Data Package

415928

Contract Lab Program Data Package
for
Du Pont Environmental Remediation Services

INORGANICS DATA PACKAGE

Christiana River Tidal Study
Grab Water Samples
Collected on 12/17/92, 12/18/92 by SN/DB
LLI Sample No. 1908350-1908359, 1908981-1908993

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SAMPLE DATA

TABLE OF CONTENTS

CASE NARRATIVE

CASE NARRATIVE FOR INORGANICS

Laboratory Name: Lancaster Laboratories

SDG Number: CRT02

Date Received: 12/19/92

Explanatory Notes:

The inappropriate error flag, % Recovery Failed, Repeat by Standard Additions, is due to a manufacturer's (Varian) software error. The software is designed to follow the analysis protocol for USEPA Contract Laboratory Program ILM01.0. Following this protocol, method of standard additions is not required for samples demonstrating sample absorbance or concentration less than 50% of the spike absorbance or concentration; however, the Varian software prints this flag, regardless of sample absorbance or concentration, whenever the percent recovery is below 85% or above 115%.

Varian has been informed of this problem and has communicated that the problem will be resolved in a future software revision. The software is needed to operate the instrument and to control the automatic spiking function of the autosampler. Since the software will not permit this error to be disabled, data that demonstrates this error will be manually corrected on the raw data by the analyst.

Calibration Standards:

Instrument calibration standards are prepared monthly from stock solutions purchased from Spex or Baker chemical.

Total ICP integration time is 24 seconds.

Case Narrative reviewed and approved by:

Eric W Cuba Date 1/8/93

Eric W. Cuba, Chemist II Coordinator
Inorganic Data Packages

2

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COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

SOW No.: 3/90_

EPA Sample No.	Lab Sample ID
_CHRDF_____	__1908989_____
_CHRDM_____	__1908988_____
_CHULF_____	__1908359_____
_CHULT_____	__1908358_____
_CRCH-_____	__1908983_____
_CRCHF_____	__1908984_____
_CRHLF_____	__1908993_____
_CRHLT_____	__1908992_____
_CRTFB_____	__1908985_____
_JAMFM_____	__1908354_____
_JAMLF_____	__1908357_____
_JAMLT_____	__1908356_____
_JAMMF_____	__1908351_____
_JAMMT_____	__1908350_____
_JAMMTD_____	__1908353_____
_JAMMTS_____	__1908352_____
_JBMTF_____	__1908987_____
_JSBFL_____	__1908982_____
_JSBHT_____	__1908981_____
_JSBLF_____	__1908991_____

Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES
If yes - were raw data generated before application of background corrections ? Yes/No NO_

Comments:

I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on floppy diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Eric W Cuba Name: Eric W. Cuba

Date: 1/8/93 Title: Chemist II Coordinator Inorganic Data Packages

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COVER PAGE - INORGANIC ANALYSES DATA PACKAGE

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

SOW No.: 3/90_

EPA Sample No.	Lab Sample ID
_JSBLT_____	__1908990_____
_JSBMT_____	__1908986_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____
_____	_____

Were ICP interelement corrections applied ? Yes/No YES

Were ICP background corrections applied ? Yes/No YES

If yes - were raw data generated before application of background corrections ? Yes/No NO_

Comments:

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I certify that this data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hardcopy data package and in the computer-readable data submitted on floppy diskette has been authorized by the Laboratory Manager or the Manager's designee, as verified by the following signature.

Signature: Eric W. Cuba Name: Eric W Cuba

Date: 1/8/93 Title: Chemist II Coordinator

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2

Sample Reference List for SDG # CRT02
with a Package Type of CLPI

Lab Sample Number	Sample Code	Client Sample Description
1908350	JAMMT	St. James St. Bridge - Mid Tide Unspiked Water
1908351	JAMMF	St. James St. Bridge - Mid Tide Unspiked Filtered
1908352	JAMMT	St. James St. Bridge - Mid Tide Matrix Spike Water
1908353	JAMMT	St. James St. Bridge - Mid Tide Duplicate Water
1908354	JAMFM	St. James St. Bridge - Mid Tide Duplicate Filtered
1908355	FLDBL	Field Blank Grab Water Sample
1908356	JAMLT	St. James St. Bridge - Low Tide Grab Water Sample
1908357	JAMLF	St. James St. Bridge - Low Tide Filtered Water
1908358	CHULT	Churchman's Road - Low Tide Grab Water Sample
1908359	CHULF	Churchman's Road - Low Tide Filtered Grab Water
1908981	JSBHT	James St. Bridge - High Tide Grab Water Sample
1908982	JSBFL	James St. Bridge - High Tide Filtered Water Sample
1908983	CRCH-	Churchman's Road - High Tide Grab Water Sample
1908984	CRCHF	Churchman's Road - High Tide Filtered Water Sample
1908985	CRTFB	Field Blank Grab Water Sample
1908986	JSBMT	James St. Bridge - Mid Tide Grab Water Sample
1908987	JBMTF	James St. Bridge - Mid Tide Filtered Water Sample
1908988	CHRDM	Churchman's Road - Mid Tide Grab Water Sample
1908989	CHRDF	Churchman's Road - Mid Tide Filtered Water Sample
1908990	JSBLT	James St. Bridge - Low Tide Grab Water Sample
1908991	JSBLF	James St. Bridge - Low Tide Filtered Water Sample
1908992	CRHLT	Churchman's Road - Low Tide Grab Water Sample
1908993	CRHLF	Churchman's Road - Low Tide Filtered Water Sample

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1
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EPA SAMPLE NO.

JAMMT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908350___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	1990			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	68.2	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	47600			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	2220		N	P
7439-92-1	Lead	2.9	B		F
7439-95-4	Magnesium	7130			P
7439-96-5	Manganese	113			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5920			P
7782-49-2	Selenium	1.0	U		F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	10400			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	74.9			P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JAMMF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908351___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: _____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum_	57.0	U		P	
7440-36-0	Antimony_	22.0	U		P	
7440-38-2	Arsenic__	1.0	U		F	
7440-39-3	Barium___	[illegible]	B		P	
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium__	3.0	U		P	
7440-70-2	Calcium__	[illegible]			P	
7440-47-3	Chromium_	8.0	U		P	
7440-48-4	Cobalt___	5.0	U		P	
7440-50-8	Copper___	5.0	U		P	
7439-89-6	Iron_____	[illegible]		N	P	
7439-92-1	Lead_____	1.0	U		F	
7439-95-4	Magnesium	[illegible]			P	
7439-96-5	Manganese	[illegible]			P	
7439-97-6	Mercury__	0.10	U		CV	
7440-02-0	Nickel___	8.0	U		P	
7440-09-7	Potassium	[illegible]			P	
7782-49-2	Selenium_	1.0	U	W	F	
7440-22-4	Silver___	11.0	U		P	
7440-23-5	Sodium___	[illegible]			P	
7440-28-0	Thallium_	2.0	U		F	
7440-62-2	Vanadium_	6.0	U		P	
7440-66-6	Zinc_____	[illegible]			P	U
_____	Cyanide__	_____	_	_____	NR	

Color Before: COLORLESS Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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EPA SAMPLE NO.

JAMFM

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908354___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum_	57.0	U		P
7440-36-0	Antimony_	22.0	U		P
7440-38-2	Arsenic__	1.0	U		F
7440-39-3	Barium___	49.6	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium__	3.0	U		P
7440-70-2	Calcium__	17000			P
7440-47-3	Chromium_	8.0	U		P
7440-48-4	Cobalt___	5.0	U		P
7440-50-8	Copper___	5.0	U		P
7439-89-6	Iron_____	153		N	P
7439-92-1	Lead_____	1.0	U		F
7439-95-4	Magnesium	6750			P
7439-96-5	Manganese	80.5			P
7439-97-6	Mercury__	0.10	U		CV
7440-02-0	Nickel___	8.0	U		P
7440-09-7	Potassium	5590			P
7782-49-2	Selenium_	1.0	U	W	F
7440-22-4	Silver___	11.0	U		P
7440-23-5	Sodium___	10200			P
7440-28-0	Thallium_	2.0	U		F
7440-62-2	Vanadium_	6.0	U		P
7440-66-6	Zinc_____	44.0			P
_____	Cyanide__	_____	_	_____	NR

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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1

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JAMLT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908356___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	1480			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	60.1	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	17800			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	1830		N	P
7439-92-1	Lead	3.2	B		F
7439-95-4	Magnesium	7090			P
7439-96-5	Manganese	109			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5400			P
7782-49-2	Selenium	1.0	U		F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9660			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	79.3			P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

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EPA SAMPLE NO.

JAMLF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908357___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	57.0	U		P	
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	R
7440-39-3	Barium	40.8	B		P	
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	17100			P	
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.0	U		P	
7439-89-6	Iron	165		N	P	
7439-92-1	Lead	1.0	U	W	F	
7439-95-4	Magnesium	6750			P	
7439-96-5	Manganese	83.3			P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	5130			P	
7782-49-2	Selenium	1.0	U	W	F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	9450			P	
7440-28-0	Thallium	2.0	U		F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	40.9			P	V
	Cyanide				NR	

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

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1

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CHULT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908358___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	1680			P	
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	R
7440-39-3	Barium	52.9	B		P	
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	10300			P	
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.0	U		P	
7439-89-6	Iron	2480		N	P	
7439-92-1	Lead	3.0			F	
7439-95-4	Magnesium	4500	B		P	
7439-96-5	Manganese	154			P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	3020	B		P	
7782-49-2	Selenium	1.0	U		F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	9110			P	
7440-28-0	Thallium	2.0	U		F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	33.5			P	V
_____	Cyanide	_____			NR	
_____	_____	_____				

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

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EPA SAMPLE NO.

CHULF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908359___

Level (low/med): LOW__ Date Received: 12/18/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	57.0	U		P	
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	R
7440-39-3	Barium	41.2	B		P	
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	9210			P	
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.0	U		P	
7439-89-6	Iron	325		N	P	
7439-92-1	Lead	1.0	U		F	
7439-95-4	Magnesium	3950	B		P	
7439-96-5	Manganese	116			P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	2720	B		P	
7782-49-2	Selenium	1.0	U		F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	8160			P	
7440-28-0	Thallium	2.0	U		F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	16.0	B		P	V
_____	Cyanide	_____	_	_____	NR	
_____	_____	_____	_	_____	__	

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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1

INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JSBHT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908981___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: _____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	2100			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	60.9	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	17800			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	2260		N	P
7439-92-1	Lead	3.9		S	F
7439-95-4	Magnesium	6980			P
7439-96-5	Manganese	102			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5270			P
7782-49-2	Selenium	1.0	U	W	F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9710			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	78.0			P
_____	Cyanide	_____	_	_____	NR

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JSBFL

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908982___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	57.0	U		P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	40.8	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	17400			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	138		N	P
7439-92-1	Lead	1.0	U	W	F
7439-95-4	Magnesium	6680			P
7439-96-5	Manganese	67.8			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5150			P
7782-49-2	Selenium	1.0	U	W	F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9640			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	34.7			P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CRCH-

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908983___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	1480			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	46.5	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	10200			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	6.0	B		P
7439-89-6	Iron	2000		N	P
7439-92-1	Lead	4.0			F
7439-95-4	Magnesium	4340	B		P
7439-96-5	Manganese	115			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	2920	B		P
7782-49-2	Selenium	1.0	U		F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	10600			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	30.1			P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

ILMO2.1

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CRCHF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908984___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	57.0	U		P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	41.6	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	10200			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	314		N	P
7439-92-1	Lead	1.0	U		F
7439-95-4	Magnesium	4290	B		P
7439-96-5	Manganese	105			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	2720	B		P
7782-49-2	Selenium	1.0	U	W	F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	10500			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	16.3	B		P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO. 02

CRTFB

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908985___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	57.0	U		P	
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	R
7440-39-3	Barium	4.0	U		P	
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	171	B		P	V
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.0	U		P	
7439-89-6	Iron	7.0	U	N	P	
7439-92-1	Lead	1.0	U		F	
7439-95-4	Magnesium	45.0	U		P	
7439-96-5	Manganese	1.0	U		P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	205	B		P	
7782-49-2	Selenium	1.0	U		F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	417	B		P	U
7440-28-0	Thallium	2.0	U		F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	4.0	U		P	
	Cyanide				NR	

Color Before: COLORLESS Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JSBMT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908986___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	1500			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	60.9	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	17800			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	2610		N	P
7439-92-1	Lead	5.0		S	F
7439-95-4	Magnesium	6980			P
7439-96-5	Manganese	126			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	6050			P
7782-49-2	Selenium	1.0	U	W	F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9240			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	89.8			P
_____	Cyanide	_____			NR

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JBMTF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908987___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	57.0	U		P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	38.4	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	17000			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	170		N	P
7439-92-1	Lead	1.0	U		F
7439-95-4	Magnesium	6460			P
7439-96-5	Manganese	70.3			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5770			P
7782-49-2	Selenium	1.0	U		F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9020			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	33.4			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CHRDM

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908988___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	2330			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	50.5	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	10400			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	2740		N	P
7439-92-1	Lead	7.6			F
7439-95-4	Magnesium	4360	B		P
7439-96-5	Manganese	117			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	2850	B		P
7782-49-2	Selenium	1.0	U	W	F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9720			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	29.3			P
	Cyanide				NR

Color Before: COLORLESS Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

FORM I - IN

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CHRDF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908989___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	57.0	U		P	
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	R
7440-39-3	Barium	[illegible]	B		P	V
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	[illegible]			P	
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.0	U		P	
7439-89-6	Iron	[illegible]		N	P	
7439-92-1	Lead	1.0	U		F	
7439-95-4	Magnesium	4310	B		P	
7439-96-5	Manganese	[illegible]			P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	2850	B		P	
7782-49-2	Selenium	1.0	U		F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	10200			P	
7440-28-0	Thallium	2.0	U		F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	16.0	B		P	V
	Cyanide				NR	

Color Before: COLORLESS Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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U.S. EPA - CLP

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JSBLT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908990___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	9290			P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	60.1	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	15900			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	12.4	B		P
7439-89-6	Iron	2840		N	P
7439-92-1	Lead	5.4			F
7439-95-4	Magnesium	6340			P
7439-96-5	Manganese	135			P
7439-97-6	Mercury	0.10	B		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5450			P
7782-49-2	Selenium	1.0	U		F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9510			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	81.6			P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

FORM I - IN

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24

U.S. EPA - CLP

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

JSBLF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908991___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M
7429-90-5	Aluminum	57.0	U		P
7440-36-0	Antimony	22.0	U		P
7440-38-2	Arsenic	1.0	U		F
7440-39-3	Barium	38.4	B		P
7440-41-7	Beryllium	1.0	U		P
7440-43-9	Cadmium	3.0	U		P
7440-70-2	Calcium	15600			P
7440-47-3	Chromium	8.0	U		P
7440-48-4	Cobalt	5.0	U		P
7440-50-8	Copper	5.0	U		P
7439-89-6	Iron	185		N	P
7439-92-1	Lead	1.0	U		F
7439-95-4	Magnesium	6050			P
7439-96-5	Manganese	91.5			P
7439-97-6	Mercury	0.10	U		CV
7440-02-0	Nickel	8.0	U		P
7440-09-7	Potassium	5180			P
7782-49-2	Selenium	1.0	U		F
7440-22-4	Silver	11.0	U		P
7440-23-5	Sodium	9640			P
7440-28-0	Thallium	2.0	U		F
7440-62-2	Vanadium	6.0	U		P
7440-66-6	Zinc	27.8			P
	Cyanide				NR

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CRHLT

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908992___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	65.1	B		P	U
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	R
7440-39-3	Barium	24.4	B		P	U
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	8390			P	
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.0	U		P	
7439-89-6	Iron	250		N	P	
7439-92-1	Lead	7.4			F	
7439-95-4	Magnesium	3320	B		P	
7439-96-5	Manganese	73.6			P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	2560	B		P	
7782-49-2	Selenium	1.0	U		F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	6850			P	
7440-28-0	Thallium	2.0	U	W	F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	11.1	B		P	U
	Cyanide				NR	

Color Before: YELLOW___ Clarity Before: CLOUDY Texture: _____

Color After: COLORLESS Clarity After: CLEAR_ Artifacts: _____

Comments:

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FORM I - IN

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13

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INORGANIC ANALYSES DATA SHEET

EPA SAMPLE NO.

CRHLF

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Lab Sample ID: 1908993___

Level (low/med): LOW__ Date Received: 12/19/92

% Solids: ____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

CAS No.	Analyte	Concentration	C	Q	M	
7429-90-5	Aluminum	3000			P	
7440-36-0	Antimony	22.0	U		P	
7440-38-2	Arsenic	1.0	U		F	
7440-39-3	Barium	45.7	B		P	V
7440-41-7	Beryllium	1.0	U		P	
7440-43-9	Cadmium	3.0	U		P	
7440-70-2	Calcium	8650			P	
7440-47-3	Chromium	8.0	U		P	
7440-48-4	Cobalt	5.0	U		P	
7440-50-8	Copper	5.3	B		P	V
7439-89-6	Iron	3240		N	P	
7439-92-1	Lead	1.0	U		F	
7439-95-4	Magnesium	3580	B		P	
7439-96-5	Manganese	107			P	
7439-97-6	Mercury	0.10	U		CV	
7440-02-0	Nickel	8.0	U		P	
7440-09-7	Potassium	2920	B		P	
7782-49-2	Selenium	1.0	U		F	
7440-22-4	Silver	11.0	U		P	
7440-23-5	Sodium	6780			P	
7440-28-0	Thallium	2.0	U		F	
7440-62-2	Vanadium	6.0	U		P	
7440-66-6	Zinc	33.0			P	V
	Cyanide				NR	

Color Before: YELLOW___ Clarity Before: CLEAR_ Texture: _____

Color After: COLORLESS Clarity After: CLOUDY Artifacts: _____

Comments:

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QUALITY CONTROL DATA

Analysis Request/Environmental Services

Lancaster Laboratories
Where quality is a science.

M08981-93

Chain of Custody

Please print. Instructions on reverse side correspond with circled numbers.

(1) Client: DuPont — Acct. #:
Project Name/#: Christina River Tidal Study
Project Manager: Joel Kalmazyn — P.O. #:
Sampler: Scott Northey / Dan Bacon — Quote #:

For LLI use only
FSC:
SCR #: 104295

AR318844

(2) Sample Identification	Date Collected	Time Collected	(3) Grab	Composite	(4) Matrix: Soil	Water	Other	Total # of Containers	(5) Analyses Requested: Metals - Total	Metals - Dissolved	(6) Remarks
James St. Bridge - High Tide	12/18/92	720	✓			✓		2	✓	✓	
Churchman's Road - High Tide	↓	800	✓			✓		2	✓	✓	
Field Blank	↓	1025	✓			✓		1	✓		
James St. Bridge - Mid Tide	↓	1030	✓			✓		2	✓	✓	

Cooler Temp. 4°C
12/19/92

(7) **Turnaround time requested** (please circle): Normal Rush
(Rush TAT is subject to LLI approval and surcharge.)
Rush results requested by (please circle):
Fax Fax #:
Phone Phone #:

(8) **Data package options** (please circle if requested):
QC Summary
Tier I (NJ)
Tier II (NJ)
EPA CLP
Site-specific QC required? Yes No
(If yes, indicate QC sample and submit triplicate volume.)
Data Package Chain of Custody required? Yes No

(9)

Relinquished by:	Date	Time	Received by:	Date	Time
Tina M. Wike	12/18/92	16:39	Signature	12/17/92	1200
Signature	12/18/92	1600			
Relinquished by:	Date	Time	Received by:	Date	Time
Relinquished by:	Date	Time	Received by:	Date	Time
Relinquished by:	Date	Time	Received for LLI by:	Date	Time
			Jeanne M. Zuckerstein	12/19/92	0900

Analysis Request/Environmental Services

Lancaster Laboratories
Where quality is a science.

