9600	0.0545	3.4203	3.0400	0.1040	-0.0684	1.6593
CaO	0.7080	1.3992	1.11	0.9906	1.10	56.0794	0.0196	1.2306	-16.1160	-0.1963	0.0246	-3.1628
Fe2O3	8.6690	1.4267	1.04	12.6044	13.08	159.6922	0.0819	5.1345	15.5000	0.7958	-0.1027	12.6925
FeO	N/A	1.2865	1.00	0.0681	0.07	71.8464	0.0094	0.0594	-14.6090	-0.0087	0.0012	-0.1364
MgO	0.7540	1.6583	1.12	1.2504	1.41	40.3114	0.0349	2.1870	-13.8880	-0.3037	0.0437	-4.8439
MnO	1.4800	1.2912	1.31	1.9110	2.50	70.9374	0.0353	2.2110	-14.8710	-0.3268	0.0442	-5.2437
Na2O	7.6970	1.3480	1.03	10.3758	10.68	61.9790	0.1723	10.8028	-28.8150	-3.1128	0.2161	-48.6443
Li2O	1.9740	2.1525	1.06	4.2490	4.50	28.8774	0.1508	9.4432	-22.7400	-2.1474	0.1889	-34.2472
NO	0.5280	0.6719	1.18	0.79	0.79	74.7084	0.0108	0.6670	-14.3470	-0.0957	0.0133	-1.5263
SiO2	23.2310	2.1393	1.10	49.6981	54.64	60.0848	0.9094	67.0230	5.9900	1.6009	0.0000	25.5310
Cr2O3	0.1270	1.4616	1.59	0.1856	0.30	151.9974	0.0019	0.1219	37.3600	0.0455	-0.0024	0.7262
B2O3	1.6930	3.2201	1.02	6.0956	6.20	69.6204	0.0891	5.5847	-9.9300	-0.5548	0.0000	-8.8443
LO2	1.1344	0.0000	1.00	0.0000	0.00	270.0388	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000
SiO	1.1379	0.0000	1.00	0.0000	0.00	284.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
ZrO	1.1826	0.0272	1.00	0.0272	0.03	103.6194	0.0003	0.0165	-24.4000	-0.0040	0.0003	-0.0640
ZrO2	0.1550	1.3508	1.00	0.2094	0.21	123.2188	0.0017	0.1063	45.1000	0.0479	0.0043	0.7644
TiO2	0.1410	1.6680	0.98	0.2352	0.23	79.8688	0.0029	0.1801	15.9900	0.0288	0.0072	0.4594
K2O	1.6500	1.2046	1.02	1.9878	1.83	94.2034	0.0194	1.2163	-41.7950	-0.5076	0.0243	-8.0959
CaO	0.3350	1.0602	0.92	0.0371	0.04	281.8094	0.0001	0.0083	-46.8200	-0.0039	0.0002	-0.0616
Si2O3	1.1970	0.0000	1.00	0.0000	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0710	1.00	0.0710	0.07	141.9370	0.0005	0.0314	-53.6200	-0.0168	0.0000	-0.2883
Na2CO3	1.1660	0.0000	1.00	0.0000	0.00	336.4762	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.0000	0.00	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.0000	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
B2O	0.0870	1.1165	1.14	0.0748	0.09	153.3394	0.0006	0.0348	-30.5700	-0.0108	0.0007	-0.1898
RO	0.1120	1.0772	2.11	0.1206	0.25	223.1900	0.0011	0.0715	-8.8100	-0.0063	0.0014	-0.1005
CaO2	1.2284	0.0000	1.00	0.0000	0.00	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MgO3	1.5003	0.0000	1.00	0.0000	0.00	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZrO	0.1720	1.2447	2.00	0.2141	0.43	61.3800	0.0053	0.3306	-2.4000	-0.0079	0.0066	-0.1265
O2O	0.1020	1.2518	1.18	0.1277	0.15	79.5454	0.0019	0.1189	1.7900	0.0021	0.0024	0.0340
Fe2+/-Fe3+	0.0060											
SLMS	52.1840			95.8059	104.1464	100.0000	1.5948	100.0000	SUM + pH S	-4.6821	0.4059	-74.6719
ALK/SiO2 =		0.377			TEST=				SUM + pH B	-5.8078		
ALK+ B2O3/SiO2 =		0.474			TIME=					-7.9471		
B2O3/SiO2 =		0.698								10.73		
VISCOSITY @TEMP(°C)		77.343			SAIV=							
LIQUIDUS RATIO =		6.699										
LIQUIDUS (°C) =		938										

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

ELEMENT	GRAV	OXIDE	FACTORS FOR	NORMOXIDE	M.W.	MOLE	NORMOXIDE	COMPONENT	FREE ENERGY	NEO	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
			BIAS CORRECT								
Al2O3	2.3780	4.4895	1.21	5.43	101.9600	0.0532	3.3642	3.0400	0.1023	-0.0873	1.6181
CaO	0.7160	1.0018	1.11	1.11	56.0784	0.0198	1.2544	-16.1160	-0.2022	0.0251	-3.1985
Fe2O3	8.6220	12.2534	1.04	12.71	159.8922	0.0786	5.0315	15.5000	0.7789	-0.1006	12.3390
FeO	N/A	0.0662	1.00	0.07	71.8464	0.0009	0.0582	-14.6090	-0.0085	0.0012	-0.1345
MgO	0.7470	1.2388	1.12	1.39	40.3114	0.0346	2.1840	-13.8880	-0.3033	0.0437	-4.7989
MnO	1.4390	1.8580	1.31	2.43	70.9374	0.0343	2.1670	-14.8710	-0.3223	0.0433	-5.0985
Na2O	7.5090	10.1221	1.03	10.42	61.9790	0.1681	10.6234	-28.8150	-3.0611	0.2125	-48.4318
Li2O	1.9700	4.2404	1.06	4.49	29.8774	0.1503	9.4986	-22.7400	-2.1602	0.1900	-34.1778
NO	0.5050	0.6427	1.18	0.76	74.7094	0.0102	0.6431	-14.3470	-0.0923	0.0129	-1.4598
SiO2	23.2600	49.7801	1.10	54.71	60.0848	0.9106	57.5520	5.5900	1.6455	0.0000	26.0335
Cr2O3	0.1260	0.1842	1.59	0.29	151.9974	0.0019	0.1219	37.3600	0.0455	-0.0024	0.7205
B2O3	1.8590	5.9862	1.02	6.09	69.8204	0.0875	5.8283	-9.9300	-0.5480	0.0000	-8.6854
Li2O	1.1344	0.0000	1.00	0.00	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SrO	0.0230	0.0272	1.00	0.03	103.8194	0.0003	0.0186	-24.4000	-0.0040	0.0003	-0.0640
ZrO2	0.0510	0.0689	1.00	0.07	123.2188	0.0006	0.0353	45.1000	0.0159	0.0014	0.2515
TiO2	0.1370	0.2285	0.98	0.22	79.8988	0.0028	0.1764	15.9900	0.0282	0.0071	0.4463
K2O	1.6450	1.9816	0.92	1.82	94.2034	0.0193	1.2224	-41.7350	-0.5102	0.0244	-8.0714
Ca2O	0.0850	0.0689	1.00	0.07	281.8094	0.0002	0.0155	-48.8200	-0.0072	0.0003	-0.1145
Si2O3	1.1970	0.0000	1.00	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0000	1.00	0.00	141.9370	0.0000	0.0000	-53.8200	0.0000	0.0000	0.0000
Mn2O3	1.1660	0.0000	1.00	0.00	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2899	0.0000	1.00	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
B2O	1.1165	0.0728	1.14	0.08	153.3394	0.0005	0.0341	-30.5700	-0.0104	0.0007	-0.1647
RO	1.0772	0.0883	2.11	0.19	223.1900	0.0008	0.0528	-8.8100	-0.0046	0.0011	-0.0736
CaO2	1.2284	0.0000	1.00	0.00	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MnO3	1.5003	0.0000	1.00	0.00	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.1992	2.00	0.40	81.3900	0.0049	0.3100	-2.4000	-0.0074	0.0062	-0.1177
QO	1.2518	0.1164	1.18	0.14	79.5454	0.0017	0.1093	1.7900	0.0020	0.0022	0.0310
Fe2+/Fe3+											
SUMS	51.4500	94.6948	102.9203	100.0000	1.5821	100.0000	SUM + pH SI	-4.6235	0.4019	-73.1512	
							SUM + pH B	-5.7250			
							pH =	-7.8374			
								10.71			