3656
1908350-59
1908335-40

Chain of Custody

Please print. Instructions on reverse side correspond with circled numbers.

(1) Client: DuPont — Acct. #:
Project Name/#: Christina River Tidal Study
Project Manager: Joel Karmazyn — P.O. #:
Sampler: S. Northey / D. Bacon — Quote #:

For LLI use only
FSC:
SCR #: 104396

(2) Sample Identification	Date Collected	Time Collected	(3) Grab	Composite	(4) Matrix: Soil	Water	Other	Total # of Containers	(5) Analyses Requested: Metals Total	Metals - Dissolved	(6) Remarks
St. James St. Bridge - High Tide	12/11/92	610	✓			✓		2	✓	✓	
Churchman's Road - High Tide		645	✓			✓		2	✓	✓	
St. James St. Bridge - Mid Tide		930	✓			✓		2	✓	✓	
St. James St Bridge - Duplicate Mid Tide		930	✓			✓		2	✓	✓	
St. James St. Bridge - Matrix Spike Mid Tide		930	✓			✓		1	✓		
Churchman's Road - Mid Tide		1000	✓			✓		2	✓	✓	
Field Blank		955	✓			✓		1	✓		
St. James St. Bridge - Low Tide		1300	✓			✓		2	✓	✓	
Churchmans Road - Low Tide		1340	✓			✓		2	✓	✓	Temp. of cooler upon receipt: 5°C. 12/15/92 DN

(7) **Turnaround time requested** (please circle): Normal Rush
(Rush TAT is subject to LLI approval and surcharge.)

Rush results requested by (please circle):
Fax Fax #:
Phone Phone #:

(8) **Data package options** (please circle if requested):
QC Summary
Tier I (NJ)
Tier II (NJ)
EPA CLP
Site-specific QC required? Yes No
(If yes, indicate QC sample and submit triplicate volume.)
Data Package Chain of Custody required? Yes No

Relinquished by:	Date	Time	Received by:	Date	Time
Gina M. Wilt	12-14-92	22:45	S. Northey	12/15/92	1200
S. Northey	12/15/92	1500			
Relinquished by:	Date	Time	Received by:	Date	Time
Relinquished by:	Date	Time	Received by:	Date	Time
Relinquished by:	Date	Time	Received by:	Date	Time
Relinquished by:	Date	Time	Received for LLI by:	Date	Time
			Deborah A. Nesluna	12/15/92	0855

(9)

2425 New Holland Pike, Lancaster, PA 17601-5994 (717) 656-2301

Copies: White and Yellow copies accompany samples to lab. Pink is retained by client.

U.S. EPA - CLP

2A
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Initial Calibration Source: LLI_____

Continuing Calibration Source: LLI_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Aluminum	25000.0	24397.00	97.6	25000.0	24141.00	96.6	23754.00	95.0	P
Antimony	5000.0	5175.00	103.5	5000.0	5114.30	102.3	5009.70	100.2	P
Arsenic	50.0	50.79	101.6	50.0	49.59	99.2	50.21	100.4	F
Barium	5000.0	5079.80	101.6	5000.0	5036.60	100.7	4933.00	98.7	P
Beryllium	500.0	489.35	97.9	500.0	485.24	97.0	475.69	95.1	P
Cadmium	5000.0	4758.82	95.2	5000.0	4793.90	95.9	4802.60	96.1	P
Calcium	50000.0	50224.00	100.4	50000.0	49634.00	99.3	48655.00	97.3	P
Chromium	5000.0	5231.00	104.6	5000.0	5193.80	103.9	5095.40	101.9	P
Cobalt	1000.0	1001.10	100.1	1000.0	989.82	99.0	966.19	96.6	P
Copper	5000.0	4951.30	99.0	5000.0	4911.00	98.2	4815.20	96.3	P
Iron	5000.0	5036.70	100.7	5000.0	4982.30	99.6	4883.20	97.7	P
Lead	20.0	20.73	103.6	20.0	20.45	102.2	21.03	105.2	F
Magnesium	25000.0	24165.00	96.7	25000.0	24000.00	96.0	23665.00	94.7	P
Manganese	5000.0	5029.10	100.6	5000.0	4996.50	99.9	4905.60	98.1	P
Mercury	2.5	2.53	101.2	1.0	1.02	102.0	1.02	102.0	CV
Nickel	10000.0	9927.10	99.3	10000.0	9856.50	98.6	9686.50	96.9	P
Potassium	50000.0	50954.00	101.9	50000.0	50480.00	101.0	49327.00	98.7	P
Selenium	20.0	20.21	101.0	20.0	19.81	99.0	19.55	97.8	F
Silver	500.0	478.06	95.6	500.0	474.81	95.0	466.69	93.3	P
Sodium	50000.0	52505.00	105.0	50000.0	52393.00	104.8	49555.00	99.1	P
Thallium	40.0	40.70	101.8	40.0	42.60	106.5	43.70	109.2	F
Vanadium	2000.0	2078.60	103.9	2000.0	2058.50	102.9	2015.70	100.8	P
Zinc	5000.0	5026.50	100.5	5000.0	5000.70	100.0	4924.90	98.5	P
Cyanide									NR

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2A
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Initial Calibration Source: LLI_____

Continuing Calibration Source: LLI_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Initial Calibration Found	Initial Calibration %R(1)	Continuing Calibration True	Continuing Calibration Found	Continuing Calibration %R(1)	Continuing Calibration Found	Continuing Calibration %R(1)	M
Aluminum				25000.0	23828.00	95.3	23911.00	95.6	P
Antimony				5000.0	5037.10	100.7	5062.50	101.2	P
Arsenic				50.0	49.75	99.5	47.40	94.8	F
Barium				5000.0	4946.20	98.9	4955.80	99.1	P
Beryllium				500.0	477.44	95.5	478.30	95.7	P
Cadmium				5000.0	4788.60	95.8	4662.30	93.2	P
Calcium				50000.0	48783.00	97.6	48794.00	97.6	P
Chromium				5000.0	5117.80	102.4	5115.40	102.3	P
Cobalt				1000.0	970.61	97.1	967.76	96.8	P
Copper				5000.0	4833.70	96.7	4835.60	96.7	P
Iron				5000.0	4905.80	98.1	4940.20	98.8	P
Lead				20.0	21.82	109.1			F
Magnesium				25000.0	23764.00	95.1	23843.00	95.4	P
Manganese				5000.0	4916.60	98.3	4925.20	98.5	P
Mercury				1.0	1.01	101.0	1.02	102.0	CV
Nickel				10000.0	9701.10	97.0	9713.60	97.1	P
Potassium				50000.0	49006.00	98.0	49147.00	98.3	P
Selenium				20.0	19.81	99.0	20.31	101.6	F
Silver				500.0	463.76	92.8	464.41	92.9	P
Sodium				50000.0	50177.00	100.4	53757.00	107.5	P
Thallium	40.0	38.50	96.2	40.0	38.40	96.0	39.60	99.0	F
Vanadium				2000.0	2020.80	101.0	2026.60	101.3	P
Zinc				5000.0	4936.00	98.7	4941.30	98.8	P
Cyanide									NR

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2A
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Initial Calibration Source: LLI_____

Continuing Calibration Source: LLI_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Initial Calibration Found	Initial Calibration %R(1)	Continuing Calibration True	Continuing Calibration Found	Continuing Calibration %R(1)	Continuing Calibration Found	Continuing Calibration %R(1)	M
Aluminum									NR
Antimony									NR
Arsenic				50.0	46.37	92.7	45.34	90.7	F
Barium									NR
Beryllium									NR
Cadmium									NR
Calcium									NR
Chromium									NR
Cobalt									NR
Copper									NR
Iron									NR
Lead	20.0	20.69	103.4	20.0	20.16	100.8	20.76	103.8	F
Magnesium									NR
Manganese									NR
Mercury									NR
Nickel									NR
Potassium									NR
Selenium				20.0	20.95	104.8	20.94	104.7	F
Silver									NR
Sodium									NR
Thallium				40.0	42.10	105.2	43.50	108.8	F
Vanadium									NR
Zinc									NR
Cyanide									NR

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2A
INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Initial Calibration Source: LLI_____

Continuing Calibration Source: LLI_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Aluminum									NR
Antimony									NR
Arsenic				50.0	45.03	90.1			F
Barium									NR
Beryllium									NR
Cadmium									NR
Calcium									NR
Chromium									NR
Cobalt									NR
Copper									NR
Iron									NR
Lead				20.0	20.88	104.4	21.47	107.4	F
Magnesium									NR
Manganese									NR
Mercury									NR
Nickel									NR
Potassium									NR
Selenium	20.0	20.77	103.8	20.0	21.13	105.6	20.05	100.2	F
Silver									NR
Sodium									NR
Thallium	40.0	42.00	105.0	40.0	42.80	107.0	41.90	104.8	F
Vanadium									NR
Zinc									NR
Cyanide									NR

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2A

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Initial Calibration Source: LLI_____

Continuing Calibration Source: LLI_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Aluminum									NR
Antimony									NR
Arsenic									NR
Barium									NR
Beryllium									NR
Cadmium									NR
Calcium									NR
Chromium									NR
Cobalt									NR
Copper									NR
Iron									NR
Lead				20.0	21.79	109.0			F
Magnesium									NR
Manganese									NR
Mercury									NR
Nickel									NR
Potassium									NR
Selenium	20.0	21.85	109.2	20.0	21.50	107.5	20.07	100.4	F
Silver									NR
Sodium									NR
Thallium									NR
Vanadium									NR
Zinc									NR
Cyanide									NR

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2A

INITIAL AND CONTINUING CALIBRATION VERIFICATION

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Initial Calibration Source: LLI_____

Continuing Calibration Source: LLI_____

Concentration Units: ug/L

Analyte	Initial Calibration True	Found	%R(1)	Continuing Calibration True	Found	%R(1)	Found	%R(1)	M
Aluminum									NR
Antimony									NR
Arsenic									NR
Barium									NR
Beryllium									NR
Cadmium									NR
Calcium									NR
Chromium									NR
Cobalt									NR
Copper									NR
Iron									NR
Lead	20.0	21.11	105.6	20.0	20.35	101.8	20.02	100.1	F
Magnesium									NR
Manganese									NR
Mercury									NR
Nickel									NR
Potassium									NR
Selenium									NR
Silver									NR
Sodium									NR
Thallium									NR
Vanadium									NR
Zinc									NR
Cyanide									NR

(1) Control Limits: Mercury 80-120; Other Metals 90-110; Cyanide 85-115

U.S. EPA - CLP

2B
CRDL STANDARD FOR AA AND ICP

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

AA CRDL Standard Source: LLI_____

ICP CRDL Standard Source: LLI_____

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP				
					Initial		Final	
	True	Found	%R	True	Found	%R	Found	%R
Aluminum								
Antimony				120.0	124.97	104.1	110.88	92.4
Arsenic	10.0	9.31	93.1					
Barium								
Beryllium				10.0	8.93	89.3	10.04	100.4
Cadmium								
Calcium								
Chromium				20.0	20.65	103.2	22.31	111.6
Cobalt				100.0	99.53	99.5	100.75	100.8
Copper				50.0	49.99	100.0	47.48	95.[illegible]
Iron								
Lead	3.0	2.79	93.0					
Magnesium								
Manganese				30.0	29.75	99.2	28.91	96.4
Mercury	0.2	0.21	105.0					
Nickel				80.0	78.42	98.0	79.62	99.5
Potassium								
Selenium	5.0	5.28	105.6					
Silver				20.0	18.18	90.9	98.25	491.2
Sodium								
Thallium	10.0	10.80	108.0					
Vanadium				100.0	100.71	100.7	98.80	98.8
Zinc				40.0	40.08	100.2	40.64	101.6

U.S. EPA - CLP

2B
CRDL STANDARD FOR AA AND ICP

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

AA CRDL Standard Source: LLI_____

ICP CRDL Standard Source: LLI_____

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP				
	True	Found	%R	True	Initial Found	%R	Final Found	%R
Aluminum								
Antimony								
Arsenic								
Barium								
Beryllium								
Cadmium				10.0	9.87	98.7	9.05	90.5
Calcium								
Chromium								
Cobalt								
Copper								
Iron								
Lead	3.0	3.15	105.0					
Magnesium								
Manganese								
Mercury								
Nickel								
Potassium								
Selenium	5.0	5.10	102.0					
Silver								
Sodium								
Thallium	10.0	8.60	86.0					
Vanadium								
Zinc								

U.S. EPA - CLP

2B
CRDL STANDARD FOR AA AND ICP

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

AA CRDL Standard Source: LLI_____

ICP CRDL Standard Source: _____

Concentration Units: ug/L

Analyte	CRDL Standard for AA			CRDL Standard for ICP				
	True	Found	%R	True	Initial Found	%R	Final Found	%R
Aluminum								
Antimony								
Arsenic								
Barium								
Beryllium								
Cadmium								
Calcium								
Chromium								
Cobalt								
Copper								
Iron								
Lead	3.0	2.15	71.7					
Magnesium								
Manganese								
Mercury								
Nickel								
Potassium								
Selenium	5.0	5.39	107.8					
Silver								
Sodium								
Thallium	10.0	10.80	108.0					
Vanadium								
Zinc								

3
BLANKS

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Preparation Blank Matrix (soil/water): WATER

Preparation Blank Concentration Units (ug/L or mg/kg): UG/L_

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prep. Blank	C	M
Aluminum	57.0	U	57.0	U	156.6	B	57.0	U	57.000	U	P
Antimony	22.0	U	22.0	U	22.0	U	22.0	U	22.000	U	P
Arsenic	-2.5	B	1.0	U	1.0	U	1.0	U	1.000	U	F
Barium	4.0	U	4.0	U	4.0	U	4.0	B	4.000	U	P
Beryllium	1.0	U	1.0	U	1.0	U	1.0	U	1.000	U	P
Cadmium	3.0	U	3.3	B	4.5	B	3.0	U	3.000	U	P
Calcium	23.0	U	23.0	U	23.0	U	40.5	B	23.000	U	P
Chromium	8.0	U	8.0	U	8.0	U	8.0	U	8.000	U	P
Cobalt	5.0	U	5.0	U	5.0	U	5.0	U	5.000	U	P
Copper	5.0	U	5.0	U	5.0	U	5.0	U	5.000	U	P
Iron	7.0	U	7.0	U	7.0	U	7.0	U	22.650	B	P
Lead	1.0	U	1.0	U	1.0	U	1.0	U	1.000	U	F
Magnesium	45.0	U	45.0	U	45.0	U	45.0	U	45.000	U	P
Manganese	1.0	U	1.0	U	1.0	U	1.2	B	1.000	U	P
Mercury	0.1	U	0.1	U	0.1	U	0.1	U	0.100	U	CV
Nickel	8.0	U	8.0	U	8.0	U	8.0	U	8.000	U	P
Potassium	132.0	U	132.0	U	132.0	U	132.0	U	132.000	U	P
Selenium	1.0	U	1.0	U	1.0	U	1.0	U	1.000	U	F
Silver	11.0	U	11.0	U	11.0	U	11.0	U	11.000	U	P
Sodium	72.0	U	72.0	U	72.0	U	72.0	U	92.460	B	P
Thallium	2.0	U	2.0	U	2.0	U			2.000	U	F
Vanadium	6.0	U	6.0	U	6.0	U	6.0	U	6.000	U	P
Zinc	4.0	U	4.0	U	13.1	B	5.0	B	4.000	U	P
Cyanide											NR

U.S. EPA - CLP

3
BLANKS

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Aluminum			73.4	B							P
Antimony			22.0	U							P
Arsenic			-1.5	B	-1.2	B	1.0	U			F
Barium			10.4	B							P
Beryllium			1.0	U							P
Cadmium			3.7	B							P
Calcium			117.8	B							P
Chromium			8.0	U							P
Cobalt			5.0	U							P
Copper			7.9	B							P
Iron			16.2	B							P
Lead	1.0	U	1.0	U	1.0	U	1.0	U			F
Magnesium			58.6	B							P
Manganese			6.9	B							P
Mercury			0.1	U							CV
Nickel			16.9	B							P
Potassium			256.2	B							P
Selenium			1.0	U	1.0	U	1.0	U			F
Silver			11.0	U							P
Sodium			92.4	B							P
Thallium	-2.2	B	2.0	U	2.0	U	2.0	U			F
Vanadium			6.0	U							P
Zinc			9.3	B							P
Cyanide											NR

FORM III - IN

ILMO2.1

U.S. EPA - CLP

3
BLANKS

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Aluminum											NR
Antimony											NR
Arsenic			1.0	U							F
Barium											NR
Beryllium											NR
Cadmium											NR
Calcium											NR
Chromium											NR
Cobalt											NR
Copper											NR
Iron											NR
Lead			1.0	U	1.0	U					F
Magnesium											NR
Manganese											NR
Mercury											NR
Nickel											NR
Potassium											NR
Selenium	1.0	U	1.0	U	1.0	U					F
Silver											NR
Sodium											NR
Thallium			2.0	U							F
Vanadium											NR
Zinc											NR
Cyanide											NR

3
BLANKS

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Preparation Blank Matrix (soil/water): _____

Preparation Blank Concentration Units (ug/L or mg/kg): _____

Analyte	Initial Calib. Blank (ug/L)	C	Continuing Calibration Blank (ug/L) 1	C	2	C	3	C	Prepa-ration Blank	C	M
Aluminum											NR
Antimony											NR
Arsenic											NR
Barium											NR
Beryllium											NR
Cadmium											NR
Calcium											NR
Chromium											NR
Cobalt											NR
Copper											NR
Iron											NR
Lead	1.0	U	1.0	U	1.0	U					F
Magnesium											NR
Manganese											NR
Mercury											NR
Nickel											NR
Potassium											NR
Selenium	1.0	U	1.0	U	1.0	U					F
Silver											NR
Sodium											NR
Thallium	2.0	U	2.0	U	2.0	U					F
Vanadium											NR
Zinc											NR
Cyanide											NR

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2360-ICP1100_ Method: P_

Start Date: 12/19/92 End Date: 12/20/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	2215		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
S	1.00	2219				_	X		_		X		X	X	_		X	_	X		_			_		X	_
S	1.00	2223		_	X	_	_	X	_	_	_	X	_	_	_	_	_	_	_	_	_	_	_	_	X	_	_
S	1.00	2228		_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	X	_	_	_	_	_
S	1.00	2231		_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	X	_	_	X	_	_	_	_
S	1.00	2233		X	_	_	_	_	_	X	_	_	_	_	_	X	_	_	_	_	_	_	_	_	_	_	_
ICV	1.00	2236		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
ICB	1.00	2240		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
ICSA	1.00	2245		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
ICSAB	1.00	2249		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CRI	1.00	2254			X	_		X	_		X	X	X		_		X	_	X		_	X		_	X	X	_
CCV	1.00	2259		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CCB (1)	1.00	2303		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
ICSW	1.00	2308		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
LCSW	1.00	2312		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMMT	1.00	2317		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMMT A	1.00	2321												X													
JAMMT D	1.00	2326		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMMT S	1.00	2331		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMMT L	5.00	2335		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMMF	1.00	2340		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMFM	1.00	2344		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
ZZZZZZ	1.00	2349																									
CCV	1.00	2353		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CCB (2)	1.00	2358		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMLT	1.00	0002		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JAMLF	1.00	0007		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CHULT	1.00	0012		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CHULF	1.00	0016		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JSBHT	1.00	0021		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JSBFL	1.00	0025		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CRCH-	1.00	0030		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
				_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_	_

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: ___________

Lab Code: _______ Case No.: _______ SAS No.: _______ SDG No.:CRT02_

Instrument ID Number: 2360-ICP1100_ Method: P_

Start Date: 12/19/92 End Date: 12/20/92

EPA Sample No.	D/F	Time	% R	Analytes																							
				A L	S B	A S	B A	B E	C D	C A	C R	C O	C U	F E	P B	M G	M N	H G	N I	K	S E	A G	N A	T L	V	Z N	C N
CRCHF	1.00	0035		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CRTFB	1.00	0039		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JSBMT	1.00	0044		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CCV	1.00	0049		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CCB (2)	1.00	0053		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JBMTF	1.00	0058		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CHRDM	1.00	0103		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CHRDF	1.00	0107		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JSBLT	1.00	0112		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
JSBLF	1.00	0117		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CRHLT	1.00	0121		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CRHLF	1.00	0126		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
ICSA	1.00	0131		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	[illegible]	[illegible]
ICSAB	1.00	0135		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	[illegible]	[illegible]
CRI	1.00	0140		_	X	_	_	X	_	_	X	X	X	_	_	_	X	_	X	_	_	X	_	_	X	X	
CCV	1.00	0145		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_
CCB (4)	7.9 1.00	0149		X	X	_	X	X	_	X	X	X	X	X	_	X	X	_	X	X	_	X	X	_	X	X	_

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2360-ICP1100_ Method: P_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1433							X																		
S	1.00	1437							X																		
S	1.00	1441																									
S	1.00	1446																									
S	1.00	1449																									
S	1.00	1451																									
ICV	1.00	1454							X																		
ICB	1.00	1458							X																		
ICSA	1.00	1503							X																		
ICSAB	1.00	1507							X																		
CRI	1.00	1512							X																		
CCV	1.00	1517							X																		
[illegible]B	1.00	1521							X																		
[illegible]W	1.00	1526							X																		
[illegible]CSW	1.00	1530							X																		
JAMMT	1.00	1535							X																		
JAMMT A	1.00	1539																									
JAMMT D	1.00	1544							X																		
JAMMT S	1.00	1548							X																		
JAMMT L	1.00	1553							X																		
JAMMF	1.00	1558							X																		
JAMFM	1.00	1602							X																		
ZZZZZZ	1.00	1607																									
CCV	1.00	1611							X																		
CCB	1.00	1616							X																		
JAMLT	1.00	1620							X																		
JAMLF	1.00	1625							X																		
CHULT	1.00	1630							X																		
CHULF	1.00	1634							X																		
JSBHT	1.00	1639							X																		
JSBFL	1.00	1643							X																		
CRCH-	1.00	1648							X																		

U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2360-ICP1100_ Method: P_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
CRCHF	1.00	1652							X																	
CRTFB	1.00	1657							X																	
JSBMT	1.00	1702							X																	
CCV	1.00	1706							X																	
CCB	1.00	1711							X																	
JBMTF	1.00	1716							X																	
CHRDM	1.00	1720							X																	
CHRDF	1.00	1725							X																	
JSBLT	1.00	1729							X																	
JSBLF	1.00	1734							X																	
CRHLT	1.00	1739							X																	
CRHLF	1.00	1743							X																	
ICSA	1.00	1748							X																	
ICSAB	1.00	1753							X																	
CRI	1.00	1757							X																	
CCV	1.00	1802							X																	
CCB	1.00	1806							X																	

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4724-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1454				X																					
S10	1.00	1458				X																					
S20	1.00	1502				X																					
S40	1.00	1506				X																					
S60	1.00	1511				X																					
S80	1.00	1515				X																					
ICV	1.00	1520				X																					
ICB -2.5	1.00	1524				X																					
CRA	1.00	1529				X																					
CCV	1.00	1533				X																					
CCB	1.00	1538				X																					
ZZZZZZ	5.00	1542																									
ZZZZZ	5.00	1547																									
W	1.00	1551				X																					
PBWA	1.00	1556	91.0			X																					
LCSW	1.00	1600				X																					
LCSWA	1.00	1605	106.6			X																					
JAMMT	1.00	1609				X																					
JAMMTA	1.00	1614	102.5			X																					
JAMMF	1.00	1618				X																					
JAMMFA	1.00	1623	96.3			X																					
CCV	1.00	1627				X																					
CCB	1.00	1632				X																					
JAMMT S	1.00	1636				X																					
ZZZZZZ	1.00	1640																									
JAMMT D	1.00	1645				X																					
JAMMT DA	1.00	1649	96.3			X																					
JAMFM	1.00	1654				X																					
JAMFMA	1.00	1658	97.0			X																					
ZZZZZZ	1.00	1702																									
ZZZZZZ	1.00	1707																									
JAMLT	1.00	1711				X																					

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4724-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
JAMLTA	1.00	1716	102.1			X																				
CCV	1.00	1720				X																				
CCB 3	1.00	1724				X																				
JAMLF	1.00	1729				X																				
JAMLFA	1.00	1733	100.2			X																				
CHULT	1.00	1737				X																				
CHULTA	1.00	1742	101.2			X																				
CHULF	1.00	1746				X																				
CHULFA	1.00	1751	90.3			X																				
JSBHT	1.00	1755				X																				
JSBHTA	1.00	1800	104.7			X																				
JSBFL	1.00	1804				X																				
JSBFLA	1.00	1808	100.4			X																				
CCV	1.00	1813				X																				
CCB 4	1.00	1817				X																				
CRCH	1.00	1821				X																				
CRCH-A	1.00	1826	94.9			X																				
CRCHF	1.00	1830				X																				
CRCHFA	1.00	1835	91.6			X																				
CRTFB	1.00	1840				X																				
CRTFBA	1.00	1844	90.4			X																				
JSBMT	1.00	1849				X																				
JSBMTA	1.00	1853	102.6			X																				
JBMTF	1.00	1858				X																				
JBMTFA	1.00	1902	92.6			X																				
CCV	1.00	1907				X																				
CCB 5	1.00	1911				X																				
CHRDM	1.00	1916				X																				
CHRDMA	1.00	1920	98.3			X																				
CHRDF	1.00	1925				X																				
CHRDFA	1.00	1929	89.4			X																				
JSBLT	1.00	1934				X																				

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4724-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
JSBLTA	1.00	1938	94.6			X																				
JSBLF	1.00	1943				X																				
JSBLFA	1.00	1947	96.8			X																				
CRHLT	1.00	1952				X																				
CRHLTA	1.00	1956	96.3			X																				
CCV	1.00	2001				X																				
CCB	1.00	2005				X																				
CRHLF	1.00	2009				X																				
CRHLFA	1.00	2014	90.5			X																				
CCV	1.00	2018				X																				
CCB	1.00	2022				X																				

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3812-V40Z GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1220													X												
S5	1.00	1226													X												
S10	1.00	1233													X												
S25	1.00	1240													X												
S40	1.00	1246													X												
S50	1.00	1253													X												
ICV	1.00	1259													X												
ICB	1.00	1305													X												
CRA	1.00	1312													X												
CRAA	1.00	1319																									
CCV	1.00	1325													X												
CCB	1.00	1331													X												
ZZZZZZ	200.00	1338																									
ZZZZZZ	200.00	1345																									
PBW	1.00	1351													X												
PBWA	1.00	1358	102.4												X												
LCSW	1.00	1404													X												
LCSW	1.00	1411	105.6												X												
JAMMT	1.00	1417													X												
JAMMTA	1.00	1424	88.6												X												
JAMMF	1.00	1430													X												
JAMMFA	1.00	1437	90.3												X												
CCV	1.00	1443													X												
CCB	1.00	1450													X												
JAMMT S	1.00	1456													X												
ZZZZZZ	1.00	1503																									
JAMMT D	1.00	1509													X												
JAMMT DA	1.00	1516	88.2												X												
JAMFM	1.00	1522													X												
JAMFMA	1.00	1529	92.3												X												
ZZZZZZ	1.00	1535																									
ZZZZZZA	1.00	1542	106.2																								

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3812-V40Z GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
JAMLT	1.00	1548													X												
JAMLTA	1.00	1555	93.2												X												
CCV	1.00	1601													X												
CCB	1.00	1607													X												

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3812-V40Z GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1808		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
S5	1.00	1815		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
S10	1.00	1822		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
S25	1.00	1828		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
S40	1.00	1835		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
S50	1.00	1841		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
ICV	1.00	1848		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
ICB	1.00	1854		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRA	1.00	1900		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRAA	1.00	1907		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
CCV	1.00	1914		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CCB	1.00	1920		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
JAMLF	1.00	1926		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
JAMLFA	1.00	1933	84.1	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CHULT	1.00	1939		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CHULTA	1.00	1946	88.2	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CHULF	1.00	1952		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CHULFA	1.00	1959	91.2	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
JSBHT	1.00	2006		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
JSBHTA	1.00	2012	84.4	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
JSBFL	1.00	2019		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
JSBFLA	1.00	2025	84.8	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CCV	1.00	2032		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CCB	1.00	2038		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRCH-	1.00	2044		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRCH-A	1.00	2051	89.5	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRCHF	1.00	2057		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRCHFA	1.00	2104	91.4	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRTFB	1.00	2110		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
CRTFBA	1.00	2117	105.8	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-
JSBMT	1.00	2123		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
JSBMTA	1.00	2130	83.8	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

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AR318868

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3812-V40Z GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
JBMTF	1.00	2136													X												
JBMTFA	1.00	2143	86.3												X												
CCV	1.00	2150													X												
CCB	1.00	2156													X												
CHRDM	1.00	2202													X												
CHRDMA	1.00	2209	90.5												X												
CHRDF	1.00	2215													X												
CHRDFA	1.00	2222	90.4												X												
JSBLT	1.00	2228													X												
JSBLTA	1.00	2235	90.0												X												
JSBLF	1.00	2241													X												
JSBLFA	1.00	2248	88.5												X												
HLT	1.00	2255													X												
HLTA	1.00	2302	93.1												X												
CCV	1.00	2308													X												
CCB	1.00	2314													X												
CRHLF	1.00	2321													X												
CRHLFA	1.00	2327	95.5												X												
CCV	1.00	2334													X												
CCB	1.00	2340													X												

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U.S. EPA - CLP

14

ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3812-V40Z GF_ Method: F_

Start Date: 12/23/92 End Date: 12/23/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1609		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
S5	1.00	1613		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
S10	1.00	1616		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
S25	1.00	1619		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
S40	1.00	1623		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
S50	1.00	1626		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
ICV	1.00	1629		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
ICB	1.00	1632		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
CRA	1.00	1635		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
CCV	1.00	1639		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
CCB	1.00	1642		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBHT0	1.00	1645		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBHT1	1.00	1648		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBHT2	1.00	1651		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBHT3	1.00	1654		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBMT0	1.00	1658		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBMT1	1.00	1701		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBMT2	1.00	1704		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
JSBMT3	1.00	1707		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ0	1.00	1710		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ1	1.00	1713		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ2	1.00	1716		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ3	1.00	1719		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ0	1.00	1722		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ1	1.00	1725		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ2	1.00	1729		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ3	1.00	1732		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ0	1.00	1735		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ1	1.00	1738		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ2	1.00	1741		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
ZZZZZZ3	1.00	1744		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-
CCV	1.00	1748		-	-	-	-	-	-	-	-	-	-	-	X	-	-	-	-	-	-	-	-	-	-	-	-

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: __________

Lab Code: ______ Case No.: ______ SAS No.: ______ SDG No.:CRT02_

Instrument ID Number: 3812-V40Z GF_ Method: F_

Start Date: 12/23/92 End Date: 12/23/92

EPA Sample No.	D/F	Time	% R	Analytes																							
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
CCB	1.00	1751													X												

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75

AR318871

U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4725-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1258																		X						
S5	1.00	1303																		X						
S15	1.00	1307																		X						
S25	1.00	1312																		X						
S40	1.00	1317																		X						
S50	1.00	1322																		X						
ICV	1.00	1327																		X						
ICB	1.00	1332																		X						
CRA	1.00	1336																		X						
CCV	1.00	1341																		X						
CCB	1.00	1346																		X						
PBW	1.00	1351																		X						
PBWA	1.00	1356	94.8																	X						
LCSW	1.00	1401																		X						
LCSWA	1.00	1406	86.6																	X						
JAMMT	1.00	1411																		X						
JAMMTA	1.00	1416	79.8																	X						
JAMMF	1.00	1421																		X						
JAMMFA	1.00	1426	70.1																	X						
JAMMT S	1.00	1431																		X						
ZZZZZZ	1.00	1436																								
CCV	1.00	1440																		X						
CCB	1.00	1445																		X						
JAMMT D	1.00	1450																		X						
JAMMT DA	1.00	1455	82.1																	X						
JAMFM	1.00	1500																		X						
JAMFMA	1.00	1505	70.0																	X						
ZZZZZZ	1.00	1509																								
ZZZZZZ	1.00	1514																								
JAMLT	1.00	1519																		X						
JAMLTA	1.00	1524	77.6																	X						
JAMLF	1.00	1529																		X						

FORM XIV - IN ILM02.1

U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4725-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
JAMLFA	1.00	1534	80.4																	X						
CCV	1.00	1539																		X						
CCB	1.00	1543																		X						
CHULT	1.00	1548																		X						
CHULTA	1.00	1553	91.7																	X						
CHULF	1.00	1557																		X						
CHULFA	1.00	1602	91.6																	X						
JSBHT	1.00	1607																		X						
JSBHTA	1.00	1612	[illegible]																	X						
JSBFL	1.00	1617																		X						
JSBFLA	1.00	1622	82.7																	X						
CRCH-	1.00	1627																		X						
CH-A	1.00	1632	89.4																	X						
V	1.00	1637																		X						
CCB	1.00	1641																		X						
CRCHF	1.00	1646																								
CRCHFA	1.00	1651	82.5																							
CRTFB	1.00	1656																		X						
CRTFBA	1.00	1701	89.5																	X						
JSBMT	1.00	1706																		X						
JSBMTA	1.00	1711	82.0																	X						
JBMTF	1.00	1716																		X						
JBMTFA	1.00	1721	86.7																	X						
CHRDM	1.00	1726																		X						
CHRDMA	1.00	1731	72.9																	X						
CCV	1.00	1736																		X						
CCB	1.00	1740																		X						
CHRDF	1.00	1745																		X						
CHRDFA	1.00	1750	99.3																	X						
JSBLT	1.00	1755																		X						
JSBLTA	1.00	1800	67.7																	X						
JSBLF	1.00	1805																								

FORM XIV - IN ILM02.1

U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4725-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
JSBLFA	1.00	1810	81.3																							
CRHLT	1.00	1815																		X						
CRHLTA	1.00	1820	90.9																	X						
CRHLF	1.00	1825																		X						
CRHLFA	1.00	1830	94.8																	X						
CCV	1.00	1835																		X						
CCB	1.00	1839																		X						

FORM XIV - IN ILM02.1

U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4725-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																							
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	2027																		X						
S5	1.00	2031																		X						
S15	1.00	2036																		X						
S25	1.00	2041																		X						
S40	1.00	2045																		X						
S50	1.00	2050																		X						
ICV	1.00	2055																		X						
ICB	1.00	2100																		X						
CRA	1.00	2105																		X						
CCV	1.00	2110																		X						
CCB	1.00	2114																		X						
CRCHF	1.00	2119																		X						
CHFA	1.00	2124	103.9																	X						
BLF	1.00	2129																								
JSBLFA	1.00	2134	0.5																							
CCV	1.00	2139																		X						
CCB	1.00	2144																		X						

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U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 4725-PE51 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/22/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	2240																		X						
S5	1.00	2245																		X						
S15	1.00	2250																		X						
S25	1.00	2254																		X						
S40	1.00	2259																		X						
S50	1.00	2304																		X						
ICV	1.00	2309																		X						
ICB	1.00	2314																		X						
CRA	1.00	2319																		X						
CCV	1.00	2323																		X						
CCB	1.00	2328																		X						
JSBLF	1.00	2333																		X						
JSBLFA	1.00	2338	86.0																	X						
ZZZZZZ	1.00	2343																								
ZZZZZZ	1.00	2348																								
ZZZZZZ	50.00	2353																								
ZZZZZZ	50.00	2358																								
ZZZZZZ	1.00	0003																								
ZZZZZZ	1.00	0008																								
ZZZZZZ	1.00	0013																								
ZZZZZZ	1.00	0018																								
CCV	1.00	0023																		X						
CCB	1.00	0027																		X						

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U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2803-V400 GF_ Method: F_

Start Date: 12/21/92 End Date: 12/21/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1954																					X			
S10	1.00	2001																					X			
S25	1.00	2007																					X			
S50	1.00	2014																					X			
S75	1.00	2021																					X			
S100	1.00	2027																					X			
ICV	1.00	2033																					X			
ICB	1.00	2040																					X			
CRA	1.00	2046																					X			
CRAA	1.00	2053																								
CCV	1.00	2059																					X			
CCB	1.00	2105																					X			
[illegible]	1.00	2112																					X			
[illegible]NA	1.00	2118	112.9																				X			
LCSW	1.00	2124																					X			
LCSWA	1.00	2131	132.2																				X			
JAMMT	1.00	2137																					X			
JAMMTA	1.00	2144	110.3																				X			
JAMMF	1.00	2150																					X			
JAMMFA	1.00	2157	117.2																				X			
JAMMT S_	1.00	2203																					X			
ZZZZZZ	1.00	2210																								
CCV	1.00	2216																					X			
CCB	1.00	2222																					X			

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2803-V400 GF_ Method: F_

Start Date: 12/22/92 End Date: 12/22/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	0119		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
S10	1.00	0125		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
S25	1.00	0132		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
S50	1.00	0139		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
S75	1.00	0146		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
S100	1.00	0152		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
ICV	1.00	0158		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
ICB	1.00	0204		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRA	1.00	0211		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRAA	1.00	0217		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-		-	-	-
CCV	1.00	0224		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CCB	1.00	0230		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CHULT	1.00	0236		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CHULTA	1.00	0243	111.0	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CHULF	1.00	0249		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CHULFA	1.00	0256	99.5	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JSBHT	1.00	0302		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JSBHTA	1.00	0309	104.1	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JSBFL	1.00	0315		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JSBFLA	1.00	0322	110.8	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRCH-	1.00	0328		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRCH-A	1.00	0335	106.4	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CCV	1.00	0341		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CCB	1.00	0347		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRCHF	1.00	0353		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRCHFA	1.00	0400	104.3	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRTFB	1.00	0406		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
CRTFBA	1.00	0413	106.2	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JSBMT	1.00	0419		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JSBMTA	1.00	0426	116.8	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JBMTF	1.00	0432		-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
JBMTFA	1.00	0438	108.7	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	X	-	-	-
				-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-	-