ALK/SiO2 = 0.371  
 ALK+ B2O3/SiO2 = 0.467  
 B2O3/SiO2 = 0.096  
 VISCOSITY @TEMP(C) = 81.796  
 LIQUIDUS RATIO = 0.087  
 LIQUIDUS (C) = 978

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	FRACTION MOLE	NORMOXIDE MOLE %	COMPONENT FREE ENERGY	FREE ENERGY HYDRATION	NBD	KCAL/KG
Al2O3	2.3690	4.4762	1.21	5.41	5.1391	101.9600	0.0531	3.3167	3.0400	0.1008	-0.0663	1.8133
CaO	0.7060	0.9878	1.11	1.10	1.0423	56.0794	0.0196	1.2231	-16.1160	-0.1971	0.0245	-3.1539
Fe2O3	8.6450	12.2880	1.04	12.75	12.1060	159.6922	0.0798	4.9884	15.5000	0.7732	-0.0998	12.3719
FeO	N/A	0.0663	1.00	0.07	0.0630	71.8464	0.0009	0.0577	-14.8090	-0.0084	0.0012	-0.1349
MgO	0.7450	1.2354	1.12	1.39	1.3194	40.3114	0.0345	2.1538	-13.8880	-0.2991	0.0431	-4.7861
MnO	1.4340	1.8516	1.31	2.42	2.3018	70.9374	0.0342	2.1352	-14.8710	-0.3175	0.0427	-5.0908
Mn2O	7.6570	10.3216	1.03	10.62	10.0889	61.9790	0.1714	10.7114	-28.8150	-3.0865	0.2142	-49.3863
Li2O	1.9720	4.2447	1.06	4.50	4.2692	29.8774	0.1505	9.4027	-22.7400	-2.1382	0.1661	-34.2125
NaO	0.5230	0.6656	1.18	0.79	0.7477	74.7094	0.0105	0.8586	-14.3470	-0.0945	0.0132	-1.5118
SiO2	2.1393	49.7216	1.10	54.67	51.9214	60.0848	0.9099	56.8629	5.5900	1.6135	0.0000	25.8177
Cr2O3	1.4616	0.2032	1.59	0.32	0.3071	151.9874	0.0021	0.1930	37.3600	0.0497	-0.0027	0.7948
B2O3	3.2201	6.0055	1.02	6.11	5.8021	69.8204	0.0877	5.4840	-9.9300	-0.5446	0.0000	-6.7134
Li2O	1.1344	0.0000	1.00	0.00	0.0000	270.0368	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	0.0000	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SO	0.0220	0.0260	1.00	0.03	0.0247	103.6184	0.0003	0.0157	-24.4000	-0.0038	0.0003	-0.0613
ZrO2	0.0510	0.0669	1.00	0.07	0.0653	123.2188	0.0006	0.0349	45.1000	0.0157	0.0014	0.2515
TiO2	0.1360	0.2268	0.98	0.22	0.2103	79.8988	0.0028	0.1732	15.9900	0.0277	0.0069	0.4431
K2O	1.6440	1.9804	0.92	1.82	1.7293	94.2034	0.0193	1.2079	-41.7350	-0.5041	0.0242	-6.0665
Ca2O	0.0650	0.0688	1.00	0.07	0.0655	281.8094	0.0002	0.0153	-46.8200	-0.0072	0.0003	-0.1145
Sn2O3	1.1970	0.0000	1.00	0.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0000	1.00	0.00	0.0000	141.9370	0.0000	0.0000	-53.6200	0.0000	0.0000	0.0000
N2O3	1.1660	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	0.0000	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	0.0640	0.0715	1.14	0.08	0.0773	153.3394	0.0005	0.0332	-30.5700	-0.0101	0.0007	-0.1622
RbO	0.0820	0.0883	2.11	0.19	0.1770	223.1900	0.0008	0.0522	-8.8100	-0.0046	0.0010	-0.0736
CaO2	1.2264	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MnO3	1.4340	2.1514	1.00	2.15	2.0433	143.9382	0.0149	0.9341	-21.7500	-0.2032	0.0000	-3.2510
ZnO	0.1550	0.1929	2.00	0.39	0.3672	81.3800	0.0048	0.2969	-2.4000	-0.0071	0.0059	-0.1140
ClO	0.0940	0.1177	1.18	0.14	0.1321	79.5454	0.0017	0.1092	1.7900	0.0020	0.0022	0.0313
Fe2+/Fe3+	0.0060											
SUMS	53.0440	97.0565		105.2904	100.0000		1.6001	100.0000		-4.8434	0.4010	-77.4990
ALK/SiO2 =	0.375				TEST=			SUM + pH SI		-5.9449		
ALK+ B2O3/SiO2 =	0.472				TIME=			SUM + pH B		-8.0573		
B2O3/SiO2 =	0.696				SA/V=			pH =		10.71		
VISCOSITY @TEMP(°C)	78.940											
LIQUIDUS RATIO =	0.087											
LIQUIDUS (°C) =	979											

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)



ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	MOLE FRACTION	NORMOXIDE MOLE %	COMPONENT FREE ENERGY	FREE ENERGY HYDRATION	NEO	KCAL/KG
Al2O3	2.4560	1.8995	4.6406	1.21	5.61	101.9600	0.0550	3.5181	3.0400	0.1070	-0.0704	1.6726
CaO	0.7340	1.3992	1.0270	1.11	1.14	56.0794	0.0203	1.3010	-16.1160	-0.2097	0.0260	-3.2790
Fe2O3	8.6660	1.4297	12.0952	1.04	12.55	159.6922	0.0786	5.0246	15.5000	0.7786	-0.1005	12.1797
FeO	N/A	1.2865	0.5224	1.00	0.52	71.8464	0.0073	0.4850	-14.8090	-0.0679	0.0093	-1.0623
MgO	0.7950	1.6583	1.2189	1.12	1.37	40.3114	0.0340	2.1741	-13.8880	-0.3019	0.0435	-4.7219
MnO	1.4720	1.9006	1.31	2.49	2.4342	70.9374	0.0351	2.2428	-14.8710	-0.3335	0.0449	-5.2154
Na2O	7.2720	1.3480	9.8027	1.03	10.09	61.9790	0.1628	10.4083	-28.8150	-2.9992	0.2082	-46.9031
Li2O	1.9160	2.1525	4.1242	1.06	4.37	29.8774	0.1462	9.3472	-22.7400	-2.1255	0.1869	-33.2410
NO	0.5910	1.2726	0.6758	1.18	0.80	74.7094	0.0107	0.8841	-14.3470	-0.0981	0.0137	-1.5349
SiO2	22.7000	2.1393	48.5821	1.10	53.39	60.0948	0.8888	56.8228	5.5900	1.5815	0.0000	24.7327
Cr2O3	1.4616	1.4616	0.2046	1.58	0.33	151.9874	0.0021	0.1370	37.3600	0.0512	-0.0027	0.8006
B2O3	3.2201	6.0216	6.0216	1.02	6.13	69.8204	0.0880	5.8280	-9.9300	-0.5587	0.0000	-8.7368
UO2	1.1344	0.0000	0.0000	1.00	0.00	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1378	0.0000	0.0000	1.00	0.00	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
ZrO2	1.1826	0.0272	0.1756	1.00	0.18	123.2188	0.0014	0.0809	45.1000	0.0410	0.0036	0.8411
TiO2	1.3508	0.2352	0.2352	0.98	0.23	79.8988	0.0029	0.1837	15.9900	0.0294	0.0073	0.4594
K2O	1.6680	2.1878	2.1878	0.92	2.01	94.2034	0.0214	1.3652	-41.7350	-0.5698	0.0273	-8.9104
CeO	1.2048	1.0602	0.0318	1.00	0.03	281.8094	0.0001	0.0072	-48.8200	-0.0034	0.0001	-0.0528
SrO	1.1970	0.0000	0.0000	1.00	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0527	0.0527	1.00	0.05	141.9370	0.0004	0.0237	-53.6200	-0.0127	0.0000	-0.1991
Nb2O5	1.1660	0.0000	0.0000	1.00	0.00	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	0.0000	1.00	0.00	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2899	0.0000	0.0000	1.00	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0737	0.1165	1.14	0.08	153.3394	0.0005	0.0350	-30.3700	-0.0107	0.0007	-0.1672
RbO	1.0772	0.1099	0.1099	2.11	0.23	223.1900	0.0010	0.0664	-8.8100	-0.0059	0.0013	-0.0915
CeO2	1.2284	0.0000	0.0000	1.00	0.00	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MgO3	1.5003	0.0000	0.0000	1.00	0.00	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.2079	0.2079	2.00	0.42	81.3800	0.0051	0.3273	-2.4000	-0.0079	0.0065	-0.1229
UO	0.1120	1.2518	0.1402	1.18	0.17	78.5454	0.0021	0.1332	1.7900	0.0024	0.0027	0.0373
Fe2+/Fe3+	0.0480											
SUMS	51.3020	94.0373	102.2054	100.0000	100.0000	1.5639	100.0000	SUM + pH SI	-4.7177	0.4089	-73.7790	
ALK/SiO2 =	0.372							SUM + pH B	-5.8434			
ALK+ B2O3/SiO2 =	0.471							pH =	-7.9827			
B2O3/SiO2 =	0.099								10.73			
VISCOSITY @TEMP(C)	79.802											
LIQUIDUS RATIO =	0.089											
LIQUIDUS (C) =	997											