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2803-V400 GF_ Method: F_

Start Date: 12/22/92 End Date: 12/22/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
CHRDM	1.00	0444																					X			
CHRDMA	1.00	0451	106.3																				X			
CCV	1.00	0457																					X			
CCB	1.00	0504																					X			
CHRDF	1.00	0510																					X			
CHRDFA	1.00	0516	106.8																				X			
JSBLT	1.00	0523																					X			
JSBLTA	1.00	0529	108.7																				X			
JSBLF	1.00	0536																					X			
JSBLFA	1.00	0542	108.5																				X			
CRHLT	1.00	0548																					X			
CRHLTA	1.00	0555	119.3																				X			
[illegible]LF	1.00	0601																					X			
[illegible]LFA	1.00	0608	114.6																				X			
CCV	1.00	0614																					X			
CCB	1.00	0621																					X			

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U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 2803-V400 GF_ Method: F_

Start Date: 12/22/92 End Date: 12/22/92

EPA Sample No.	D/F	Time	% R	Analytes																									
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	1709																					X			
S10	1.00	1716																					X			
S25	1.00	1723																					X			
S50	1.00	1729																					X			
S75	1.00	1736																					X			
S100	1.00	1742																					X			
ICV	1.00	1749																					X			
ICB	1.00	1755																					X			
CRA	1.00	1801																					X			
CRAA	1.00	1808																								
CCV	1.00	1814																					X			
CCB	1.00	1821																					X			
JAMMT D	1.00	1827																					X			
JAMMT DA	1.00	1833	115.3																				X			
JAMFM	1.00	1840																					X			
JAMFMA	1.00	1846	107.6																				X			
ZZZZZZ	1.00	1853																								
ZZZZZZ	1.00	1859																								
JAMLT	1.00	1905																					X			
JAMLTA	1.00	1912	110.4																				X			
JAMLF	1.00	1918																					X			
JAMLFA	1.00	1925	103.0																				X			
CCV	1.00	1931																					X			
CCB	1.00	1938																					X			

FORM XIV - IN ILM02.1

84

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14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3998-V20 AA__ Method: CV

Start Date: 12/22/92 End Date: 12/22/92

EPA Sample No.	D/F	Time	% R	Analytes																								
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
S0	1.00	0900																X									
S0.2	1.00	0901																X									
S0.5	1.00	0903																X									
S1.0	1.00	0904																X									
S2.5	1.00	0906																X									
S5.0	1.00	0907																X									
ICV	1.00	0909																X									
ICB	1.00	0910																X									
CRA	1.00	0912																X									
CCV	1.00	0913																X									
CCB	1.00	0915																X									
PBW	1.00	0916																X									
SW	1.00	0918																X									
MMT	1.00	0919																X									
JAMMF	1.00	0921																X									
JAMMT S	1.00	0922																X									
JAMMT D	1.00	0924																X									
JAMFM	1.00	0925																X									
ZZZZZZ	1.00	0927																									
JAMLT	1.00	0928																X									
JAMLF	1.00	0930																X									
CCV	1.00	0931																X									
CCB	1.00	0933																X									
CHULT	1.00	0934																X									
CHULF	1.00	0936																X									
JSBHT	1.00	0937																X									
JSBFL	1.00	0939																X									
CRCH-	1.00	0940																X									
CRCHF	1.00	0942																X									
CRTFB	1.00	0943																X									
JSBMT	1.00	0945																X									
JBMTF	1.00	0946																X									

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U.S. EPA - CLP

14
ANALYSIS RUN LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.:CRT02_

Instrument ID Number: 3998-V20 AA__ Method: CV

Start Date: 12/22/92 End Date: 12/22/92

EPA Sample No.	D/F	Time	% R	Analytes																										
				AL	SB	AS	BA	BE	CD	CA	CR	CO	CU	FE	PB	MG	MN	HG	NI	K	SE	AG	NA	TL	V	ZN	CN
CHRDM	1.00	0948																X									
CCV	1.00	0949																X									
CCB	1.00	0951																X									
CHRDF	1.00	0952																X									
JSBLT	1.00	0954																X									
JSBLF	1.00	0955																X									
CRHLT	1.00	0957																X									
CRHLF	1.00	0958																X									
ZZZZZZ	1.00	1000																									
ZZZZZZ	1.00	1001																									
ZZZZZZ	1.00	1003																									
ZZZZZZ	1.00	1004																									
ZZZZZZ	1.00	1006																									
CCV	1.00	1007																X									
CCB	1.00	1009																X									

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4

ICP INTERFERENCE CHECK SAMPLE

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 ICS Source: LLI_____

Concentration Units: ug/L

Analyte	True Sol. A	True Sol. AB	Initial Found Sol. A	Initial Found Sol. AB	Initial Found %R	Final Found Sol. A	Final Found Sol. AB	Final Found %R
Aluminum	500000	500000	475510	479310.0	95.9	467040	470970.0	94.2
Antimony								
Arsenic								
Barium	0	500	7	488.3	97.7	6	480.2	96.0
Beryllium	0	500	1	460.5	92.1	1	450.9	90.2
Cadmium								
Calcium	500000	500000	474510	476850.0	95.4	461270	464760.0	93.0
Chromium	0	500	-17	446.7	89.3	-16	438.7	87.7
Cobalt	0	500	2	448.6	89.7	1	429.3	85.9
Copper	0	500	4	468.8	93.8	1	458.4	91.7
Iron	200000	200000	181120	182200.0	91.1	176770	178300.0	89.2
Lead								
Magnesium	500000	500000	489920	493300.0	98.7	482410	485870.0	97.2
Manganese	0	500	-13	441.4	88.3	-12	432.4	86.5
Mercury								
Nickel	0	1000	-25	845.2	84.5	-26	830.0	83.0
Potassium								
Selenium								
Silver	0	1000	-33	901.6	90.2	-34	880.0	88.0
Sodium								
Thallium								
Vanadium	0	500	4	474.2	94.8	4	466.4	93.3
Zinc	0	1000	-8	907.6	90.8	-10	896.2	89.6

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4

ICP INTERFERENCE CHECK SAMPLE

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 ICS Source: LLI_____

Concentration Units: ug/L

Analyte	True Sol. A	True Sol. AB	Initial Found Sol. A	Initial Found Sol. AB	Initial Found %R	Final Found Sol. A	Final Found Sol. AB	Final Found %R
Aluminum								
Antimony								
Arsenic								
Barium								
Beryllium								
Cadmium	0	1000	29	869.2	86.9	29	860.3	86.0
Calcium								
Chromium								
Cobalt								
Copper								
Iron								
Lead								
Magnesium								
Manganese								
Mercury								
Nickel								
Potassium								
Selenium								
Silver								
Sodium								
Thallium								
Vanadium								
Zinc								

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5A
SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

JAMMT S

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER_ Level (low/med): LOW___

% Solids for Sample: __0.0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Spike Added (SA)	%R	Q	M
Aluminum	75-125	3540.1000		1989.0000		2000.00	77.6		P
Antimony	75-125	496.7000		22.0000	U	500.00	99.3		P
Arsenic	75-125	43.2400		1.0000	U	40.00	108.1		F
Barium	75-125	1960.9000		68.8800	B	2000.00	94.6		P
Beryllium	75-125	48.7100		1.0000	U	50.00	97.4		P
Cadmium	75-125	47.0500		3.0000	U	50.00	94.1		P
Calcium									NR
Chromium	75-125	205.3600		8.0000	U	200.00	102.7		P
Cobalt	75-125	484.1700		5.0000	U	500.00	96.8		P
Copper	75-125	247.6200		5.0000	U	250.00	99.0		P
Iron	75-125	2841.2000		2221.1000		1000.00	62.0	N	P
Lead	75-125	20.0800		2.8700	B	20.00	86.0		F
Magnesium									NR
Manganese	75-125	581.8100		112.5400		500.00	93.9		P
Mercury	75-125	1.0100		0.1000	U	1.00	101.0		CV
Nickel	75-125	494.5100		8.0000	U	500.00	98.9		P
Potassium									NR
Selenium	75-125	9.2900		1.0000	U	10.00	92.9		F
Silver	75-125	43.6700		11.0000	U	50.00	87.3		P
Sodium									NR
Thallium	75-125	53.1000		2.0000	U	50.00	106.2		F
Vanadium	75-125	481.3000		6.0000	U	500.00	96.3		P
Zinc	75-125	558.5400		74.8600		500.00	96.7		P
Cyanide									NR

Comments:

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5B

POST DIGEST SPIKE SAMPLE RECOVERY

EPA SAMPLE NO.

JAMMT A

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water) : WATER_ Level (low/med): LOW___

Concentration Units: ug/L

Analyte	Control Limit %R	Spiked Sample Result (SSR)	C	Sample Result (SR)	C	Added (SA)	%R	Q	M
Aluminum									NR
Antimony									NR
Arsenic									NR
Barium									NR
Beryllium									NR
Cadmium									NR
Calcium									NR
Chromium									NR
Cobalt									NR
Copper									NR
Iron		2465.50		2221.10		500.0	48.9		P
Lead									NR
Magnesium									NR
Manganese									NR
Mercury									NR
Nickel									NR
Potassium									NR
Selenium									NR
Silver									NR
Sodium									NR
Thallium									NR
Vanadium									NR
Zinc									NR
Cyanide									NR

Comments:

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6
DUPLICATES

EPA SAMPLE NO.

JAMMT D

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Level (low/med): _LOW__

% Solids for Sample: ____0 % Solids for Duplicate: _____0

Concentration Units (ug/L or mg/kg dry weight): UG/L_

Analyte	Control Limit	Sample (S)	C	Duplicate (D)	C	RPD	Q	M
Aluminum		1989.0000		2121.7000		6.5		P
Antimony		22.0000	U	22.0000	U			P
Arsenic		1.0000	U	1.0000	U			F
Barium		68.8800	B	69.6900	B	1.2		P
Beryllium		1.0000	U	1.0000	U			P
Cadmium		3.0000	U	3.0000	U			P
Calcium	5000.0	17637.0000		18093.0000		2.6		P
Chromium		8.0000	U	8.0000	U			P
Cobalt		5.0000	U	5.0000	U			P
Copper		5.0000	U	5.0000	U			P
Iron		2221.1000		2372.3000		6.6		P
Lead	3.0	2.8700	B	5.8100		67.7		F
Magnesium	5000.0	7131.8000		7331.4000		2.8		P
Manganese		112.5400		116.6000		3.5		P
Mercury		0.1000	U	0.1000	U			CV
Nickel		8.0000	U	8.0000	U			P
Potassium	5000.0	5921.4000		6049.5000		2.1		P
Selenium		1.0000	U	1.0000	U			F
Silver		11.0000	U	11.0000	U			P
Sodium	5000.0	10436.0000		10601.0000		1.6		P
Thallium		2.0000	U	2.0000	U			F
Vanadium		6.0000	U	6.0000	U			P
Zinc	20.0	74.8600		78.5600		4.8		P
Cyanide								NR

LSCRDL

7

LABORATORY CONTROL SAMPLE

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Solid LCS Source: _____

Aqueous LCS Source: LLI_____

Analyte	Aqueous (ug/L) True	Found	%R	Solid (mg/kg) True	Found	C	Limits		%R
Aluminum	2000.0	1920.60	96.0						
Antimony	500.0	521.37	104.3						
Arsenic	40.0	40.66	101.6						
Barium	2000.0	2024.00	101.2						
Beryllium	50.0	52.59	105.2						
Cadmium	50.0	47.79	95.6						
Calcium	4000.0	4075.90	101.9						
Chromium	200.0	216.52	108.3						
Cobalt	500.0	511.08	102.2						
Copper	250.0	260.88	104.4						
Iron	1000.0	1040.60	104.1						
Lead	20.0	20.79	104.0						
Magnesium	2000.0	1916.60	95.8						
Manganese	500.0	518.94	103.8						
Mercury	1.0	1.02	102.0						
Nickel	500.0	524.62	104.9						
Potassium	4000.0	4020.40	100.5						
Selenium	10.0	10.17	101.7						
Silver	50.0	45.92	91.8						
Sodium	4000.0	4421.70	110.5						
Thallium	50.0	51.20	102.4						
Vanadium	500.0	506.61	101.3						
Zinc	500.0	528.94	105.8						
Cyanide									

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8

STANDARD ADDITION RESULTS

Lab Name: LANCASTER_LABORATORIES___ Contract:_____

Lab Code: _____ Case No.: _____ SAS No.:_____ SDG No.:CRT02_

Concentration Units: ug/L

EPA Sample No.	An	0 ADD ABS	1 ADD CON	1 ADD ABS	2 ADD CON	2 ADD ABS	3 ADD CON	3 ADD ABS	Final Conc.	r	Q
JSBHT	PB	0.011	5.00	0.028	10.00	0.044	15.00	0.057	3.9	0.9982	
JSBMT	PB	0.014	5.00	0.028	10.00	0.044	15.00	0.056	5.0	0.9986	

FORM VIII - IN ILM02.1

U.S. EPA - CLP

9

ICP SERIAL DILUTION

EPA SAMPLE NO.

JAMMT L

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

Matrix (soil/water): WATER Level (low/med): LOW__

Concentration Units: ug/L

Analyte	Initial Sample Result (I)	C	Serial Dilution Result (S)	C	% Difference	Q	M
Aluminum	1989.00		2225.35		11.9		P
Antimony	22.00	U	110.00	U			P
Arsenic							
Barium	68.88	B	72.05	B	4.6		P
Beryllium	1.00	U	5.00	U			P
Cadmium	3.00	U	15.00	U			P
Calcium	17637.00		17581.50	B	0.3		P
Chromium	8.00	U	40.00	U			P
Cobalt	5.00	U	25.00	U			P
Copper	5.00	U	25.00	U			P
Iron	2221.10		2217.95		0.1		P
Lead							
Magnesium	7131.80		7040.50	B	1.3		P
Manganese	112.54		109.65		2.6		P
Mercury							
Nickel	8.00	U	40.00	U			P
Potassium	5921.40		6023.50	B	1.7		P
Selenium							
Silver	11.00	U	55.00	U			P
Sodium	10436.00		10744.50	B	3.0		P
Thallium							
Vanadium	6.00	U	30.00	U			P
Zinc	74.86		121.00		61.6		P

FORM IX - IN

ILM02.1

VERIFICATION OF INSTRUMENT PARAMETERS

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 Date: 10/15/92

Flame AA ID Number : _____

Furnace AA ID Number : _____

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum	308.22		200	57.0	P
Antimony	206.84		60	22.0	P
Arsenic			10		NR
Barium	493.41		200	4.0	P
Beryllium	313.04		5	1.0	P
Cadmium	226.50		5	3.0	P
Calcium	317.93		5000	23.0	P
Chromium	267.72		10	8.0	P
Cobalt	228.62		50	5.0	P
Copper	324.75		25	5.0	P
Iron	259.94		100	7.0	P
Lead			3		NR
Magnesium	279.08		5000	45.0	P
Manganese	257.61		15	1.0	P
Mercury			0.2		NR
Nickel	231.60		40	8.0	P
Potassium	766.49		5000	132.0	P
Selenium			5		NR
Silver	328.07		10	11.0	P
Sodium	589.59		5000	72.0	P
Thallium			10		NR
Vanadium	292.40		50	6.0	P
Zinc	213.86		20	4.0	P

Comments:

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: _____ Date: 10/15/92

Flame AA ID Number : 3998-V20_AA_

Furnace AA ID Number : _____

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury	253.65		0.2	0.1	CV
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: _____ Date: 10/15/92

Flame AA ID Number : _____

Furnace AA ID Number : 2803-V400_GF

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium	276.80	BD	10	2.0	F
Vanadium			50		NR
Zinc			20		NR

Comments:

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: _____ Date: 10/15/92

Flame AA ID Number : _____

Furnace AA ID Number : 3812-V40Z_GF

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead	283.30	BZ	3	1.0	F
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

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U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: _____ Date: 10/15/92

Flame AA ID Number : _____

Furnace AA ID Number : 4724-PE51_GF

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic	193.70	BZ	10	1.0	F
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium			5		NR
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

U.S. EPA - CLP

10

Instrument Detection Limits (Quarterly)

Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: _____ Date: 10/15/92

Flame AA ID Number : _____

Furnace AA ID Number : 4725-PE51_GF

Analyte	Wave-length (nm)	Back-ground	CRDL (ug/L)	IDL (ug/L)	M
Aluminum			200		NR
Antimony			60		NR
Arsenic			10		NR
Barium			200		NR
Beryllium			5		NR
Cadmium			5		NR
Calcium			5000		NR
Chromium			10		NR
Cobalt			50		NR
Copper			25		NR
Iron			100		NR
Lead			3		NR
Magnesium			5000		NR
Manganese			15		NR
Mercury			0.2		NR
Nickel			40		NR
Potassium			5000		NR
Selenium	196.00	BZ	5	1.0	F
Silver			10		NR
Sodium			5000		NR
Thallium			10		NR
Vanadium			50		NR
Zinc			20		NR

Comments:

FORM X - IN ILMO2.1

U.S. EPA - CLP

11A

ICP INTERELEMENT CORRECTION FACTORS (ANNUALLY)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 Date: 01/15/92

Analyte	Wave-length (nm)	Interelement Correction Factors for : Al	Ca	Fe	Mg	BE_
Aluminum_	308.22	0.0000000	0.0001640	0.0000000	0.0002340	0.0000000
Antimony_	206.84	-0.0000120	0.0000000	0.0001220	0.0000000	0.0000000
Arsenic__	193.70	-0.0071120	0.0000000	-0.0071120	0.0000000	0.0000000
Barium___	493.41	0.0000000	0.0000000	-0.0000320	0.0000000	0.0000000
Beryllium	313.04	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Cadmium__	226.50	-0.0000390	0.0000000	0.0000710	0.0000090	0.0000000
Calcium__	317.93	0.0000000	0.0000000	-0.0000170	0.0003960	0.0000000
Chromium_	267.72	0.0000000	0.0000000	0.0000050	0.0000080	0.0000000
Cobalt___	228.62	0.0000000	0.0000000	-0.0000040	-0.0000020	0.0000000
Copper___	324.75	0.0000000	0.0000000	0.0000110	0.0000000	0.0000000
Iron_____	259.94	0.0000000	0.0000320	0.0000000	0.0000600	0.0000000
Lead_____	220.35	-0.0009830	-0.0000110	0.0000000	-0.0000190	0.0000000
Magnesium	279.08	0.0000000	0.0000890	0.0002080	0.0000000	-0.003141
Manganese	257.61	0.0005890	0.0000000	0.0000460	0.0000280	0.000000
Mercury__						
Nickel___	231.60	0.0000100	0.0000000	0.0000810	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0002140	0.0000000	-0.0000270	0.0000000
Selenium_	196.03	0.0002130	0.0000170	0.0000420	-0.0000810	0.0000000
Silver___	328.07	0.0000000	0.0000170	0.0000630	0.0000000	0.0000000
Sodium___	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium_	377.57	0.0006790	-0.0002420	0.0000460	0.0000110	0.0000000
Vanadium_	292.40	0.0000000	0.0000000	-0.0000300	0.0000000	0.0000000
Zinc_____	213.86	0.0000350	0.0000000	0.0001160	0.0000380	0.0000000