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

ELEMENT	GRAV FACTOR	OXIDE WT%	FACTORS FOR BIAS CORRECT	BIAS CORRECT	BIAS CORRECT OXIDE WT%	NORMOXIDE WT%	M.W.	MOLE FRACTION	NORMOXIDE MOLE %	COMPONENT FREE ENERGY	FREE ENERGY HYDRATION	NBO	KCAL/KG
Al2O3	2.4690	1.8995	4.6652	1.21	5.64	5.5129	101.9600	0.0553	3.5345	3.0400	0.1074	-0.0707	1.6814
CaO	0.7290	1.3992	1.0200	1.11	1.13	1.1078	56.0784	0.0202	1.2913	-16.1160	-0.2081	0.0258	-3.2566
Fe2O3	8.6510	1.4297	12.0747	1.04	12.53	12.2460	159.6922	0.0764	5.0129	15.5000	0.7770	-0.1003	12.1591
FeO	N/A	1.2665	0.5215	1.00	0.52	0.5098	71.8464	0.0073	0.4639	-14.6090	-0.0678	0.0093	-1.0605
MgO	0.7350	1.6583	1.2189	1.12	1.37	1.3398	40.3114	0.0340	2.1727	-13.8880	-0.3017	0.0435	-4.7219
MnO	1.4610	1.2912	1.8664	1.31	2.47	2.4138	70.9374	0.0348	2.2244	-14.8710	-0.3308	0.0445	-5.1764
Na2O	7.2930	1.3490	9.8310	1.03	10.12	9.8908	61.9790	0.1632	10.4318	-28.8150	-3.0059	0.2086	-47.0366
Li2O	1.9110	2.1525	4.1134	1.06	4.36	4.2593	29.6774	0.1458	9.3169	-22.7400	-2.1187	0.1863	-33.1542
NO	0.5350	1.2726	0.6808	1.18	0.81	0.7872	74.7094	0.0108	0.6868	-14.3470	-0.0988	0.0138	-1.5465
SiO2	22.6930	2.1393	48.5471	1.10	53.38	52.1790	60.0848	0.8884	56.7691	5.5900	1.5777	0.0000	24.6889
Cr2O3	0.1430	1.4816	0.2090	1.59	0.33	0.3252	151.9974	0.0022	0.1399	37.3600	0.0523	-0.0028	0.8177
B2O3	1.8820	3.2201	6.0602	1.02	6.16	6.0264	69.6204	0.0885	5.6585	-9.9300	-0.5619	0.0000	-8.7929
UO2	1.1344	0.0000	0.0000	1.00	0.00	0.0000	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	0.0000	1.00	0.00	0.0000	284.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
ZrO2	0.0230	1.1628	0.1756	1.00	0.18	0.1712	123.2188	0.0014	0.0808	45.1000	0.0410	0.0038	-0.0640
TiO2	0.1440	1.6680	2.2402	0.98	2.03	2.2292	79.8988	0.0029	1.3786	-41.7350	-0.5754	0.0075	0.4691
K2O	1.8350	1.2046	2.2104	0.92	2.03	1.9887	94.2034	0.0216	1.3786	-41.7350	-0.5754	0.0276	-9.0037
CaO	0.0300	1.0602	0.0318	1.00	0.03	0.0311	291.8094	0.0001	0.0072	-46.8200	-0.0034	0.0001	-0.0528
Sn2O3	1.1970	0.0000	0.0000	1.00	0.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0344	0.0344	1.00	0.03	0.0336	141.8370	0.0002	0.0155	-53.6200	-0.0083	0.0000	-0.1298
Na2O3	1.1660	0.0000	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	0.0000	1.00	0.00	0.0000	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0748	0.1099	1.14	0.09	0.0832	153.3394	0.0008	0.0355	-30.5700	-0.1008	0.0007	-0.1698
RbO	0.1020	1.0772	0.1099	2.11	0.23	0.2266	223.1900	0.0010	0.0684	-8.8100	-0.0058	0.0013	-0.0915
CaO2	1.2284	0.0000	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
Mn2O3	1.5003	0.0000	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.2290	0.2290	2.00	0.46	0.4487	81.3600	0.0058	0.3604	-2.4000	-0.0087	0.0072	-0.1354
QO	0.1150	1.2518	0.1440	1.16	0.17	0.1663	79.5454	0.0021	0.1367	1.7900	0.0024	0.0027	0.0383
Fe2+/Fe3+	0.0480												
SUMS	51.3470	94.1056	102.2959	100.0000	102.2959	100.0000	100.0000	1.5649	100.0000		-4.7224	0.4092	-73.8988

ALK/SiO2 = 0.372  
 ALK+ B2O3/SiO2 = 0.472  
 B2O3/SiO2 = 0.100  
 VISCOSITY @TEMP(C) = 79.697  
 LIQUIDUS RATIO = 0.089  
 LIQUIDUS (C) = 997

SUM + pH SI = -5.8461  
 SUM + pH B = -7.9874  
 pH = 10.73

TEST =  
 TIME =  
 SA/IV =

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

ELEMENT	GRAV	OXIDE	FACTORS FOR	BIAS CORRECT	NORMOXIDE	M.W.	MOLE	NORMOXIDE	COMPONENT	FREE ENERGY	NEO	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	OXIDE WT%	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
Al2O3	2.5110	1.8895	1.21	5.74	5.5807	101.9600	0.0563	3.5914	3.0400	0.1092	-0.0718	1.7100
CaO	0.7570	1.3992	1.11	1.18	1.1450	56.0794	0.0210	1.3397	-16.1160	-0.2159	0.0268	-3.3817
Fe2O3	9.2210	1.4297	1.04	12.79	12.4494	159.6922	0.0801	5.1153	15.5000	0.7929	-0.1023	12.4165
FeO	N/A	1.2865	1.00	0.7657	0.7450	71.8464	0.0107	0.8804	-14.6090	-0.0994	0.0136	-1.5570
MgO	0.7280	1.6583	1.12	1.36	1.3209	40.3114	0.0337	2.1501	-13.8980	-0.2986	0.0430	-4.6769
MnO	1.5260	1.9704	1.31	2.58	2.5095	70.9374	0.0364	2.3213	-14.8710	-0.3452	0.0464	-5.4067
Na2O	7.1490	1.3480	1.03	8.92	9.6504	61.9790	0.1600	10.2166	-28.8150	-2.9439	0.2043	-46.1098
Li2O	2.1525	1.8800	1.06	4.29	4.1698	29.8774	0.1434	9.1575	-22.7400	-2.0824	0.1832	-32.8164
NiO	0.5240	1.2726	1.18	0.79	0.7675	74.7094	0.0108	0.8741	-14.3470	-0.0967	0.0135	-1.5147
SiO2	22.5960	2.1393	1.10	53.15	51.7154	60.0848	0.8846	56.4757	5.5900	1.5579	0.0000	24.4004
Cr2O3	0.1270	1.4816	1.59	0.30	0.2875	151.9974	0.0019	0.1241	37.3600	0.0464	-0.0025	0.7262
B2O3	3.2201	6.1407	1.02	6.25	6.0782	69.6204	0.0897	5.7285	-9.9300	-0.5688	0.0000	-8.9097
UO2	1.1344	0.0000	1.00	0.00	0.0000	270.0388	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	0.0000	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SnO	0.0250	1.1826	1.00	0.03	0.0288	103.6194	0.0003	0.0182	-24.4000	-0.0044	0.0004	-0.0696
ZrO2	0.1290	1.3508	1.00	0.17	0.1691	123.2188	0.0014	0.0901	45.1000	0.0406	0.0036	0.8362
TiO2	0.1500	1.6680	0.98	0.24	0.2376	79.8988	0.0031	0.1951	15.9900	0.0312	0.0078	0.4867
K2O	2.0500	1.2046	0.92	2.27	2.2092	94.2034	0.0241	1.5388	-41.7350	-0.8422	0.0308	-10.0586
Ce2O	0.0400	1.0602	1.00	0.04	0.0413	281.8094	0.0002	0.0096	-46.8200	-0.0045	0.0002	-0.0705
Sn2O3	0.0100	1.1970	1.00	0.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0229	1.00	0.02	0.0223	141.9370	0.0002	0.0103	-53.6200	-0.0055	0.0000	-0.0665
Nb2O3	0.0030	1.1660	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0035	1.00	0.00	0.0034	325.8100	0.0000	0.0007	-31.1400	-0.0002	0.0000	-0.0034
Y2O3	1.2699	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0782	1.14	0.09	0.0866	153.3394	0.0006	0.0370	-30.5700	-0.0113	0.0007	-0.1774
RbO	1.0772	0.1034	2.11	0.22	0.2123	223.1900	0.0010	0.0624	-8.8100	-0.0055	0.0012	-0.0661
CeO2	1.2284	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
Mn2O3	1.5003	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	-2.4000	-0.0075	0.0000	0.0000
ZnO	1.2447	6.1979	2.00	0.40	0.3859	81.3800	0.0049	0.3112	1.7900	0.0027	0.0062	-0.1170
UO	1.2518	0.1802	1.18	0.19	0.1842	79.5454	0.0024	0.1520	0.0000	0.0000	0.0030	0.0428
Fe2+/Fe3+												
SLIMS	51.7860	94.6277	102.7718	100.0000	100.0000	1.5663	100.0000	SUM + pH SI	-4.7514	0.4082	-74.4192	
			TEST=					SUM + pH B	-5.8528			
			TIME=					pH=	-7.9653			
			SA/VE=						10.71			

ALK/SIO2 = 0.370  
 ALK+ B2O3/SIO2 = 0.472  
 B2O3/SIO2 = 0.101  
 VISCOSITY @TEMP(C) = 78.459  
 LIQUIDUS RATIO = 0.091  
 LIQUIDUS (C) = 1012

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

GLASS ID= HG3-2B (SPREADSHEET REVISION 5.6, OCTOBER 23, 1990)

GLASS ID= HG3-2B

ELEMENT	GRAV	OXIDE	FACTORS FOR	BIAS CORRECT	NORM. OXIDE	M.W.	MOLE	NORM. OXIDE	COMPONENT	FREE ENERGY	NBD	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	OXIDE WT%	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
Al2O3	1.8895	4.7399	1.21	5.73	5.5665	101.9600	0.0562	3.5924	3.0400	0.1089	-0.0716	1.7080
CaO	1.3992	1.0662	1.11	1.18	1.1510	56.0784	0.0211	1.3468	-16.1160	-0.2171	0.0269	-3.4040
Fe2O3	1.4297	12.4447	1.04	12.91	12.5459	159.6922	0.0808	5.1552	15.5000	0.7991	-0.1031	12.9316
FeO	1.2865	0.7727	1.00	0.77	0.7508	71.8464	0.0108	0.6657	-14.6090	-0.1002	0.0137	-1.5711
MgO	1.6583	1.2205	1.12	1.37	1.3336	40.3114	0.0340	2.1709	-13.8880	-0.3015	0.0434	-4.7283
MnO	1.5400	1.2912	1.31	2.60	2.5292	70.9374	0.0367	2.3395	-14.8710	-0.3479	0.0468	-5.4563
Na2O	1.3480	9.5182	1.03	9.60	9.5188	61.9790	0.1561	10.0777	-28.8150	-2.9039	0.2016	-45.5422
Li2O	1.8950	4.0790	1.06	4.32	4.1974	29.8774	0.1446	9.2186	-22.7400	-2.0963	0.1844	-32.8768
NiO	1.2726	0.6656	1.18	0.79	0.7650	74.7094	0.0105	0.6719	-14.3470	-0.0964	0.0134	-1.5118
SiO2	22.6250	46.4017	1.10	53.22	51.7121	60.0948	0.8957	56.4745	5.5900	1.5615	0.0000	24.4996
Cr2O3	1.4616	0.1769	1.59	0.26	0.2735	151.9674	0.0019	0.1181	37.3600	0.0441	-0.0024	0.9919
B2O3	3.2201	6.1987	1.02	6.31	6.1273	69.6204	0.0906	5.7751	-9.9300	-0.5735	0.0000	-6.9938
UO2	1.1344	0.0000	1.00	0.00	0.0000	270.0388	0.0000	0.0000	-6.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	0.0000	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SnO	1.1826	0.0296	1.00	0.03	0.0287	103.6194	0.0003	0.0182	-24.4000	-0.0044	0.0004	-0.0696
ZrO2	1.3508	0.1743	1.00	0.17	0.1689	123.2188	0.0014	0.0899	45.1000	0.0406	0.0036	0.6362
TiO2	1.6880	0.2502	0.98	0.24	0.2373	79.8988	0.0031	0.1949	15.9900	0.0312	0.0078	0.4867
K2O	1.2046	2.4020	0.92	2.21	2.1459	94.2034	0.0234	1.4948	-41.7350	-0.8238	0.0299	-9.7838
Ca2O	1.0602	0.0424	1.00	0.04	0.0412	281.8094	0.0002	0.0096	-46.8200	-0.0045	0.0002	-0.0705
P2O5	2.2910	0.0252	1.00	0.03	0.0245	141.9370	0.0002	0.0113	-53.8200	-0.0061	0.0000	-0.0952
Na2O3	1.1660	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	0.0000	326.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.00	0.0000	225.9082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0770	1.14	0.09	0.0852	153.3394	0.0008	0.0365	-30.5700	-0.0111	0.0007	-0.1748
RO	0.1040	0.1120	2.11	0.24	0.2297	223.1900	0.0011	0.0675	-6.8100	-0.0059	0.0014	-0.0933
CaO2	1.2284	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MgO3	1.5003	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.1992	2.00	0.40	0.3878	81.3800	0.0049	0.3127	-2.4000	-0.0075	0.0063	-0.1177
Q.O	1.2518	0.1565	1.18	0.18	0.1797	79.5454	0.0023	0.1462	1.7900	0.0027	0.0030	0.0416
Fe2+/Fe3+												
SUMS	51.8080	94.7397	102.9103	100.0000	100.0000	1.5683	100.0000				0.4062	-73.9015

ALK/SiO2 = 0.368  
 ALK+ B2O3/SiO2 = 0.471  
 B2O3/SiO2 = 0.102  
 VISCOSITY @TEMP(C) = 78.946  
 LIQUIDUS RATIO = 0.092  
 LIQUIDUS (°C) = 1914

SUM + pH SI  
 SUM + pH B  
 PH = 10.71

TEST =  
 TIME =  
 SA/V =

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

GLASS ID= HG3-3A (SPREADSHEET REVISION 5.6, OCTOBER 23, 1990)