Comments:

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U.S. EPA - CLP

11B
ICP INTERELEMENT CORRECTION FACTORS (ANNUALLY)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 Date: 01/15/92

Analyte	Wavelength (nm)	Interelement Correction Factors for : CO_	CR_	CU_	K__	MN_
Aluminum_	308.22	0.0001860	0.0002450	0.0000420	0.0000000	0.0013450
Antimony_	206.84	-0.0038310	0.0070110	-0.0000150	0.0000000	-0.0001910
Arsenic__	193.70	-0.0030800	0.0009500	0.0000180	0.0000000	-0.0004130
Barium___	493.41	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Beryllium	313.04	0.0000000	0.0000180	0.0000000	0.0000000	0.0000000
Cadmium__	226.50	0.0000320	-0.0000120	0.0000120	0.0000000	0.0000000
Calcium__	317.93	0.0000000	0.0009140	0.0000000	0.0000000	0.0003530
Chromium_	267.72	-0.0000760	0.0000000	-0.0000150	0.0000000	-0.0011240
Cobalt___	228.62	0.0000000	0.0000750	0.0000000	0.0000000	0.0000000
Copper___	324.75	0.0000570	0.0000000	0.0000000	0.0000000	0.0000390
Iron_____	259.94	0.0000810	0.0001150	0.0000000	0.0000000	0.0008980
Lead_____	220.35	0.0002700	-0.0010390	-0.0000110	0.0000000	0.0000920
[illegible]nesium	279.08	-0.0018810	-0.0007860	0.0000000	0.0001870	-0.0080910
[illegible]ganese	257.61	0.0000000	0.0000000	-0.0000770	0.0000000	0.0000000
Mercury__						
Nickel___	231.60	0.0015590	0.0000000	0.0000000	0.0000000	0.0000000
Potassium	766.49	0.0000000	0.0000000	0.0001330	0.0000000	0.0000000
Selenium_	196.03	0.0005990	-0.0002980	0.0004380	0.0000000	0.0002750
Silver___	328.07	0.0000120	0.0000420	0.0000740	0.0000000	0.0001340
Sodium___	589.59	0.0000000	0.0000000	0.0000000	0.0000000	0.0000000
Thallium_	377.57	0.0001330	-0.0007970	0.0000000	0.0000000	-0.0002470
Vanadium_	292.40	0.0000000	-0.0022240	-0.0000640	0.0000000	-0.0001760
Zinc_____	213.86	0.0000000	0.0000620	0.0043410	0.0000000	0.0000000

Comments:

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U.S. EPA - CLP

11B
ICP INTERELEMENT CORRECTION FACTORS (ANNUALLY)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 Date: 01/15/92

Analyte	Wave-length (nm)	Interelement Correction Factors for : MO_	SI_	SR_	TI_	V__
Aluminum_	308.22	0.0089550	0.0105440	0.0000000	0.0024810	0.0228200
Antimony_	206.84	0.0048890	0.0000000	0.0000000	0.0011430	-0.0017710
Arsenic__	193.70	0.0010720	0.0000000	0.0000000	-0.0000120	0.0055100
Barium___	493.41	0.0000120	0.0000000	0.0000000	0.0000000	0.0000000
Beryllium	313.04	0.0000190	0.0000000	0.0000000	-0.0000350	0.0029800
Cadmium__	226.50	0.0001370	0.0000000	0.0000000	0.0002920	-0.0000150
Calcium__	317.93	0.0007860	0.0000000	0.0000000	0.0007140	0.0008280
Chromium_	267.72	-0.0012330	0.0000000	0.0000000	0.0001710	-0.0880700
Cobalt___	228.62	-0.0017320	0.0000000	0.0000000	0.0017060	0.0000300
Copper___	324.75	0.0006890	0.0000000	0.0000000	0.0003140	0.0000620
Iron_____	259.94	0.0003620	0.0000000	0.0000000	0.0002420	0.0002960
Lead_____	220.35	-0.0019300	0.0000000	0.0000000	0.0003340	-0.0003050
Magnesium	279.08	-0.0215500	0.0000000	0.0001940	-0.0036750	-0.0003240
Manganese	257.61	-0.0001860	0.0000000	0.0000000	0.0000430	-0.0001500
Mercury__						
Nickel___	231.60	-0.0026600	0.0000000	0.0000000	0.0000000	-0.0004560
Potassium	766.49	0.0018950	0.0000000	0.0000000	0.0002270	0.0002670
Selenium_	196.03	0.0001070	0.0000000	0.0000000	-0.0003270	0.0002030
Silver___	328.07	0.0006400	0.0000000	0.0000000	0.0002520	0.0002910
Sodium___	589.59	0.0004883	0.0000000	0.0000000	-0.0049410	0.0007290
Thallium_	377.57	0.0000000	0.0000000	0.0023740	-0.0550300	-0.0042380
Vanadium_	292.40	0.0000000	0.0000000	0.0000000	0.0003140	0.0000000
Zinc_____	213.86	0.0000000	0.0000000	0.0000000	0.0000360	0.0002510

Comments:

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U.S. EPA - CLP

12
ICP LINEAR RANGES (QUARTERLY)

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.: _____ SAS No.: _____ SDG No.: CRT02_

ICP ID Number: 2360-ICP1100 Date: 10/15/92

Analyte	Integ. Time (sec.)	Concentration (ug/L)	M
Aluminum	6.00	600000.0	P
Antimony	6.00	200000.0	P
Arsenic	6.00	200000.0	P
Barium	6.00	100000.0	P
Beryllium	6.00	10000.0	P
Cadmium	6.00	100000.0	P
Calcium	6.00	600000.0	P
Chromium	6.00	100000.0	P
Cobalt	6.00	100000.0	P
Copper	6.00	100000.0	P
Iron	6.00	250000.0	P
Lead	6.00	100000.0	P
Magnesium	6.00	800000.0	P
Manganese	6.00	100000.0	P
Mercury			NR
Nickel	6.00	100000.0	P
Potassium	6.00	900000.0	P
Selenium	6.00	200000.0	P
Silver	6.00	50000.0	P
Sodium	6.00	900000.0	P
Thallium	6.00	200000.0	P
Vanadium	6.00	100000.0	P
Zinc	6.00	100000.0	P

Comments:

PREPARATION AND RUN LOGS

U.S. EPA - CLP

13
PREPARATION LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.:_____ SAS No.: _____ SDG No.:CRT02_

Method: P_

EPA Sample No.	Preparation Date	Weight (gram)	Volume (mL)
CHRDF	12/19/92		100
CHRDM	12/19/92		100
CHULF	12/19/92		100
CHULT	12/19/92		100
CRCH-	12/19/92		100
CRCHF	12/19/92		100
CRHLF	12/19/92		100
CRHLT	12/19/92		100
CRTFB	12/19/92		100
JAMFM	12/19/92		100
JAMLF	12/19/92		100
JAMLT	12/19/92		100
JAMMF	12/19/92		100
JAMMT	12/19/92		100
JAMMT_D	12/19/92		100
JAMMT_S	12/19/92		100
JBMTF	12/19/92		100
JSBFL	12/19/92		100
JSBHT	12/19/92		100
JSBLF	12/19/92		100
JSBLT	12/19/92		100
JSBMT	12/19/92		100
LCSW	12/19/92		100
PBW	12/19/92		100

FORM XIII - IN ILM02.1

U.S. EPA - CLP

13
PREPARATION LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.:_____ SAS No.: _____ SDG No.:CRT02_

Method: F_

EPA Sample No.	Preparation Date	Weight (gram)	Volume (mL)
CHRDF	12/19/92		100
CHRDM	12/19/92		100
CHULF	12/19/92		100
CHULT	12/19/92		100
CRCH-	12/19/92		100
CRCHF	12/19/92		100
CRHLF	12/19/92		100
CRHLT	12/19/92		100
CRTFB	12/19/92		100
JAMFM	12/19/92		100
JAMLF	12/19/92		100
JAMLT	12/19/92		100
JAMMF	12/19/92		100
JAMMT	12/19/92		100
JAMMT_D	12/19/92		100
JAMMT_S	12/19/92		100
JBMTF	12/19/92		100
JSBFL	12/19/92		100
JSBHT	12/19/92		100
JSBLF	12/19/92		100
JSBLT	12/19/92		100
JSBMT	12/19/92		100
LCSW	12/19/92		100
PBW	12/19/92		100

FORM XIII - IN ILM02.1

U.S. EPA - CLP

13
PREPARATION LOG

Lab Name: LANCASTER_LABORATORIES___ Contract: _____

Lab Code: _____ Case No.:_____ SAS No.: _____ SDG No.:CRT02_

Method: CV

EPA Sample No.	Preparation Date	Weight (gram)	Volume (mL)
CHRDF	12/21/92		100
CHRDM	12/21/92		100
CHULF	12/21/92		100
CHULT	12/21/92		100
CRCH-	12/21/92		100
CRCHF	12/21/92		100
CRHLF	12/21/92		100
CRHLT	12/21/92		100
CRTFB	12/21/92		100
JAMFM	12/21/92		100
JAMLF	12/21/92		100
JAMLT	12/21/92		100
JAMMF	12/21/92		100
JAMMT	12/21/92		100
JAMMT_D	12/21/92		100
JAMMT_S	12/21/92		100
JBMTF	12/21/92		100
JSBFL	12/21/92		100
JSBHT	12/21/92		100
JSBLF	12/21/92		100
JSBLT	12/21/92		100
JSBMT	12/21/92		100
LCSW	12/21/92		100
PBW	12/21/92		100

FORM XIII - IN ILM02.1

RAW DATA

ICP DATA

Run Name: 9235402I11
Data Reviewed By: NR/142 12-20-92
Data Verified By: Max Dravely 12-21-92

LANCASTER LABORATORIES

PAGE: 1 INSTRUMENT ID: 02360

1 SO 50 12/19/92 22:15

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	24.00000	8.333	22.0000	26.0000	24.0000		
__Al	554.00000	0.361	556.000	552.000	554.000		
__As	23.66666	6.454	22.0000	24.0000	25.0000		
__B	5.33333	28.64	5.00000	7.00000	4.00000		
__Ba	50.00000	0.000	50.0000	50.0000	50.0000		
__Be	22.00000	0.000	22.0000	22.0000	22.0000		
__Ca	574.00000	0.000	574.000	574.000	574.000		
__Cd	-0.66666	346.4	-2.0000	2.00000	-2.0000		
__Co	-0.14814	1875.	1.11111	1.77777	-3.3333		
__Cr	10.85186	51.66	17.2222	6.66668	8.66668		
__Cu	12.00000	16.66	14.0000	10.0000	12.0000		
__Fe	10.00000	20.00	8.00000	12.0000	10.0000		
__K	254.66667	0.453	254.000	256.000	254.000		
__Li	0.00000	0.000	0.00000	0.00000	0.00000		
__Mg	8.00000	50.00	12.0000	8.00000	4.00000		
__Mn	4.33333	35.25	3.00000	4.00000	6.00000		
__Mo	4.00000	50.00	6.00000	4.00000	2.00000		
__Na	143.00000	0.699	143.000	142.000	144.000		
__Ni	3.66666	68.63	1.00000	6.00000	4.00000		
__P	4.31034	55.25	6.00000	1.58621	5.34483		
__Pb	3.99998	34.64	4.79998	4.79998	2.39999		
__Sb	0.00000	0.000	-2.0000	2.00000	0.00000		
__Se	8.66666	81.04	16.0000	8.00000	2.00000		
__Si	32.66666	9.352	30.0000	32.0000	36.0000		
__Sn	0.66666	173.2	2.00000	0.00000	0.00000		
__Sr	1.00000	100.0	1.00000	0.00000	2.00000		
__Ti	68.00000	7.781	72.0000	62.0000	70.0000		
__Tl	19.33333	48.88	30.0000	16.0000	12.0000		
__V	2.66666	43.30	2.00000	2.00000	4.00000		
__Zn	42.00000	0.000	42.0000	42.0000	42.0000		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 2

INSTRUMENT ID: 02360

2 S1 S 12/19/92 22:19

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ba	8381.33300	1.088	8421.00	8446.00	8277.00	_____	_____
__Cd	15219.00000	0.783	15276.0	15299.0	15082.0	_____	_____
__Cr	21792.14843	0.900	21815.2	21975.8	21585.3	_____	_____
__Cu	16009.00000	1.028	16068.0	16136.0	15823.0	_____	_____
__Fe	11339.00000	0.915	11386.0	11411.0	11220.0	_____	_____
__Mn	8186.33349	0.882	8216.00	8239.00	8104.00	_____	_____
__Mo	70637.00000	0.869	70912.0	71066.0	69933.0	_____	_____
__Ni	21818.33398	0.825	21860.0	21974.0	21621.0	_____	_____
__Pb	18771.19921	0.911	18840.5	18896.5	18576.4	_____	_____
__Zn	10730.00000	0.643	10760.0	10779.0	10651.0	_____	_____

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 3

INSTRUMENT ID: 02360

3 S2 S 12/19/92 22:23

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__As	1663.66662	1.249	1640.00	1672.00	1679.00	_____	_____
__Be	1537.33337	0.762	1524.00	1542.00	1546.00	_____	_____
__Co	3447.81469	1.276	3401.33	3453.22	3488.88	_____	_____
__Li	16497.33398	0.474	16407.0	16538.0	16547.0	_____	_____
__P	5553.24121	0.692	5522.48	5540.89	5596.34	_____	_____
__Sb	9947.66699	0.537	9886.00	9980.00	9977.00	_____	_____
__Se	4008.66674	0.923	3966.00	4028.00	4032.00	_____	_____
__Si	6539.33349	1.139	6464.00	6541.00	6613.00	_____	_____
__Sn	3900.00000	0.712	3868.00	3914.00	3918.00	_____	_____
__Sr	9360.00000	0.669	9288.00	9390.00	9402.00	_____	_____
__Ti	24050.66601	0.674	23880.0	24069.0	24203.0	_____	_____
__Tl	1245.66662	0.858	1240.00	1239.00	1258.00	_____	_____
__V	3479.66674	0.692	3452.00	3491.00	3496.00	_____	_____

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 4 INSTRUMENT ID: 02360

4 . 53 S 12/19/92 22:28

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	1048.66662	1.588	1062.00	1054.00	1030.00	_____	_____
__B	1864.00000	0.107	1862.00	1866.00	1864.00	_____	_____

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 5

INSTRUMENT ID: 02360

5 S4 S 12/19/92 22:31

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__K	13258.66699	0.533	13298.0	13301.0	13177.0	_____	_____
__Na	59632.33203	0.401	59612.0	59881.0	59404.0	_____	_____

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 6 INSTRUMENT ID: 02360

6 S5 S 12/19/92 22:33

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Al	26733.33398	0.925	27014.0	26547.0	26639.0	_____	_____
__Ca	189651.67187	0.946	191691.	188306.	188958.	_____	_____
__Mg	30736.00000	0.900	31034.0	30487.0	30687.0	_____	_____

94

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 7

INSTRUMENT ID: 02360

7 ICV 12/19/92 22:36

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.47806	0.847	0.48130	0.47935	0.47352		
__Al	24.39757	1.063	24.4893	24.5987	24.1045		
__As	2.10778	1.424	2.08332	2.14128	2.09872		
__B	1.00395	1.214	1.00750	1.01398	0.99038		
__Ba	5.07988	1.339	5.10589	5.13109	5.00267		
__Be	0.48935	1.291	0.49084	0.49479	0.48242		
__Ca	50.22468	0.984	50.4205	50.5913	49.6621		
__Cd	4.92920	0.883	4.92393	4.97516	4.88850		
__Co	1.00118	1.135	1.00269	1.01172	0.98914		
__Cr	5.23104	0.980	5.25303	5.26768	5.17243		
__Cu	4.95136	1.256	4.97755	4.99620	4.88032		
__Fe	5.03671	1.053	5.05165	5.08071	4.97777		
__K	50.95428	1.664	51.0310	51.7613	50.0704		
__Li	25.90547	1.461	26.0590	26.1832	25.4740		
__Mg	24.16545	1.188	24.3045	24.3567	23.8351		
__Mn	5.02919	1.029	5.05406	5.06386	4.96966		
__Mo	25.03213	1.037	25.1080	25.2454	24.7428		
__Na	52.50534	0.590	52.5585	52.7854	52.1719		
__Ni	9.92718	1.075	9.96375	10.0108	9.80693		
__P	25.63397	0.610	25.5574	25.8140	25.5304		
__Pb	25.40632	0.962	25.4491	25.6266	25.1431		
__Sb	5.17505	0.965	5.19895	5.20860	5.11761		
__Se	2.07569	3.915	2.08066	2.15437	1.99205		
__Si	1.94346	0.844	1.94521	1.95894	1.92625		
__Sn	5.00657	0.806	5.03725	5.02166	4.96081		
__Sr	1.04500	1.319	1.05020	1.05544	1.02937		
__Ti	2.07159	1.358	2.07938	2.09502	2.04037		
__Tl	5.38027	0.813	5.38778	5.41979	5.33323		
__V	2.07869	1.464	2.09529	2.09722	2.04356		
__Zn	5.02654	0.840	5.04003	5.06039	4.97921		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 8

INSTRUMENT ID: 02360

8 ICB 12/19/92 22:40

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	-0.00162	151.1	0.00097	-0.0039	-0.0019		
Al	-0.00884	107.6	-0.0000	-0.0189	-0.0075		
As	-0.01115	27.38	-0.0111	-0.0080	-0.0142		
B	-0.00070	176.1	-0.0014	-0.0014	0.00072		
Ba	-0.00000	34.09	-0.0000	-0.0000	-0.0000		
Be	0.00001	35.15	0.00001	0.00002	0.00001		
Ca	0.00105	360.5	0.00423	-0.0031	0.00211		
Cd	0.00087	86.64	0.00043	0.00175	0.00043		
Co	-0.00085	100.7	-0.0010	0.00009	-0.0015		
Cr	-0.00310	79.00	-0.0015	-0.0059	-0.0018		
Cu	-0.00000	13841	0.00124	-0.0018	0.00062		
Fe	-0.00176	100.0	0.00000	-0.0035	-0.0017		
K	-0.00000	1009	-0.0256	-0.0256	0.05126		
Li	0.00000	103.1	0.00000	0.00000	0.00000		
Mg	-0.01953	0.080	-0.0195	-0.0195	-0.0195		
Mn	-0.00081	229.2	-0.0004	-0.0028	0.00081		
Mo	0.00259	103.1	0.00566	0.00071	0.00141		
Na	0.04202	40.01	0.04201	0.02520	0.05884		
Ni	0.00031	590.4	0.00215	-0.0015	0.00031		
P	-0.06683	26.01	-0.0863	-0.0531	-0.0609		
Pb	-0.00462	212.9	0.00106	0.00105	-0.0159		
Sb	0.00603	173.2	0.01205	0.01208	-0.0060		
Se	-0.00666	184.9	-0.0008	-0.0208	0.00166		
Si	0.00255	120.1	0.00563	-0.0005	0.00253		
Sn	-0.00342	228.7	0.00341	-0.0017	-0.0119		
Sr	-0.00010	200.4	-0.0001	-0.0003	0.00010		
Ti	-0.00069	249.8	-0.0012	-0.0020	0.00125		
Tl	0.01080	571.3	0.05421	-0.0598	0.03810		
V	-0.00479	34.88	-0.0038	-0.0067	-0.0038		
Zn	0.00031	169.1	0.00092	0.00001	-0.0000		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 9