ELEMENT	GRAV	OXIDE	FACTORS FOR	NORMOXIDE	M.W.	MOLE	NORMOXIDE	COMPONENT	FREE ENERGY	NEO	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
Al2O3	2.5300	4.7804	1.21	5.78	101.9600	0.0567	3.5927	3.0400	0.1062	-0.0719	1.7230
CaO	0.7950	1.1124	1.11	1.24	56.0764	0.0220	1.3969	-16.1160	-0.2251	0.0279	-3.6515
Fe2O3	9.3100	12.1336	1.04	12.59	159.6922	0.0788	4.9969	15.5000	0.7745	-0.0999	12.2184
FeO	N/A	1.0591	1.00	1.06	71.8464	0.0147	0.9344	-14.6090	-0.1365	0.0167	-2.1535
MgO	0.7420	1.2305	1.12	1.38	40.3114	0.0343	2.1758	-13.8680	-0.3022	0.0435	-4.7668
MnO	1.5560	2.0091	1.31	2.63	70.9374	0.0371	2.3500	-14.8710	-0.3495	0.0470	-5.5130
Ni2O3	6.9670	9.3915	1.03	9.67	91.9790	0.1559	9.8654	-22.8150	-2.8485	0.1977	-44.9959
NiO	1.9070	4.1048	1.06	4.35	29.8774	0.1455	9.2227	-20.7400	-2.0972	0.1945	-33.0848
SiO2	0.5250	0.6681	1.18	0.79	74.7084	0.0106	0.6705	-14.3470	-0.0962	0.0134	-1.5176
SiO2	22.6950	48.5514	1.10	53.38	60.0848	0.8884	56.3160	5.5900	1.5407	0.0000	24.3048
Cr2O3	0.1050	1.4616	1.59	0.24	151.9974	0.0016	0.1019	37.3600	0.0391	-0.0020	0.6004
B2O3	1.9680	6.3307	1.02	6.44	69.6204	0.0925	5.8636	-9.9300	-0.5623	0.0000	-9.1853
UO2	1.1344	0.0000	1.00	0.00	270.0388	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SrO	1.1628	0.0296	1.00	0.03	103.6194	0.0003	0.0181	-24.4000	-0.0044	0.0004	-0.0696
ZrO2	0.1170	0.1580	1.00	0.16	123.2188	0.0013	0.0811	45.1000	0.0368	0.0032	0.5770
TiO2	0.1550	0.2585	0.98	0.25	79.8988	0.0032	0.2002	15.9900	0.0320	0.0080	0.5050
K2O	2.1270	2.5622	0.92	2.36	94.2034	0.0250	1.5851	-41.7350	-0.6616	0.0317	-10.4364
CaO	0.0100	0.0106	1.00	0.01	281.8094	0.0000	0.0024	-46.6200	-0.0011	0.0000	-0.0176
Sr2O3	1.1970	0.0000	1.00	0.00	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0344	1.00	0.03	141.9370	0.0002	0.0153	-53.6200	-0.0082	0.0000	-0.1298
Nd2O3	1.1660	0.0000	1.00	0.00	336.4762	0.0000	0.0000	-26.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.00	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
BaO	1.1165	0.0759	1.14	0.09	153.3394	0.0008	0.0357	-30.5700	-0.0109	0.0007	-0.1723
PbO	1.0772	0.1077	2.11	0.23	223.1900	0.0010	0.0645	-8.8100	-0.0057	0.0013	-0.0897
CaO2	1.2284	0.0000	1.00	0.00	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
Mn2O3	1.5003	0.0000	1.00	0.00	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.2141	2.00	0.43	61.3800	0.0053	0.3342	-2.4000	-0.0080	0.0067	-0.1265
Fe2+/Fe3+	1.2518	0.1640	1.18	0.19	79.5454	0.0024	0.1544	1.7900	0.0028	0.0031	0.0436
SUMS	52.0180	95.1400	103.3205	100.0000	1.5775	100.0000	SUM + pHSi	-4.8636	0.4140	-75.7765	
					1.4988		SUM + pHB	-5.8726			
							pH =	-7.9490			
								10.683			

ALK/SIO2 = 0.367  
 ALK+ B2O3/SIO2 = 0.472  
 B2O3/SIO2 = 0.104  
 VISCOSITY @TEMP(C) = 79.361  
 LIQUIDUS RATIO = 0.096  
 LIQUIDUS (C) = 1014

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1990)

GLASS ID= HG3-3B (SPREADSHEET REVISION 5.0, OCTOBER 23, 1998)

ELEMENT	GRAV	OXIDE	FACTORS FOR	BIAS CORRECT	NORMOXIDE	M.W.	MOLE	NORMOXIDE	COMPONENT	FREE ENERGY	NEO	KCAL/KG
WT%	FACTOR	WT%	BIAS CORRECT	OXIDE WT%	WT%		FRACTION	MOLE %	FREE ENERGY	HYDRATION		
Al2O3	1.6695	4.7916	1.21	5.79	5.6259	101.9600	0.0588	3.6070	3.0400	0.1097	-0.0721	1.7271
CaO	1.3992	1.1152	1.11	1.24	1.2033	56.0794	0.0221	1.4027	-16.1160	-0.2281	0.0281	-3.5604
Fe2O3	1.4297	11.9289	1.04	12.36	12.0201	159.6922	0.0775	4.9205	15.5000	0.7827	-0.0984	12.0123
FeO	1.2865	1.0412	1.00	1.04	1.0113	71.8464	0.0145	0.9201	-14.6090	-0.1344	0.0184	-2.1172
MgO	1.6583	1.2305	1.12	1.36	1.3438	40.3114	0.0343	2.1792	-13.8680	-0.3028	0.0436	-4.7688
MnO	1.2912	1.8077	1.31	2.37	2.2981	70.9374	0.0334	2.1178	-14.8710	-0.3149	0.0424	-4.9603
Na2O	1.3480	9.4643	1.03	9.74	9.4602	61.9790	0.1572	9.9779	-22.8150	-2.8751	0.1998	-45.2842
Li2O	2.1525	4.1091	1.06	4.35	4.2263	28.8774	0.1456	9.2471	-22.7400	-2.1028	0.1849	-33.1195
NiO	1.2726	0.6630	1.18	0.78	0.7617	74.7094	0.0105	0.8665	-14.3470	-0.0956	0.0133	-1.5060
SiO2	2.1393	49.6691	1.10	53.51	51.9724	60.0848	0.8908	56.5444	5.5900	1.5811	0.0000	24.5893
Cr2O3	1.4616	0.1549	1.59	0.25	0.2395	151.9974	0.0018	0.1030	37.3600	0.0385	-0.0021	0.6081
B2O3	3.2201	6.3148	1.02	6.42	6.2369	69.6204	0.0923	5.8580	-9.9300	-0.5817	0.0000	-9.1620
UO2	1.1344	0.0000	1.00	0.00	0.0000	270.0388	0.0000	0.0000	-8.8000	0.0000	0.0000	0.0000
ThO2	1.1379	0.0000	1.00	0.00	0.0000	264.0400	0.0000	0.0000	-2.5330	0.0000	0.0000	0.0000
SiO	1.1826	0.0296	1.00	0.03	0.0287	103.6194	0.0003	0.0181	-24.4000	-0.0644	0.0004	-0.0696
ZrO2	1.3508	0.1580	1.00	0.16	0.1531	123.2188	0.0013	0.0812	45.1000	0.0368	0.0032	0.5770
TiO2	1.6680	0.2502	0.98	0.24	0.2372	79.8988	0.0031	0.1940	15.9900	0.0310	0.0078	0.4887
K2O	1.2046	2.5441	0.92	2.34	2.2718	94.2034	0.0248	1.5765	-41.7350	-0.6579	0.0315	-10.3628
CaO	1.0602	0.0106	1.00	0.01	0.0103	281.8094	0.0000	0.0024	-46.8200	-0.0011	0.0000	-0.0176
Sn2O3	1.1970	0.0000	1.00	0.00	0.0000	291.4982	0.0000	0.0000	0.0000	0.0000	0.0000	0.0000
P2O5	2.2910	0.0000	1.00	0.00	0.0000	141.9370	0.0000	0.0000	-53.6200	0.0000	0.0000	0.0000
Mg2O3	1.1680	0.0000	1.00	0.00	0.0000	336.4782	0.0000	0.0000	-28.3200	0.0000	0.0000	0.0000
La2O3	1.1728	0.0000	1.00	0.00	0.0000	325.8100	0.0000	0.0000	-31.1400	0.0000	0.0000	0.0000
Y2O3	1.2699	0.0000	1.00	0.00	0.0000	225.8082	0.0000	0.0000	-12.9500	0.0000	0.0000	0.0000
B2O	1.1165	0.0782	1.14	0.09	0.0864	153.3394	0.0006	0.0368	-30.5700	-0.0113	0.0007	-0.1774
RO	1.0772	0.1034	2.11	0.22	0.2119	223.1900	0.0010	0.0621	-8.8100	-0.0055	0.0012	-0.0861
CaO2	1.2284	0.0000	1.00	0.00	0.0000	172.1200	0.0000	0.0000	11.3450	0.0000	0.0000	0.0000
MgO3	1.5003	0.0000	1.00	0.00	0.0000	143.9382	0.0000	0.0000	-21.7500	0.0000	0.0000	0.0000
ZnO	1.2447	0.2104	2.00	0.42	0.4095	81.3600	0.0052	0.3289	-2.4000	-0.0079	0.0066	-0.1243
CaO	1.2518	0.1652	1.18	0.20	0.1896	79.5454	0.0025	0.1558	1.7900	0.0028	0.0031	0.0439
Fe2+/Fe3+												
SUMS	51.7770	94.8400		102.9606	100.0000		1.5750	100.0000		-4.7799	0.4122	-75.2708
ALK/SiO2 =	0.368			TEST=				SUM + pH SI		-5.8480		
ALK+ B2O3/SiO2 =	0.472			TIME=				SUM + pH B		-7.9244		
B2O3/SiO2 =	0.104			SA/V =				pH =		10.683		
VISCOSITY @TEMP(C)	80.313											
LIQUIDUS RATIO =	0.088											
LIQUIDUS (°C) =	1008											