INSTRUMENT ID: 02360

9 ICSA 12/19/92 22:45

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	-0.03274	3.847	-0.0321	-0.0341	-0.0318		
__Al	475.51550	0.629	474.111	473.482	478.952		
__As	0.20417	7.132	0.20383	0.18978	0.21890		
__B	0.28479	0.756	0.28612	0.28231	0.28596		
__Ba	0.00659	20.80	0.00818	0.00577	0.00583		
__Be	0.00130	0.651	0.00130	0.00130	0.00131		
__Ca	474.51574	0.531	473.605	472.575	477.366		
__Cd	0.02987	2.190	0.02921	0.03052	0.02989		
__Co	0.00199	84.08	0.00142	0.00387	0.00067		
__Cr	-0.01667	14.05	-0.0155	-0.0151	-0.0193		
__Cu	0.00427	50.79	0.00677	0.00303	0.00300		
__Fe	181.12896	0.575	180.633	180.427	182.326		
__K	0.21836	0.036	0.21832	0.21831	0.21845		
__Li	-0.00000	1.732	-0.0000	-0.0000	-0.0000		
__Mg	489.92709	0.620	488.422	487.931	493.427		
__Mn	-0.01281	1.047	-0.0127	-0.0127	-0.0129		
__Mo	-0.04863	1.732	-0.0487	-0.0494	-0.0477		
__Na	0.14296	11.75	0.15977	0.14296	0.12616		
__Ni	-0.02545	8.519	-0.0229	-0.0266	-0.0268		
__P	-1.72598	4.243	-1.6789	-1.8103	-1.6886		
__Pb	0.02710	48.02	0.02285	0.04172	0.01674		
__Sb	-0.00341	610.4	0.01908	-0.0221	-0.0072		
__Se	-0.07799	36.69	-0.0623	-0.1110	-0.0605		
__Si	0.00225	441.9	0.01375	-0.0038	-0.0031		
__Sn	0.08762	13.03	0.07457	0.09242	0.09585		
__Sr	0.01791	1.377	0.01805	0.01763	0.01805		
__Ti	0.01222	3.948	0.01250	0.01166	0.01250		
__Tl	-0.40034	40.44	-0.2145	-0.4750	-0.5114		
__V	0.00427	66.81	0.00713	0.00425	0.00142		
__Zn	-0.00842	17.09	-0.0073	-0.0100	-0.0078		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 10 INSTRUMENT ID: 02360

10 . ICSAB 12/19/92 22:49

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.90156	1.686	0.91844	0.88892	0.89732		
__Al	479.31124	0.919	478.367	475.454	484.112		
__As	0.22422	25.29	0.21272	0.17414	0.28581		
__B	0.28968	0.437	0.29098	0.28959	0.28845		
__Ba	0.48834	0.992	0.48834	0.48350	0.49319		
__Be	0.46052	0.857	0.46052	0.45657	0.46447		
__Ca	476.85308	0.779	476.735	473.195	480.628		
__Cd	0.90317	1.295	0.90710	0.89001	0.91241		
__Co	0.44859	1.088	0.44585	0.44570	0.45423		
__Cr	0.44670	1.415	0.44599	0.44075	0.45334		
__Cu	0.46880	0.667	0.46922	0.46548	0.47170		
__Fe	182.20707	0.819	182.012	180.820	183.788		
__K	0.18030	76.82	0.21873	0.29555	0.02663		
__Li	-0.00000	4.572	-0.0000	-0.0000	-0.0000		
__Mg	493.30447	0.843	492.547	489.576	497.789		
__Mn	0.44139	0.931	0.44183	0.43708	0.44527		
__Mo	-0.04666	4.572	-0.0442	-0.0475	-0.0482		
__Na	0.14822	6.551	0.14261	0.15943	0.14262		
__Ni	0.84522	0.807	0.84373	0.83926	0.85267		
__P	-1.73421	4.069	-1.6527	-1.7747	-1.7751		
__Pb	0.95757	1.551	0.96172	0.94108	0.96992		
__Sb	0.00359	351.8	-0.0094	0.01576	0.00450		
__Se	-0.04829	37.03	-0.0335	-0.0431	-0.0681		
__Si	0.01049	101.0	0.02193	0.00856	0.00099		
__Sn	0.08589	9.752	0.08057	0.09554	0.08154		
__Sr	0.01831	1.346	0.01845	0.01803	0.01845		
__Ti	0.01167	11.48	0.01014	0.01264	0.01223		
__Tl	-0.50778	16.02	-0.4201	-0.5809	-0.5222		
__V	0.47421	0.618	0.47421	0.47128	0.47714		
__Zn	0.90761	0.776	0.90644	0.90121	0.91516		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 11

INSTRUMENT ID: 02360

11 CRI 12/19/92 22:54

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.01818	12.38	0.01558	0.01948	0.01948		
__Al	0.00518	254.7	0.02043	-0.0025	-0.0023		
__As	0.01571	19.43	0.01571	0.01876	0.01265		
__B	-0.00073	169.8	-0.0014	0.00070	-0.0014		
__Ba	0.00080	173.1	0.00000	0.00240	-0.0000		
__Be	0.00893	0.055	0.00893	0.00894	0.00894		
__Ca	0.01198	143.7	0.03173	0.00423	0.00000		
__Cd	0.01139	6.660	0.01095	0.01095	0.01226		
__Co	0.09953	0.984	0.09867	0.09932	0.10060		
__Cr	0.02065	2.991	0.01994	0.02101	0.02100		
__Cu	0.04999	2.501	0.04874	0.04998	0.05124		
__Fe	0.00464	153.6	0.01228	0.00346	-0.0018		
__K	0.15380	76.36	0.05128	0.28196	0.12817		
__Li	0.00101	173.2	0.00303	0.00000	-0.0000		
__Mg	0.00264	864.3	0.02652	0.00046	-0.0190		
__Mn	0.02975	4.739	0.03138	0.02894	0.02894		
__Mo	-0.00048	609.8	-0.0014	0.00282	-0.0028		
__Na	0.06437	15.07	0.05876	0.07557	0.05877		
__Ni	0.07842	2.698	0.08087	0.07721	0.07719		
__P	-0.06688	105.7	-0.1191	-0.0950	0.01360		
__Pb	-0.00035	175.4	0.00001	-0.0000	-0.0010		
__Sb	0.12497	7.937	0.13170	0.12964	0.11358		
__Se	0.00072	1726.	-0.0109	0.01406	-0.0009		
__Si	0.00885	43.58	0.00886	0.01269	0.00498		
__Sn	-0.00683	75.11	-0.0016	-0.0068	-0.0119		
__Sr	-0.00025	96.80	-0.0005	-0.0001	-0.0001		
__Ti	0.00140	123.9	0.00001	0.00084	0.00334		
__Tl	-0.00179	3445.	-0.0454	-0.0289	0.06902		
__V	0.10071	1.645	0.10262	0.09976	0.09974		
__Zn	0.04008	2.695	0.04070	0.04071	0.03883		

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Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 12

INSTRUMENT ID: 02360

12 CCV 12/19/92 22:59

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.47481	0.853	0.47935	0.47351	0.47156		
__Al	24.14152	0.931	24.4006	24.0246	23.9992		
__As	2.10576	0.711	2.12297	2.09868	2.09564		
__B	0.99685	0.919	1.00651	0.99577	0.98827		
__Ba	5.03667	1.151	5.10349	5.00747	4.99906		
__Be	0.48524	1.237	0.49217	0.48176	0.48177		
__Ca	49.63462	1.139	50.2862	49.3458	49.2718		
__Cd	4.89986	0.968	4.95414	4.87930	4.86615		
__Co	0.98982	0.679	0.99756	0.98652	0.98539		
__Cr	5.19383	1.128	5.26087	5.16816	5.15246		
__Cu	4.91104	1.031	4.96935	4.88596	4.87779		
__Fe	4.98237	1.172	5.04984	4.94952	4.94775		
__K	50.48028	0.891	50.9924	50.1473	50.3010		
__Li	25.71655	1.162	26.0590	25.5831	25.5074		
__Mg	24.00026	0.933	24.2589	23.8677	23.8740		
__Mn	4.99658	1.047	5.05652	4.97332	4.95989		
__Mo	24.90167	1.065	25.2064	24.7237	24.7747		
__Na	52.39329	2.086	53.5082	52.3485	51.3230		
__Ni	9.85656	1.204	9.99335	9.78030	9.79602		
__P	25.29309	0.478	25.4095	25.1678	25.3018		
__Pb	25.15381	0.906	25.4171	25.0226	25.0215		
__Sb	5.11438	0.969	5.16442	5.06525	5.11347		
__Se	2.06327	2.102	2.11318	2.04207	2.03457		
__Si	1.93001	0.834	1.94678	1.91468	1.92855		
__Sn	4.97253	0.725	5.01143	4.96594	4.94021		
__Sr	1.03663	1.142	1.05031	1.03012	1.02948		
__Ti	2.05713	1.209	2.08585	2.04266	2.04287		
__Tl	5.36261	2.069	5.48516	5.26896	5.33373		
__V	2.05852	1.381	2.09132	2.04348	2.04074		
__Zn	5.00075	0.778	5.04560	4.97552	4.98113		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 13 INSTRUMENT ID: 02360

13 CCB 12/19/92 23:03

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	-0.00097	100.1	0.00000	-0.0009	-0.0019		
__Al	-0.00126	867.8	-0.0076	0.01139	-0.0075		
__As	-0.01065	8.254	-0.0111	-0.0111	-0.0096		
__B	0.00179	59.68	0.00287	0.00179	0.00072		
__Ba	0.00160	86.60	0.00240	0.00240	-0.0000		
__Be	0.00001	1.341	0.00001	0.00001	0.00001		
__Ca	0.00176	150.9	0.00423	0.00211	-0.0010		
__Cd	0.00218	34.64	0.00306	0.00175	0.00175		
__Co	0.00056	211.9	-0.0008	0.00137	0.00112		
__Cr	-0.00184	81.52	-0.0010	-0.0035	-0.0009		
__Cu	-0.00083	86.37	-0.0000	-0.0012	-0.0012		
__Fe	-0.00117	86.63	0.00000	-0.0017	-0.0017		
__K	0.12816	60.00	0.05126	0.20506	0.12816		
__Li	0.00101	173.2	0.00303	-0.0000	0.00000		
__Mg	-0.00325	172.8	0.00324	-0.0065	-0.0065		
__Mn	-0.00040	0.088	-0.0004	-0.0004	-0.0004		
__Mo	0.00236	91.66	0.00424	-0.0000	0.00283		
__Na	0.04763	20.38	0.04201	0.04203	0.05283		
__Ni	0.00092	115.6	0.00215	0.00030	0.00030		
__P	-0.08612	63.97	-0.0824	-0.1429	-0.0329		
__Pb	0.00159	722.4	-0.0047	0.01492	-0.0053		
__Sb	0.01809	38.58	0.01406	0.02616	0.01406		
__Se	-0.00916	173.2	0.00916	-0.0183	-0.0183		
__Si	0.00536	195.2	0.00718	0.01483	-0.0059		
__Sn	0.00512	208.3	-0.0068	0.00855	0.01367		
__Sr	-0.00024	49.52	-0.0001	-0.0003	-0.0003		
__Ti	0.00027	623.7	-0.0016	0.00166	0.00083		
__Tl	-0.04895	101.5	0.00526	-0.0597	-0.0923		
__V	-0.00383	0.284	-0.0038	-0.0038	-0.0038		
__Zn	0.00156	34.98	0.00093	0.00187	0.00187		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 14 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

14 PBW PBW 12/19/92 23:08

****,100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
✓Ag	-0.00064	174.7	-0.0019	0.00000	0.00000		
✓Al	0.02682	21.84	0.02170	0.03321	0.02555		
As	-0.01721	30.68	-0.0202	-0.0111	-0.0202		
B	0.01001	12.37	0.00929	0.00930	0.01144		
✓Ba	0.00240	0.004	0.00240	0.00240	0.00240		
✓Be	0.00001	1.071	0.00001	0.00001	0.00001		
✓Ca	0.00246	24.74	0.00211	0.00317	0.00211		
Cd	0.00087	173.2	0.00175	-0.0008	0.00175	D	
✓Co	-0.00288	22.48	-0.0022	-0.0028	-0.0035		
✓Cr	-0.00256	82.14	-0.0003	-0.0045	-0.0028		
✓Cu	-0.00104	242.7	-0.0006	-0.0037	0.00125		
✓Fe	0.02265	12.52	0.01942	0.02383	0.02471		
K	0.07690	115.4	0.12816	-0.0256	0.12817		
Li	0.00101	173.2	0.00303	-0.0000	-0.0000		
✓Mg	-0.00869	188.4	0.00649	-0.0260	-0.0065		
✓Mn	-0.00000	36654	-0.0004	0.00081	-0.0004		
Mo	-0.00236	91.55	-0.0000	-0.0042	-0.0028		
✓Na	0.09246	36.35	0.09245	0.12608	0.05884		
✓Ni	-0.00031	680.1	-0.0015	-0.0015	0.00213		
P	-0.07763	48.22	-0.0347	-0.1038	-0.0943		
✓Pb ms 12-21-92(s)	-0.01348	65.78	-0.0063	-0.0234	-0.0106		
✓Sb	0.00134	374.2	0.00600	-0.0040	0.00203		
Se	0.00916	178.9	-0.0058	0.00665	0.02666		
Si	0.12395	0.011	0.12395	0.12395	0.12393		
Sn	0.00513	251.5	-0.0068	0.01881	0.00342		
Sr	-0.00017	183.5	-0.0001	-0.0005	0.00010		
Ti	0.00097	98.99	0.00041	0.00041	0.00208		
Tl	-0.02171	511.1	-0.0597	0.10331	-0.1087		
✓V	-0.00384	0.308	-0.0038	-0.0038	-0.0038		
✓Zn	0.00312	35.07	0.00374	0.00376	0.00185		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 15 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

15 LCSW LCSW 12/19/92 23:12

****,1-1,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
✓Ag	0.04592	4.910	0.04462	0.04461	0.04852		
✓Al	1.92062	0.535	1.91221	1.93209	1.91754		
As	2.09550	0.587	2.08230	2.10665	2.09754		
B	2.08356	0.671	2.06994	2.09790	2.08284		
✓Ba	2.02409	0.707	2.01769	2.04049	2.01409		
✓Be	0.05259	0.028	0.05260	0.05258	0.05260		
✓Ca	4.07592	0.765	4.05124	4.11100	4.06552		
Cd	0.05198	1.458	0.05110	0.05242	0.05242	D	
✓Co	0.51108	1.153	0.50807	0.51787	0.50729		
✓Cr	0.21652	0.929	0.21484	0.21875	0.21597		
✓Cu	0.26088	0.134	0.26068	0.26129	0.26068		
✓Fe	1.04067	0.514	1.03714	1.04683	1.03802		
✓K	4.02045	1.104	4.04608	3.96917	4.04609		
Li	2.03267	0.688	2.02459	2.04883	2.02459		
✓Mg	1.91667	0.856	1.89928	1.93190	1.91882		
✓Mn	0.51894	0.719	0.51568	0.52302	0.51813		
Mo	2.05343	0.430	2.04942	2.06357	2.04730		
✓Na	4.42179	2.878	4.33213	4.36578	4.56746		
✓Ni	0.52462	2.138	0.51484	0.53686	0.52217		
P	2.01438	1.763	2.03143	2.03815	1.97356		
Pb	0.49401	1.619	0.48477	0.49864	0.49863		
✓Sb	0.52137	4.090	0.52466	0.54086	0.49858		
Se	2.15630	0.241	2.16214	2.15463	2.15214		
Si	5.03867	0.725	5.00103	5.07398	5.04100		
Sn	4.07029	0.693	4.05149	4.10275	4.05662		
Sr	1.02030	0.710	1.01770	1.02850	1.01471		
Ti	1.04651	0.535	1.04414	1.05290	1.04248		
Tl	2.11171	1.602	2.10099	2.08452	2.14962		
✓V	0.50661	0.989	0.50371	0.51239	0.50371		
✓Zn	0.52894	0.403	0.52647	0.53015	0.53019		

103

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Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 16 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

16 1908350 JAmnt 12/19/92 23:17

J***,100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00284	79.04	-0.0015	-0.0015	-0.0054		
X Al	1.98903	0.320	1.99634	1.98454	1.98623		
As	-0.00948	103.3	-0.0166	0.00168	-0.0134		
B	0.05021	6.258	0.05383	0.04842	0.04836		
X Ba	0.06888	2.012	0.06728	0.06968	0.06968		
X Be	0.00000	324.8	-0.0000	-0.0000	0.00001		
X Ca	17.63702	0.872	17.4790	17.6456	17.7863		
Cd	0.00201	37.49	0.00158	0.00158	0.00289	D	
X Co	0.00175	84.63	0.00124	0.00059	0.00343		
X Cr	-0.00046	215.0	-0.0000	0.00023	-0.0015		
X Cu	0.00160	118.7	0.00369	-0.0000	0.00119		
Fe	2.22117	0.735	2.20441	2.22206	2.23706		
K	5.92142	1.499	5.97268	5.97269	5.81889		
Li	-0.00101	173.2	-0.0030	-0.0000	0.00000		
X Mg	7.13184	0.760	7.08844	7.11448	7.19260		
X Mn	0.11254	1.251	0.11172	0.11172	0.11416		
Mo	0.00064	127.4	0.00111	-0.0003	0.00111		
X Na	10.43650	1.219	10.3804	10.3468	10.5821		
X Ni	0.00071	644.6	0.00010	0.00560	-0.0035		
P	0.13549	25.81	0.17542	0.11034	0.12070		
Pb	-0.00966	101.0	-0.0181	0.00099	-0.0117		
X Sb	0.00504	23.11	0.00369	0.00574	0.00568		
Se	-0.01827	23.71	-0.0207	-0.0207	-0.0132		
Si	7.66735	0.405	7.63586	7.66813	7.69807		
Sn	0.00492	216.7	-0.0070	0.01347	0.00833		
Sr	0.08932	1.022	0.08847	0.08921	0.09028		
Ti	0.09256	1.351	0.09131	0.09257	0.09381		
Tl	-0.05100	129.0	-0.0891	0.02497	-0.0888		
X V	0.00005	5670.	0.00197	0.00197	-0.0037		
X Zn	0.07486	1.436	0.07423	0.07610	0.07425		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 17

INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

17 1908350 SAMPLE 12/19/92 23:21

UP**,100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.04192	2.684	0.04321	0.04126	0.04127		
__Al	4.11933	0.282	4.11143	4.11384	4.13271		
__As	0.20218	7.970	0.21436	0.20826	0.18390		
__B	0.14747	0.409	0.14784	0.14678	0.14780		
__Ba	0.55858	0.496	0.55698	0.55698	0.56178		
__Be	0.05080	0.007	0.05080	0.05081	0.05080		
__Ca	20.64989	0.324	20.6407	20.5878	20.7211		
__Cd	0.49583	0.305	0.49670	0.49408	0.49670	D	
__Co	0.09827	1.231	0.09950	0.09708	0.09823		
__Cr	0.52252	0.440	0.52329	0.51994	0.52435		
__Cu	0.49607	0.384	0.49440	0.49565	0.49815		
__Fe	2.46555	0.472	2.46320	2.45525	2.47820		
__K	10.22351	0.868	10.2747	10.1209	10.2747		
__Li	2.54286	0.630	2.53073	2.53680	2.56104		
__Mg	8.79067	0.272	8.77222	8.78200	8.81779		
__Mn	0.59740	0.204	0.59618	0.59862	0.59740		
__Mo	2.49351	0.182	2.48832	2.49681	2.49539		
__Na	14.36657	0.722	14.3385	14.2797	14.4814		
__Ni	0.99527	0.296	0.99403	0.99314	0.99863		
__P	2.55089	0.359	2.56101	2.54850	2.54314		
__Pb	2.48866	0.208	2.48990	2.48298	2.49311		
__Sb	0.49804	2.409	0.49437	0.48831	0.51145		
__Se	0.20839	0.599	0.20839	0.20714	0.20964		
__Si	7.07451	0.332	7.09932	7.05251	7.07170		
__Sn	0.48200	1.064	0.48201	0.47688	0.48713		
__Sr	0.18202	0.432	0.18173	0.18141	0.18291		
__Ti	0.28521	1.500	0.28993	0.28159	0.28409		
__Tl	0.42445	39.49	0.60959	0.28297	0.38079		
__V	0.20850	0.683	0.20993	0.20707	0.20852		
__Zn	0.56704	0.576	0.56674	0.56393	0.57045		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 18

INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

18 1908353 5 Amount D 12/19/92 23:26

D***,100=100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00216	52.02	-0.0015	-0.0015	-0.0034		
X Al	2.12175	1.269	2.11207	2.15219	2.10100		
As	-0.00618	113.3	0.00191	-0.0102	-0.0102		
B	0.04740	2.524	0.04602	0.04805	0.04812		
X Ba	0.06969	0.001	0.06969	0.06969	0.06969		
X Be	0.00000	1932.	-0.0000	-0.0000	0.00000		
X Ca	18.09362	0.792	17.9645	18.2480	18.0682		
Cd	0.00069	108.8	0.00026	0.00157	0.00025	D	
X Co	0.00147	89.57	-0.0000	0.00213	0.00233		
X Cr	0.00093	358.2	-0.0024	0.00094	0.00426		
X Cu	0.00348	27.40	0.00244	0.00369	0.00431		
X Fe	2.37239	0.923	2.35445	2.39681	2.36592		
K	6.04959	1.271	5.97269	6.12650	6.04959		
Li	-0.00101	173.2	-0.0000	-0.0030	-0.0000		
X Mg	7.33144	0.889	7.26418	7.39436	7.33579		
X Mn	0.11660	0.002	0.11660	0.11660	0.11660		
Mo	-0.00079	103.4	-0.0003	-0.0017	-0.0003		
X Na	10.60180	1.093	10.5149	10.7334	10.5569		
Ni	0.00253	41.78	0.00192	0.00191	0.00375		
P	0.03331	242.3	0.12608	-0.0213	-0.0048		
Pb	-0.00888	86.72	-0.0000	-0.0127	-0.0138		
X Sb	0.00302	166.6	-0.0023	0.00369	0.00769		
Se	0.00337	561.2	0.01170	0.01670	-0.0182		
Si	7.94088	1.440	7.89250	8.07149	7.85865		
Sn	0.00321	159.5	-0.0019	0.00321	0.00833		
Sr	0.09139	0.994	0.09071	0.09242	0.09103		
Ti	0.09416	2.889	0.09110	0.09632	0.09507		
Tl	-0.02932	361.7	-0.0240	-0.1379	0.07399		
X V	0.00102	162.0	0.00197	0.00197	-0.0008		
X Zn	0.07856	0.675	0.07795	0.07887	0.07887		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 19

INSTRUMENT ID: 02360

Ag Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

19 1908352 5 Amnts 12/19/92 23:31

R***,100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	0.04367	2.586	0.04302	0.04302	0.04498		
X Al	3.54015	0.414	3.52691	3.53762	3.55591		
As	1.96960	0.267	1.96652	1.97569	1.96658		
B	1.96004	0.712	1.94645	1.95934	1.97433		
X Ba	1.96093	0.775	1.94933	1.95533	1.97814		
X Be	0.04871	0.010	0.04872	0.04871	0.04871		
X Ca	19.16196	0.682	19.0524	19.1265	19.3068		
Cd	0.04705	5.576	0.04705	0.04443	0.04968	C.	
X Co	0.48417	1.376	0.48942	0.47667	0.48641		
X Cr	0.20536	1.952	0.20865	0.20089	0.20653		
X Cu	0.24762	0.434	0.24824	0.24637	0.24823		
X Fe	2.84128	0.786	2.82422	2.83305	2.86657		
X K	8.86550	1.001	8.96804	8.81425	8.81422		
Li	1.85688	0.753	1.84880	1.84880	1.87305		
X Mg	8.02625	0.614	7.99152	8.00454	8.08270		
X Mn	0.58181	0.737	0.57773	0.58140	0.58629		
Mo	1.92791	0.430	1.92461	1.92178	1.93735		
X Na	12.94166	1.134	12.7791	12.9808	13.0649		
X Ni	0.49451	0.535	0.49602	0.49145	0.49606		
P	1.93416	3.937	1.99903	1.95314	1.85032		
Pb	0.47564	1.150	0.46960	0.48026	0.47707		
X Sb	0.49670	1.066	0.49068	0.49872	0.50068		
Se	2.04515	0.736	2.04348	2.06098	2.03098		
Si	11.05332	0.583	10.9851	11.0613	11.1134		
Sn	3.84726	0.581	3.84128	3.82847	3.87203		
Sr	1.03829	0.719	1.03266	1.03544	1.04677		
Ti	1.05568	0.570	1.05915	1.04873	1.05915		
Tl	1.96766	1.857	1.92978	2.00271	1.97050		
X V	0.48130	0.351	0.48033	0.48031	0.48325		
X Zn	0.55854	0.510	0.55604	0.55793	0.56164		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 20 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

20 1908350 5Amnt L 12/19/92 23:35

UL**,100-100,DF5,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	-0.00284	34.30	-0.0038	-0.0018	-0.0028		
Al	0.44507	0.844	0.44137	0.44496	0.44889		
As	-0.00432	282.0	-0.0043	-0.0165	0.00788		
B	0.01708	3.633	0.01672	0.01671	0.01779		
Ba	0.01441	0.000	0.01441	0.01441	0.01441		
Be	0.00000	78.05	0.00000	0.00000	0.00001		
Ca	3.51636	0.525	3.49591	3.52130	3.53188		
Cd	0.00084	238.0	0.00303	-0.0009	0.00040	C	
Co	0.00020	599.2	-0.0011	0.00037	0.00134		
Cr	-0.00021	696.6	0.00014	0.00105	-0.0018		
Cu	-0.00001	22349	0.00248	-0.0000	-0.0025		
Fe	0.44359	0.828	0.43946	0.44653	0.44476		
K	1.20479	6.382	1.20478	1.12789	1.28169		
Li	-0.00000	28546	0.00303	-0.0000	-0.0030		
Mg	1.40813	0.706	1.41899	1.39945	1.40596		
Mn	0.02193	3.219	0.02275	0.02153	0.02152		
Mo	-0.00100	214.7	0.00135	-0.0028	-0.0014		
Na	2.14894	0.451	2.14333	2.16015	2.14334		
Ni	-0.00126	233.2	0.00210	-0.0024	-0.0034		
P	-0.03478	146.3	-0.0767	-0.0494	0.02183		
Pb	-0.00119	802.0	-0.0104	0.00874	-0.0019		
Sb	-0.00744	102.3	-0.0040	-0.0161	-0.0020		
Se	0.00083	962.8	0.00666	0.00416	-0.0083		
Si	1.61013	0.856	1.59451	1.62063	1.61525		
Sn	-0.00515	151.9	-0.0119	0.00339	-0.0068		
Sr	0.01755	0.702	0.01763	0.01763	0.01741		
Ti	0.02001	3.606	0.01918	0.02043	0.02043		
Tl	-0.02098	441.4	0.05498	0.00625	-0.1242		
V	-0.00191	87.17	-0.0009	-0.0009	-0.0038		
Zn	0.02420	0.077	0.02418	0.02421	0.02422		

108

AR318927

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 21 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

21 1908351 YAmmT 12/19/92 23:40
****,100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00126	88.81	-0.0019	-0.0019	0.00003		
X Al	0.02135	36.26	0.01343	0.02890	0.02170		
As	-0.00470	171.3	-0.0138	0.00138	-0.0016		
B	0.05044	2.460	0.05113	0.04900	0.05118		
X Ba	0.07322	1.639	0.07442	0.07322	0.07202		
X Be	0.00000	57.54	0.00001	0.00000	0.00001		
X Ca	18.28296	0.989	18.4548	18.2998	18.0941		
Cd	0.00035	367.8	-0.0009	0.00035	0.00167	D	
X Co	-0.00069	344.8	-0.0024	0.00203	-0.0017		
X Cr	-0.00305	35.92	-0.0031	-0.0041	-0.0019		
X Cu	0.00145	194.1	-0.0012	0.00124	0.00437		
X Fe	0.21662	1.627	0.22015	0.21662	0.21310		
X K	6.07524	0.730	6.12651	6.04961	6.04961		
Li	-0.00101	173.2	-0.0030	-0.0000	-0.0000		
X Mg	7.27630	0.607	7.32296	7.27088	7.23506		
X Mn	0.08492	2.876	0.08737	0.08492	0.08248		
Mo	0.00045	474.3	0.00281	-0.0000	-0.0014		
X Na	11.11404	0.698	11.2036	11.0692	11.0692		
X Ni	-0.00215	49.29	-0.0015	-0.0033	-0.0015		
P	-0.03611	67.27	-0.0081	-0.0521	-0.0479		
Pb	-0.00265	265.5	0.00533	-0.0053	-0.0079		
X Sb	0.01472	31.43	0.02007	0.01206	0.01204		
Se	-0.00528	263.4	-0.0202	0.00721	-0.0027		
Si	5.05740	0.537	5.08200	5.06202	5.02818		
Sn	0.00409	180.6	0.00836	0.00837	-0.0044		
Sr	0.09149	0.884	0.09242	0.09114	0.09092		
Ti	0.00041	199.5	0.00041	-0.0004	0.00125		
Tl	0.01064	88.29	0.02149	0.00520	0.00523		
X V	-0.00285	57.95	-0.0038	-0.0009	-0.0038		
X Zn	0.04835	0.022	0.04836	0.04835	0.04834		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 22 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni Li K Mo Na V Zn

22 1908354 TAMFW 12/19/92 23:44

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00225	66.04	-0.0019	-0.0009	-0.0038		
X Al	0.00628	65.10	0.00587	0.00241	0.01056		
As	-0.01040	59.19	-0.0032	-0.0139	-0.0139		
B	0.04931	0.070	0.04927	0.04931	0.04934		
X Ba	0.04961	2.793	0.05041	0.05041	0.04801		
X Be	0.00000	104.4	0.00000	0.00000	0.00001		
X Ca	16.96410	0.590	17.0596	16.9729	16.8597		
Cd	-0.00050	149.0	-0.0009	-0.0009	0.00036	C	
X Co	0.00125	10.42	0.00138	0.00112	0.00125		
X Cr	-0.00235	123.4	0.00068	-0.0026	-0.0051		
X Cu	0.00124	100.2	0.00249	0.00124	-0.0000		
X Fe	0.15311	1.148	0.15486	0.15311	0.15135		
X K	5.58820	1.376	5.51130	5.58821	5.66510		
Li	-0.00000	99.22	-0.0000	-0.0000	-0.0000		
X Mg	6.75233	1.162	6.83478	6.74365	6.67856		
X Mn	0.08046	0.874	0.08127	0.08005	0.08005		
Mo	-0.00142	99.22	-0.0000	-0.0028	-0.0014		
X Na	10.17829	0.165	10.1951	10.1782	10.1614		
X Ni	-0.00032	1174.	-0.0033	-0.0015	0.00395		
P	-0.04148	127.9	-0.0454	-0.0924	0.01345		
Pb	0.00160	734.0	0.00053	0.01385	-0.0095		
X Sb	0.00806	74.57	0.00801	0.01409	0.00207		
Se	0.00051	2668.	-0.0128	-0.0003	0.01467		
Si	4.68676	0.732	4.71930	4.69010	4.65088		
Sn	-0.01041	75.20	-0.0018	-0.0121	-0.0172		
Sr	0.08444	1.022	0.08537	0.08430	0.08366		
Ti	0.00027	311.0	0.00000	-0.0004	0.00125		
Tl	0.00246	503.0	-0.0110	0.01334	0.00509		
X V	-0.00238	104.5	-0.0009	-0.0009	-0.0052		
X Zn	0.04401	2.459	0.04463	0.04463	0.04276		

110

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 23 INSTRUMENT ID: 02360

A ISb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00195	0.059	-0.0019	-0.0019	-0.0019		
X Al	0.02367	32.27	0.03131	0.02367	0.01603		
As	-0.01421	42.97	-0.0203	-0.0142	-0.0081		
B	0.01218	10.17	0.01147	0.01362	0.01147		
X Ba	0.00240	0.002	0.00240	0.00240	0.00240		
X Be	0.00000	56.51	0.00000	0.00001	0.00001		
X Ca	0.15795	1.393	0.15866	0.15972	0.15549		
Cd	0.00043	599.1	0.00306	0.00043	-0.0021	D	
X Co	0.00156	25.76	0.00111	0.00188	0.00169		
X Cr	-0.00153	85.56	-0.0029	-0.0003	-0.0013		
X Cu	-0.00270	58.05	-0.0012	-0.0024	-0.0043		
X Fe	0.00000	25089	-0.0017	0.00176	0.00000		
X K	0.17943	24.74	0.20506	0.20506	0.12816		
Li	-0.00101	173.2	-0.0000	-0.0000	-0.0030		
X Mg	0.01301	0.007	0.01301	0.01301	0.01301		
X Mn	-0.00040	0.123	-0.0004	-0.0004	-0.0004		
Mo	-0.00094	86.41	-0.0000	-0.0014	-0.0014		
X Na	0.36141	9.303	0.39503	0.36141	0.32778		
X Ni	-0.00184	115.1	-0.0006	-0.0042	-0.0006		
P	-0.08179	90.55	-0.1672	-0.0359	-0.0421		
Pb	-0.00425	156.6	0.00321	-0.0095	-0.0063		
X Sb	0.00000	81925	0.00202	-0.0000	-0.0019		
Se	0.00166	1653.	0.02916	-0.0258	0.00166		
Si	0.66085	0.673	0.65571	0.66341	0.66342		
Sn	0.00000	23646	-0.0017	0.01624	-0.0145		
Sr	0.00078	20.81	0.00096	0.00074	0.00064		
Ti	0.00041	360.6	0.00208	-0.0000	-0.0008		
Tl	-0.00810	252.0	-0.0270	-0.0107	0.01353		
X V	-0.00288	57.58	-0.0009	-0.0038	-0.0038		
X Zn	0.00188	0.471	0.00187	0.00189	0.00189		

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Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 24

INSTRUMENT ID: 02360

24 CCV 12/19/92 23:53

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.46669	1.269	0.47058	0.46961	0.45987		
__Al	23.75483	1.323	24.0136	23.8460	23.4048		
__As	2.05502	2.160	2.08650	2.07432	2.00424		
__B	0.97538	1.588	0.98830	0.97964	0.95820		
__Ba	4.93305	1.537	4.99906	4.94985	4.85023		
__Be	0.47569	1.559	0.48178	0.47787	0.46743		
__Ca	48.65549	1.363	49.1998	48.8497	47.9168		
__Cd	4.82938	1.286	4.87928	4.84908	4.75979		
__Co	0.96619	1.610	0.96953	0.97981	0.94924		
__Cr	5.09544	1.415	5.15176	5.12041	5.01414		
__Cu	4.81527	1.446	4.87526	4.83166	4.73890		
__Fe	4.88320	1.451	4.94156	4.90372	4.80431		
__K	49.32746	1.126	49.6857	49.6092	48.6873		
__Li	25.44579	1.476	25.8014	25.4831	25.0527		
__Mg	23.66529	1.463	23.9000	23.8282	23.2675		
__Mn	4.90566	1.301	4.95623	4.92685	4.83388		
__Mo	24.50979	1.497	24.8228	24.6006	24.1058		
__Na	49.55537	1.686	50.2977	49.7178	48.6505		
__Ni	9.68654	1.376	9.78883	9.73505	9.53575		
__P	25.00031	1.560	25.2629	25.1861	24.5519		
__Pb	24.98950	1.241	25.2379	25.0886	24.6419		
__Sb	5.00971	1.089	5.04274	5.03966	4.94673		
__Se	2.03343	1.605	2.01833	2.07090	2.01105		
__Si	1.96752	0.810	1.97993	1.97309	1.94953		
__Sn	4.88842	1.102	4.93243	4.90453	4.82828		
__Sr	1.01032	1.611	1.02467	1.01366	0.99262		
__Ti	2.03118	1.608	2.05705	2.04203	1.99447		
__Tl	5.21679	0.602	5.23685	5.18060	5.23293		
__V	2.01575	1.315	2.03802	2.02282	1.98642		
__Zn	4.92499	1.134	4.96806	4.94503	4.86187		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 25