NOTE LIQUIDUS AND VISCOSITY FORMULAS CHANGED TO CALCULATE ALL FE AS FE2O3 (OCTOBER 23, 1998)

## APPENDIX II Leachate Multielement Standard Analyses

Element	Concentration of Multielement Standard ug/mL)	Run before Samples-1 ug/mL	Run before Samples-2 ug/mL	Run after 5 Samples-1 ug/mL	Run after 5 Samples-2 ug/mL	Run after 10 Samples-1 ug/mL	Run after 10 Samples-2 ug/mL	Run after 15 Samples-1 ug/mL	Run after 15 Samples-2 ug/mL
Na (ICP)	50.00	49.72	49.86	50.48	50.14	49.59	49.62	48.97	49.40
Na (AA)		51.82	52.32	52.16	52.82	51.49	51.49	49.51	51.00
Si	50.10	48.41	48.67	49.79	49.06	48.78	48.54	48.25	48.42
B	20.10	20.23	20.29	20.57	20.38	20.17	20.17	19.82	19.97
K	10.00	9.99	9.92	9.74	10.00	9.85	9.66	9.24	9.77
Li	10.00	10.01	10.03	10.16	10.08	9.97	9.98	9.79	9.88
Al	4.00	3.64	3.66	3.72	3.72	3.64	3.65	3.62	3.63
Fe	4.00	3.92	3.94	3.99	4.00	3.92	3.92	3.91	3.93

Element	Concentration of Multielement Standard ug/mL	Average Bias before Leachate Analysis ug/mL	Average Bias after 5 Leachate Analyses ug/mL	Average Bias after 10 Leachate Analyses ug/mL	Average Bias after 15 Leachate Analyses ug/mL
Na (ICP)	50.00	-0.21	0.31	-0.40	-0.82
Na (AA)		2.07	2.49	1.49	0.26
Si	50.10	-1.56	-0.68	-1.44	-1.77
B	20.10	0.16	0.38	0.07	-0.21
K	10.00	-0.05	-0.13	-0.25	-0.50
Li	10.00	0.02	0.12	-0.02	-0.17
Al	4.00	-0.35	-0.28	-0.36	-0.38
Fe	4.00	-0.07	0.00	-0.08	-0.08

Element	Concentration of Multielement Standard ug/mL)	Run after 20 Samples-1 ug/mL	Run after 20 Samples-2 ug/mL	Run after 25 Samples-1 ug/mL	Run after 25 Samples-2 ug/mL	ELEMENT TOTAL AVERAGE BIAS ug/mL
Na (ICP)	50.00	49.43	49.31	49.69	49.48	Na (ICP) -0.36
Na (AA)		51.33	51.00	51.82	50.00	Na (AA) 1.40
Si	50.10	48.79	48.41	48.34	48.14	Si -1.47
B	20.10	19.90	19.84	19.93	19.85	B -0.01
K	10.00	9.87	10.35	9.9	10.03	K -0.14
Li	10.00	9.88	9.85	9.94	9.89	Li -0.05
Al	4.00	3.58	3.60	3.55	3.56	Al -0.37
Fe	4.00	3.91	3.96	3.9	3.9	Fe -0.07

Element	Concentration of Multielement Standard ug/mL	Average Bias after 20 Leachate Analyses ug/mL	Average Bias after 25 Leachate Analyses ug/mL
Na (ICP)	50.00	-0.63	-0.42
Na (AA)		1.16	0.91
Si	50.10	-1.50	-1.86
B	20.10	-0.23	-0.21
K	10.00	0.11	-0.04
Li	10.00	-0.14	-0.09
Al	4.00	-0.41	-0.45
Fe	4.00	-0.07	-0.10

**This page intentionally left blank.**



### APPENDIX III Leachate Data for the IDMS Hg Campaigns 1, 2, 3

GLASS ANALYSIS ID'S	ADS #	NOTEBOOK/REF #	CALCULATED	AVG. CALC.	CALCULATED	CALCULATED	AVG. CALC.
			FREE ENERGY HYDRATION w/o pH	FREE ENERGY HYDRATION w/o pH	FREE ENERGY HYDRATION SI DISS. pH A + B	FREE ENERGY HYDRATION DISS. pH	FREE ENERGY HYDRATION WITH pH
HG-1-1A	200057660	DPSTN-4771	-4.62		-5.48	-7.20	
HG-1-1B	200057661	DPSTN-4771	-4.82		-5.50	-7.25	
HG-1-1-AVG				-4.72	-5.52	-7.29	-7.25
HG-1-2A	200057663	DPSTN-4771	-4.75		-5.34	-7.05	
HG-1-2B	200057664	DPSTN-4771	-4.42		-5.33	-7.02	
HG-1-2-AVG				-4.58	-5.32	-7.00	-7.02
HG-1-3-7-1	200057665	DPSTN-4771	-4.54		-5.31	-7.00	
HG-1-3-7-2	200057666	DPSTN-4771	-4.58		-5.31	-7.00	
HG-1-3-AVG				-4.56	-5.31	-7.00	-7.00
HG-2-1-7-1	200057668	DPSTN-4771	-4.41		-5.23	-6.98	
HG-2-1-7-2	200057669	DPSTN-4771	-4.46		-5.22	-6.96	
HG-2-1-AVG				-4.43	-5.23	-6.98	-6.98
HG-2-2-7-1	200057671	DPSTN-4771	-4.61		-5.73	-7.83	
HG-2-2-7-2	200057673	DPSTN-4771	-4.68		-5.78	-7.93	
HG-2-2-AVG				-4.64	-5.78	-7.93	-7.90
HG-2-3-7-1	200057675	DPSTN-4771	-4.62		-5.83	-7.95	
HG-2-3-7-2	200057676	DPSTN-4771	-4.84		-5.85	-7.97	
HG-2-3-AVG				-4.73	-5.81	-7.90	-7.94
HG-3-1-7-1	200057677	DPSTN-4771	-4.72		-5.83	-7.96	
HG-3-1-7-2	200057678	DPSTN-4771	-4.72		-5.85	-7.98	
HG-3-1-AVG				-4.72	-5.86	-8.01	-7.98
HG-3-2-7-1	200057679	DPSTN-4771	-4.75		-5.82	-7.92	
HG-3-2-7-2	200057680	DPSTN-4771	-4.71		-5.82	-7.92	
HG-3-2-AVG				-4.73	-5.86	-8.00	-7.95
HG-3-3-7-1	200057791	DPSTN-4771	-4.80		-5.86	-7.93	
HG-3-3-7-2	200057682	DPSTN-4771	-4.78		-5.86	-7.93	
HG-3-3-AVG				-4.79	-5.87	-7.95	-7.94

<b>GLASS ANALYSIS ID'S</b>	<b>B WT. % IN GLASS</b>	<b>LI WT.% IN GLASS</b>	<b>Na WT. % IN GLASS</b>	<b>SI WT. % IN GLASS</b>	<b>K WT.% IN GLASS</b>
HG-1-1A	2.17	2.08	6.78	25.33	2.22
HG-1-1B	2.21	2.03	6.85	23.53	2.22
HG-1-1-AVG	2.19	2.05	6.82	24.43	2.22
HG-1-2A	2.16	2.15	6.61	25.86	2.91
HG-1-2B	2.17	2.15	6.43	25.89	2.18
HG-1-2-AVG	2.17	2.15	6.52	25.87	2.55
HG-1-3-7-1	2.18	2.17	6.59	25.92	2.15
HG-1-3-7-2	2.16	2.34	6.56	25.71	2.17
HG-1-3-AVG	2.17	2.26	6.57	25.81	2.16
HG-2-1-7-1	2.13	2.15	6.55	25.94	2.03
HG-2-1-7-2	2.13	2.15	6.59	25.83	2.05
HG-2-1-AVG	2.13	2.15	6.57	25.89	2.04
HG-2-2-7-1	1.92	2.10	7.80	25.72	1.56
HG-2-2-7-2	1.93	2.09	7.93	25.55	1.52
HG-2-2-AVG	1.93	2.10	7.87	25.63	1.54
HG-2-3-7-1	1.90	2.09	7.73	25.59	1.51
HG-2-3-7-2	1.90	2.09	7.89	25.56	1.51
HG-2-3-AVG	1.90	2.09	7.81	25.58	1.51
HG-3-1-7-1	1.91	2.03	7.49	24.97	1.67
HG-3-1-7-2	1.92	2.03	7.51	24.96	1.67
HG-3-1-AVG	1.92	2.03	7.50	24.96	1.68
HG-3-2-7-1	1.95	1.99	7.36	24.86	1.89
HG-3-2-7-2	1.96	2.01	7.27	24.89	1.83
HG-3-2-AVG	1.95	2.00	7.32	24.87	1.86
HG-3-3-7-1	2.01	2.02	7.18	24.96	1.96
HG-3-3-7-2	2.00	2.02	7.23	25.03	1.94
HG-3-3-AVG	2.00	2.02	7.20	24.99	1.95

PCT LEACHATE IDS	ADS #	RESEARCHER	NOTEBOOK/REF #	PCT VERSION	MEASURED pH	MEASURED B (ICP) (ppm)	NORMALIZED B (ICP) (g/L)	LOG NORM. B (ICP) (g/L)
HG-1-1-7-1	200052676	B.J. Waters	DPSTN-4789	3.0	10.41	14.37	0.66	-0.18
HG-1-1-7-2	200052677	B.J. Waters	DPSTN-4789	3.0	10.43	14.99	0.69	-0.16
HG-1-1-7-3	200052678	B.J. Waters	DPSTN-4789	3.0	10.45	15.32	0.70	-0.15
HG-1-1-AVG					10.43	14.89	0.68	-0.17
HG-1-2-7-1	200052679	B.J. Waters	DPSTN-4789	3.0	10.40	15.65	0.72	-0.14
HG-1-2-7-2	200052680	B.J. Waters	DPSTN-4789	3.0	10.39	15.71	0.73	-0.14
HG-1-2-7-3	200052681	B.J. Waters	DPSTN-4789	3.0	10.38	15.91	0.73	-0.13
HG-1-2-AVG					10.39	15.76	0.73	-0.14
HG-1-3-7-1	200052684	B.J. Waters	DPSTN-4789	3.0	10.39	14.81	0.68	-0.17
HG-1-3-7-2	200052685	B.J. Waters	DPSTN-4789	3.0	10.39	14.74	0.68	-0.17
HG-1-3-7-3	200052686	B.J. Waters	DPSTN-4789	3.0	10.39	14.93	0.69	-0.16
HG-1-3-AVG					10.39	14.83	0.68	-0.17
HG-2-1-7-1	200052687	B.J. Waters	WSRC-NB-90-271	3.0	10.44	14.15	0.66	-0.18
HG-2-1-7-2	200052688	B.J. Waters	WSRC-NB-90-271	3.0	10.43	14.75	0.69	-0.16
HG-2-1-7-3	200052689	B.J. Waters	WSRC-NB-90-271	3.0	10.44	15.20	0.71	-0.15
HG-2-1-AVG					10.44	14.70	0.69	-0.16
HG-2-2-7-1	200052692	B.J. Waters	WSRC-NB-90-271	3.0	10.70	14.85	0.77	-0.11
HG-2-2-7-2	200052693	B.J. Waters	WSRC-NB-90-271	3.0	10.74	14.87	0.77	-0.11
HG-2-2-7-3	200052694	B.J. Waters	WSRC-NB-90-271	3.0	10.74	14.86	0.77	-0.11
HG-2-2-AVG					10.73	14.86	0.77	-0.11
HG-2-3-7-1	200052695	B.J. Waters	WSRC-NB-90-271	3.0	10.71	15.02	0.79	-0.10
HG-2-3-7-2	200052696	B.J. Waters	WSRC-NB-90-271	3.0	10.72	14.94	0.79	-0.10
HG-2-3-7-3	200052697	B.J. Waters	WSRC-NB-90-271	3.0	10.69	14.96	0.79	-0.10
HG-2-3-AVG					10.71	14.97	0.79	-0.10
HG-3-1-7-1	200052700	B.J. Waters	WSRC-NB-90-271	3.0	10.72	14.08	0.74	-0.13
HG-3-1-7-2	200052701	B.J. Waters	WSRC-NB-90-271	3.0	10.73	14.36	0.75	-0.13
HG-3-1-7-3	200052702	B.J. Waters	WSRC-NB-90-271	3.0	10.74	14.86	0.78	-0.11
HG-3-1-AVG					10.73	14.43	0.75	-0.12
HG-3-2-7-1	200052703	B.J. Waters	WSRC-NB-90-271	3.0	10.70	15.36	0.79	-0.10
HG-3-2-7-2	200052704	B.J. Waters	WSRC-NB-90-271	3.0	10.70	15.04	0.77	-0.11
HG-3-2-7-3	200052705	B.J. Waters	WSRC-NB-90-271	3.0	10.73	15.19	0.78	-0.11
HG-3-2-AVG					10.71	15.20	0.78	-0.11
HG-3-3-7-1	200052709	B.J. Waters	WSRC-NB-90-271	3.0	10.68	14.24	0.71	-0.15
HG-3-3-7-2	200052710	B.J. Waters	WSRC-NB-90-271	3.0	10.68	14.60	0.73	-0.14
HG-3-3-7-3	200052712	B.J. Waters	WSRC-NB-90-271	3.0	10.69	15.09	0.75	-0.12
HG-3-3-AVG					10.68	14.64	0.73	-0.14

PCT LEACHATE IDS	MEASURED LI (ICP) (ppm)	NORMALIZED LI (ICP) (g/L)	LOG NORM. LI (ICP) (g/L)	MEASURED K (AA) (ppm)	NORMALIZED K (AA) (g/L)	LOG NORM. K (AA) (g/L)
HG-1-1-7-1	13.14	0.64	-0.19	8.80	0.40	-0.40
HG-1-1-7-2	13.64	0.66	-0.18	9.43	0.42	-0.37
HG-1-1-7-3	13.89	0.68	-0.17	9.65	0.43	-0.36
HG-1-1-AVG	13.56	0.66	-0.18	9.29	0.42	-0.38
HG-1-2-7-1	14.44	0.67	-0.17	9.29	0.36	-0.44
HG-1-2-7-2	14.46	0.67	-0.17	9.25	0.36	-0.44
HG-1-2-7-3	14.63	0.68	-0.17	8.95	0.35	-0.45
HG-1-2-AVG	14.51	0.67	-0.17	9.16	0.36	-0.44
HG-1-3-7-1	14.00	0.62	-0.21	8.94	0.41	-0.38
HG-1-3-7-2	13.90	0.62	-0.21	8.90	0.41	-0.38
HG-1-3-7-3	14.07	0.62	-0.20	8.95	0.41	-0.38
HG-1-3-AVG	13.99	0.62	-0.21	8.93	0.41	-0.38
HG-2-1-7-1	13.39	0.62	-0.21	8.08	0.40	-0.40
HG-2-1-7-2	13.88	0.65	-0.19	8.30	0.41	-0.39
HG-2-1-7-3	14.28	0.66	-0.18	8.72	0.43	-0.37
HG-2-1-AVG	13.85	0.64	-0.19	8.37	0.41	-0.39
HG-2-2-7-1	15.20	0.73	-0.14	7.03	0.46	-0.34
HG-2-2-7-2	15.25	0.73	-0.14	7.01	0.46	-0.34
HG-2-2-7-3	15.21	0.73	-0.14	6.90	0.45	-0.35
HG-2-2-AVG	15.22	0.73	-0.14	6.98	0.45	-0.34
HG-2-3-7-1	15.12	0.72	-0.14	7.38	0.49	-0.31
HG-2-3-7-2	14.92	0.71	-0.15	6.41	0.42	-0.37
HG-2-3-7-3	14.90	0.71	-0.15	6.65	0.44	-0.36
HG-2-3-AVG	14.98	0.72	-0.14	6.81	0.45	-0.35
HG-3-1-7-1	14.22	0.70	-0.15	7.38	0.44	-0.36
HG-3-1-7-2	14.51	0.71	-0.15	7.71	0.46	-0.34
HG-3-1-7-3	14.90	0.73	-0.13	8.16	0.49	-0.31
HG-3-1-AVG	14.54	0.72	-0.14	7.75	0.46	-0.34
HG-3-2-7-1	14.57	0.73	-0.14	10.16	0.55	-0.26
HG-3-2-7-2	14.46	0.72	-0.14	6.24	0.44	-0.35
HG-3-2-7-3	14.63	0.73	-0.14	8.28	0.45	-0.35
HG-3-2-AVG	14.55	0.73	-0.14	8.89	0.48	-0.32
HG-3-3-7-1	13.45	0.67	-0.18	8.30	0.43	-0.37
HG-3-3-7-2	13.78	0.68	-0.17	8.79	0.45	-0.35
HG-3-3-7-3	14.12	0.70	-0.16	8.98	0.46	-0.34
HG-3-3-AVG	13.78	0.68	-0.17	8.69	0.45	-0.35

PCT LEACHATE IDS	MEASURED Na (ICP) (ppm)	NORMALIZED Na (ICP) (g/L)	LOG NORM. Na (ICP) (g/L)	MEASURED Na (AA) (ppm)	NORMALIZED Na (AA) (g/L)	LOG NORM. Na (AA) (g/L)	MEASURED SI (ICP) (ppm)	NORMALIZED SI (ICP) (g/L)	LOG NORM. SI (ICP) (g/L)
HG-1-1-7-1	43.81	0.64	-0.19	45.70	0.67	-0.17	91.22	0.37	-0.43
HG-1-1-7-2	45.60	0.67	-0.17	47.68	0.70	-0.16	94.34	0.39	-0.41
HG-1-1-7-3	46.76	0.69	-0.16	49.18	0.72	-0.14	95.70	0.39	-0.41
HG-1-1-AVG	45.39	0.67	-0.18	47.52	0.70	-0.16	93.75	0.38	-0.42
HG-1-2-7-1	45.29	0.69	-0.16	46.36	0.71	-0.15	100.06	0.39	-0.41
HG-1-2-7-2	45.11	0.69	-0.16	47.85	0.73	-0.13	99.67	0.39	-0.41
HG-1-2-7-3	45.51	0.70	-0.16	48.02	0.74	-0.13	100.58	0.39	-0.41
HG-1-2-AVG	45.30	0.69	-0.16	47.41	0.73	-0.14	100.10	0.39	-0.41
HG-1-3-7-1	42.99	0.65	-0.18	44.04	0.67	-0.17	96.23	0.37	-0.43
HG-1-3-7-2	42.62	0.65	-0.19	44.70	0.68	-0.17	94.12	0.36	-0.44
HG-1-3-7-3	43.17	0.66	-0.18	44.54	0.68	-0.17	95.19	0.37	-0.43
HG-1-3-AVG	42.93	0.65	-0.19	44.43	0.68	-0.17	95.18	0.37	-0.43
HG-2-1-7-1	42.61	0.65	-0.19	44.37	0.68	-0.17	93.23	0.36	-0.44
HG-2-1-7-2	44.45	0.68	-0.17	46.53	0.71	-0.15	96.86	0.37	-0.43
HG-2-1-7-3	45.76	0.70	-0.16	46.69	0.71	-0.15	99.86	0.39	-0.41
HG-2-1-AVG	44.27	0.67	-0.17	45.86	0.70	-0.16	96.65	0.37	-0.43
HG-2-2-7-1	62.82	0.80	-0.10	64.90	0.83	-0.08	105.23	0.41	-0.39
HG-2-2-7-2	62.84	0.80	-0.10	65.07	0.83	-0.08	105.36	0.41	-0.39
HG-2-2-7-3	62.57	0.80	-0.10	63.41	0.81	-0.09	104.75	0.41	-0.39
HG-2-2-AVG	62.74	0.80	-0.10	64.46	0.82	-0.09	105.11	0.41	-0.39
HG-2-3-7-1	61.76	0.79	-0.10	62.59	0.80	-0.10	105.40	0.41	-0.38
HG-2-3-7-2	61.31	0.79	-0.11	62.09	0.80	-0.10	106.75	0.42	-0.38
HG-2-3-7-3	61.29	0.78	-0.11	61.76	0.79	-0.10	106.34	0.42	-0.38
HG-2-3-AVG	61.45	0.79	-0.10	62.15	0.80	-0.10	106.16	0.42	-0.38
HG-3-1-7-1	56.49	0.75	-0.12	57.95	0.77	-0.11	98.54	0.39	-0.40
HG-3-1-7-2	57.79	0.77	-0.11	60.10	0.80	-0.10	100.20	0.40	-0.40
HG-3-1-7-3	59.48	0.79	-0.10	60.76	0.81	-0.09	102.09	0.41	-0.39
HG-3-1-AVG	57.92	0.77	-0.11	59.60	0.79	-0.10	100.28	0.40	-0.40
HG-3-2-7-1	56.73	0.78	-0.11	58.45	0.80	-0.10	100.08	0.40	-0.40
HG-3-2-7-2	56.18	0.77	-0.11	56.79	0.78	-0.11	99.61	0.40	-0.40
HG-3-2-7-3	56.78	0.78	-0.11	58.12	0.79	-0.10	99.89	0.40	-0.40
HG-3-2-AVG	56.56	0.77	-0.11	57.79	0.79	-0.10	99.86	0.40	-0.40
HG-3-3-7-1	51.57	0.72	-0.15	53.31	0.74	-0.13	92.48	0.37	-0.43
HG-3-3-7-2	53.00	0.74	-0.13	54.97	0.76	-0.12	95.15	0.38	-0.42
HG-3-3-7-3	54.66	0.76	-0.12	54.14	0.75	-0.12	97.49	0.39	-0.41
HG-3-3-AVG	53.08	0.74	-0.13	54.14	0.75	-0.12	95.04	0.38	-0.42

**This page intentionally left blank.**

**END**

**DATE  
FILMED**

**3 / 16 / 93**