INSTRUMENT ID: 02360

25 CCB 12/19/92 23:58

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	-0.00130	86.64	-0.0019	0.00000	-0.0019		
__Al	0.15662	6.425	0.16029	0.14524	0.16433		
__As	-0.00710	98.85	0.00100	-0.0111	-0.0111		
__B	-0.00035	305.0	0.00072	-0.0014	-0.0003		
__Ba	0.00160	86.60	0.00240	-0.0000	0.00240		
__Be	0.00000	175.7	-0.0000	0.00001	0.00001		
__Ca	0.00458	35.25	0.00634	0.00423	0.00317		
__Cd	-0.00021	532.7	0.00044	0.00044	-0.0015		
__Co	0.00064	149.1	0.00176	0.00008	0.00009		
__Cr	-0.00145	143.9	0.00091	-0.0030	-0.0022		
__Cu	0.00041	571.5	0.00312	-0.0012	-0.0006		
__Fe	-0.00206	89.06	-0.0000	-0.0035	-0.0026		
__K	0.10252	43.30	0.05125	0.12816	0.12815		
__Li	0.00202	86.60	0.00303	0.00303	0.00000		
__Mg	-0.00434	173.1	0.00000	-0.0130	-0.0000		
__Mn	-0.00040	0.205	-0.0004	-0.0004	-0.0004		
__Mo	0.00400	62.07	0.00636	0.00140	0.00423		
__Na	0.01401	138.5	0.02521	-0.0084	0.02522		
__Ni	0.00092	413.7	0.00398	-0.0033	0.00214		
__P	-0.03752	78.68	-0.0532	-0.0558	-0.0034		
__Pb	-0.00382	167.7	0.00221	-0.0031	-0.0105		
__Sb	0.00468	321.6	-0.0040	0.02210	-0.0040		
__Se	-0.00336	128.6	-0.0058	0.00163	-0.0058		
__Si	0.00636	55.35	0.01020	0.00560	0.00328		
__Sn	-0.00167	529.9	-0.0068	-0.0068	0.00858		
__Sr	-0.00003	918.7	0.00032	-0.0003	-0.0001		
__Ti	0.00111	120.5	0.00208	-0.0004	0.00166		
__Tl	-0.00280	2037.	0.06241	-0.0272	-0.0435		
__V	-0.00190	174.6	0.00193	-0.0038	-0.0038		
__Zn	0.01308	24.61	0.01681	0.01123	0.01121		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 26 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

26 1908356 JAWLT 12/20/92 00:02
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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00291	38.51	-0.0035	-0.0016	-0.0035		
X Al	1.39524	0.658	1.40107	1.40000	1.38465		
As	-0.00403	265.4	0.00716	-0.0050	-0.0141		
B	0.03605	6.154	0.03385	0.03828	0.03603		
X Ba	0.06007	0.001	0.06007	0.06007	0.06007		
X Be	-0.00000	111.5	-0.0000	0.00000	-0.0000		
X Ca	17.76254	0.994	17.9127	17.5679	17.8069		
Cd	-0.00059	126.9	-0.0010	-0.0010	0.00027	D	
X Co	-0.00163	55.94	-0.0006	-0.0018	-0.0024		
X Cr	-0.00055	511.1	0.00196	0.00000	-0.0036		
X Cu	0.00204	93.60	0.00370	0.00245	-0.0000		
X Fe	1.83486	1.123	1.85310	1.81251	1.83898		
X K	5.39595	0.712	5.43440	5.39594	5.35750		
Li	-0.00101	173.2	-0.0030	0.00000	-0.0000		
X Mg	7.09392	1.353	7.16335	6.98435	7.13405		
X Mn	0.10889	0.644	0.10929	0.10808	0.10930		
Mo	-0.00188	108.5	-0.0030	0.00047	-0.0030		
X Na	9.65750	1.058	9.70793	9.53983	9.72475		
X Ni	0.00075	141.5	0.00013	0.00198	0.00013		
P	0.08581	28.51	0.11402	0.07017	0.07326		
Pb	-0.00907	58.83	-0.0142	-0.0035	-0.0094		
X Sb	-0.00627	84.63	-0.0002	-0.0103	-0.0082		
Se	-0.00022	7003.	0.00560	0.01184	-0.0181		
Si	6.96614	0.621	7.01610	6.94002	6.94231		
Sn	-0.00793	130.6	-0.0173	-0.0096	0.00317		
Sr	0.08729	1.605	0.08857	0.08579	0.08750		
Ti	0.06372	2.178	0.06254	0.06338	0.06525		
Tl	-0.13642	43.27	-0.2044	-0.1065	-0.0982		
X V	0.00148	55.42	0.00196	0.00053	0.00195		
X Zn	0.07929	1.352	0.07991	0.07992	0.07805		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 27 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag V Zn

27 1908357 JAMF 12/20/92 00:07

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00127	88.11	-0.0019	0.00002	-0.0019		
X Al	0.01747	76.06	0.02297	0.00231	0.02713		
As	0.00333	190.1	0.01045	-0.0017	0.00129		
B	0.08373	0.027	0.08370	0.08374	0.08374		
X Ba	0.04081	0.000	0.04081	0.04081	0.04081		
X Be	0.00000	87.62	0.00000	0.00001	0.00000		
X Ca	17.13140	0.457	17.2167	17.1146	17.0628		
Cd	0.00189	20.00	0.00167	0.00167	0.00233	D	
X Co	0.00224	63.01	0.00306	0.00061	0.00306		
X Cr	-0.00211	162.7	-0.0001	-0.0061	-0.0001		
X Cu	0.00499	25.00	0.00624	0.00374	0.00499		
X Fe	0.16546	1.845	0.16899	0.16370	0.16370		
X K	5.12681	1.499	5.04991	5.20370	5.12680		
Li	0.00101	346.4	-0.0030	0.00303	0.00303		
X Mg	6.75127	0.665	6.80227	6.73390	6.71764		
X Mn	0.08331	1.692	0.08494	0.08250	0.08250		
Mo	-0.00048	168.3	-0.0000	-0.0014	-0.0000		
X Na	9.44988	0.447	9.48910	9.45548	9.40505		
X Ni	0.00089	580.8	0.00670	-0.0033	-0.0006		
P	0.00880	185.8	0.02532	0.00845	-0.0073		
Pb	-0.01064	86.71	-0.0159	-0.0159	0.00001		
X Sb	0.01476	34.28	0.01412	0.02011	0.01005		
Se	-0.00323	726.1	0.00842	-0.0303	0.01217		
Si	4.89166	0.326	4.90985	4.88525	4.87987		
Sn	0.00326	157.3	0.00326	-0.0018	0.00839		
Sr	0.08263	0.583	0.08312	0.08259	0.08216		
Ti	0.00153	15.66	0.00125	0.00166	0.00166		
Tl	0.02696	303.8	0.10288	-0.0598	0.03788		
X V	-0.00190	87.75	-0.0009	-0.0038	-0.0009		
X Zn	0.04087	0.054	0.04084	0.04088	0.04087		

135

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Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 28 INSTRUMENT ID: 02360

Al Sb Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

28 1908358 CHUIX 12/20/92 00:12

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00214	138.9	-0.0053	-0.0014	0.00045		
X Al	1.68506	0.458	1.67615	1.68910	1.68995		
As	-0.00701	129.9	-0.0161	-0.0070	0.00210		
B	0.11946	2.060	0.11801	0.12230	0.11806		
X Ba	0.05289	0.001	0.05289	0.05289	0.05289		
X Be	0.00000	222.7	0.00000	-0.0000	0.00000		
X Ca	10.30687	0.546	10.3206	10.3550	10.2449		
Cd	0.00026	491.2	0.00026	0.00157	-0.0010	D	
X Co	0.00140	250.6	-0.0011	-0.0000	0.00541		
X Cr	0.00050	151.2	-0.0002	0.00050	0.00127		
X Cu	0.00078	244.5	-0.0013	0.00244	0.00119		
X Fe	2.47669	0.689	2.48699	2.48611	2.45698		
X K	3.02482	1.467	2.97355	3.05045	3.05045		
Li	0.00101	173.2	-0.0000	0.00303	-0.0000		
X Mg	4.50476	0.661	4.49825	4.53730	4.47874		
X Mn	0.15375	0.458	0.15456	0.15334	0.15334		
Mo	-0.00218	99.18	-0.0045	-0.0017	-0.0002		
X Na	9.10840	0.910	9.06919	9.20368	9.05235		
X Ni	0.00374	84.82	0.00007	0.00558	0.00558		
P	0.03796	144.1	-0.0247	0.06292	0.07575		
Pb	-0.00661	116.1	-0.0151	-0.0002	-0.0044		
X Sb	0.00440	115.1	0.00370	-0.0002	0.00978		
Se	-0.00803	85.61	-0.0084	-0.0147	-0.0009		
Si	7.07873	0.512	7.08283	7.11276	7.04061		
Sn	-0.00110	583.8	-0.0019	0.00572	-0.0070		
Sr	0.06179	1.111	0.06207	0.06229	0.06101		
Ti	0.06727	6.827	0.06755	0.07172	0.06254		
Tl	-0.00325	2512.	-0.0901	0.00780	0.07253		
X V	0.00007	2244.	-0.0008	0.00199	-0.0008		
X Zn	0.03346	1.574	0.03316	0.03407	0.03314		

136

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 29

INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

29 1908359 CH WLF 12/20/92 00:16
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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00222	66.89	-0.0038	-0.0009	-0.0019		
X Al	0.04276	13.52	0.04402	0.04782	0.03645		
As	-0.00148	356.6	0.00461	-0.0045	-0.0045		
B	0.03769	5.931	0.03518	0.03948	0.03840		
X Ba	0.03121	0.000	0.03121	0.03121	0.03121		
X Be	0.00000	0.890	0.00000	0.00000	0.00000		
X Ca	9.21279	0.261	9.19833	9.24064	9.19939		
Cd	-0.00005	1311.	-0.0009	0.00038	0.00038	D	
X Co	0.00120	256.7	0.00434	0.00111	-0.0018		
X Cr	-0.00104	200.9	-0.0010	0.00104	-0.0031		
X Cu	-0.00063	98.74	-0.0012	-0.0000	-0.0006		
X Fe	0.32507	0.313	0.32625	0.32448	0.32448		
X K	2.71722	1.633	2.74285	2.74285	2.66595		
Li	-0.00101	173.2	-0.0000	-0.0030	-0.0000		
X Mg	3.95383	0.342	3.93865	3.95817	3.96468		
X Mn	0.11598	0.608	0.11557	0.11557	0.11679		
Mo	-0.00096	84.27	-0.0000	-0.0014	-0.0014		
X Na	8.16114	0.000	8.16114	8.16114	8.16113		
X Ni	0.00180	77.70	0.00302	0.00027	0.00211		
P	-0.06979	75.31	-0.1222	-0.0171	-0.0699		
Pb	-0.01170	32.81	-0.0106	-0.0159	-0.0085		
X Sb	0.00067	685.9	-0.0019	-0.0020	0.00604		
Se	0.00068	1143.	-0.0080	0.00693	0.00318		
Si	4.05889	0.115	4.05992	4.06298	4.05378		
Sn	-0.00098	988.9	0.00585	-0.0120	0.00329		
Sr	0.05420	0.113	0.05427	0.05417	0.05417		
Ti	0.00306	28.35	0.00333	0.00375	0.00208		
Tl	-0.06531	112.7	-0.0598	0.00544	-0.1414		
X V	-0.00093	0.537	-0.0009	-0.0009	-0.0009		
X Zn	0.01602	6.742	0.01477	0.01665	0.01665		

AR318936

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 30

INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

30 1908981 JSBHT 12/20/92 00:21

****.100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00349	55.86	-0.0015	-0.0054	-0.0034		
X Al	2.09601	0.623	2.10742	2.08176	2.09886		
As	-0.00433	122.0	-0.0012	-0.0012	-0.0104		
B	0.04483	2.732	0.04411	0.04625	0.04414		
X Ba	0.06088	1.138	0.06128	0.06128	0.06008		
X Be	0.00000	134.6	-0.0000	0.00000	0.00000		
X Ca	17.76747	0.537	17.8180	17.8270	17.6572		
Cd	-0.00061	124.3	-0.0010	-0.0010	0.00026	D	
X Co	0.00157	162.2	-0.0013	0.00288	0.00321		
X Cr	0.00130	28.02	0.00154	0.00148	0.00088		
X Cu	0.00452	15.99	0.00368	0.00494	0.00494		
X Fe	2.25768	0.453	2.26709	2.25915	2.24679		
X K	5.26778	2.107	5.20369	5.39595	5.20369		
Li	0.00101	173.2	-0.0000	-0.0000	0.00303		
X Mg	6.98422	0.493	6.97120	7.02327	6.95819		
X Mn	0.10235	0.689	0.10317	0.10194	0.10195		
Mo	-0.00432	77.27	-0.0031	-0.0080	-0.0017		
X Na	9.71374	0.780	9.79221	9.70812	9.64090		
X Ni	-0.00052	204.3	0.00009	-0.0017	0.00009		
P	0.11803	36.64	0.16787	0.09583	0.09038		
Pb	-0.00570	224.6	0.00744	-0.0064	-0.0181		
X Sb	0.00335	105.8	-0.0003	0.00671	0.00368		
Se	-0.00580	310.6	-0.0008	0.00919	-0.0258		
Si	7.73612	0.673	7.79627	7.70414	7.70794		
Sn	-0.01215	0.046	-0.0121	-0.0121	-0.0121		
Sr	0.08857	0.638	0.08900	0.08879	0.08793		
Ti	0.10466	4.382	0.10924	0.10007	0.10466		
Tl	-0.08842	66.65	-0.0718	-0.1538	-0.0395		
X V	0.00004	3996.	0.00196	-0.0009	-0.0009		
X Zn	0.07797	0.001	0.07797	0.07797	0.07797		

128

AR318937

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 31 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

31 1908982 JSBFL 12/20/92 00:25
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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00128	175.6	0.00001	0.00001	-0.0038		
X Al	-0.00354	103.0	-0.0033	0.00002	-0.0072		
As	-0.01093	100.4	-0.0231	-0.0017	-0.0078		
B	0.03501	3.536	0.03429	0.03644	0.03430		
X Ba	0.04081	0.000	0.04081	0.04081	0.04081		
X Be	0.00000	55.06	0.00000	0.00000	0.00000		
X Ca	17.36940	0.503	17.3489	17.4653	17.2939		
Cd	0.00058	323.2	0.00168	-0.0016	0.00168		
X Co	-0.00106	198.9	-0.0022	-0.0023	0.00138		
X Cr	-0.00403	27.02	-0.0052	-0.0037	-0.0031		
X Cu	0.00083	114.8	0.00062	0.00187	-0.0000		
X Fe	0.13841	0.737	0.13900	0.13723	0.13900		
X K	5.15244	1.723	5.20371	5.20371	5.04991		
Li	0.00101	173.2	0.00303	-0.0000	-0.0000		
X Mg	6.68280	0.313	6.69147	6.69798	6.65894		
X Mn	0.06783	0.000	0.06783	0.06783	0.06783		
Mo	-0.00284	49.82	-0.0014	-0.0028	-0.0042		
X Na	9.64039	0.628	9.62357	9.70763	9.58995		
X Ni	-0.00154	119.0	0.00029	-0.0015	-0.0033		
P	-0.02702	121.1	-0.0646	-0.0119	-0.0045		
Pb	-0.01527	76.83	-0.0074	-0.0287	-0.0095		
X Sb	0.00839	100.3	0.00405	0.00302	0.01810		
Se	0.00468	423.5	-0.0178	0.01218	0.01968		
Si	4.89603	0.431	4.87759	4.91906	4.89143		
Sn	0.00155	761.0	0.00839	0.00839	-0.0121		
Sr	0.08569	0.623	0.08569	0.08622	0.08515		
Ti	0.00111	114.4	0.00083	0.00250	0.00000		
Tl	-0.06811	141.2	-0.1741	-0.0435	0.01341		
X V	-0.00144	57.77	-0.0009	-0.0009	-0.0024		
X Zn	0.03466	1.546	0.03528	0.03434	0.03436		

AR318938

Run Name: 9235402111

LANCASTER LABORATORIES

PAGE: 32 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00288	39.07	-0.0035	-0.0015	-0.0035		
X Al	1.47778	0.241	1.47814	1.48116	1.47404		
As	-0.00476	64.02	-0.0047	-0.0078	-0.0017		
B	0.03196	3.857	0.03054	0.03268	0.03267		
X Ba	0.04647	2.982	0.04807	0.04567	0.04567		
X Be	0.00000	1.399	0.00000	0.00000	0.00000		
X Ca	10.18752	0.317	10.1683	10.1693	10.2248		
Cd	-0.00014	539.2	-0.0010	0.00029	0.00029		
X Co	0.00161	229.8	-0.0021	0.00168	0.00528		
X Cr	-0.00009	778.3	0.00031	0.00032	-0.0009		
X Cu	0.00599	26.18	0.00433	0.00745	0.00620		
Fe	1.99715	0.192	1.99450	1.99538	2.00156		
K	2.92228	3.038	2.81975	2.97354	2.97355		
Li	-0.00000	170.5	-0.0000	0.00000	-0.0000		
X Mg	4.33749	0.439	4.31579	4.35159	4.34509		
X Mn	0.11507	0.612	0.11548	0.11426	0.11548		
Mo	-0.00166	170.5	-0.0044	0.00117	-0.0016		
X Na	10.59601	0.391	10.6156	10.5483	10.6240		
X Ni	0.00256	41.49	0.00195	0.00379	0.00194		
P	0.00754	547.4	-0.0343	0.00884	0.04819		
Pb	-0.00885	52.48	-0.0120	-0.0035	-0.0109		
X Sb	-0.00191	199.3	-0.0062	0.00076	-0.0002		
Sc	0.00326	563.6	-0.0108	0.02410	-0.0033		
Si	6.15160	0.584	6.12010	6.19078	6.14392		
Sn	0.00066	384.3	-0.0018	0.00322	0.00066		
Sr	0.06336	0.337	0.06336	0.06314	0.06357		
Ti	0.05684	1.120	0.05670	0.05754	0.05629		
Tl	0.01265	74.31	0.00723	0.00722	0.02351		
X V	-0.00090	1.041	-0.0009	-0.0008	-0.0009		
X Zn	0.03008	3.592	0.02946	0.02944	0.03132		

100

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 33 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

33 1908984 CRCHK 12/20/92 00:35

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	-0.00385	0.010	-0.0038	-0.0038	-0.0038		
Al	0.04697	11.33	0.04113	0.04820	0.05157		
As	-0.00046	1007.	0.00462	-0.0045	-0.0014		
B	0.02752	10.29	0.02431	0.02861	0.02965		
Ba	0.03361	0.000	0.03361	0.03361	0.03361		
Be	0.00001	34.66	0.00002	0.00001	0.00001		
Ca	10.21079	1.050	10.0984	10.2217	10.3121		
Cd	-0.00049	152.3	-0.0009	-0.0009	0.00037	D	
Co	0.00000	5034.	0.00008	0.00022	-0.0002		
Cr	-0.00355	30.91	-0.0046	-0.0024	-0.0035		
Cu	0.00082	115.4	0.00061	0.00186	-0.0000		
Fe	0.31388	1.123	0.31035	0.31388	0.31741		
K	2.71722	1.633	2.74286	2.74285	2.66596		
Li	-0.00101	173.2	-0.0000	-0.0030	-0.0000		
Mg	4.28569	0.749	4.27049	4.26399	4.32258		
Mn	0.10537	1.338	0.10456	0.10456	0.10700		
Mo	-0.00049	329.0	-0.0014	0.00139	-0.0014		
Na	10.54812	1.304	10.3968	10.5817	10.6657		
Ni	-0.00002	6496.	0.00210	-0.0015	-0.0006		
P	-0.03520	107.2	-0.0686	0.00576	-0.0427		
Pb	-0.00388	192.6	0.00321	-0.0031	-0.0116		
Sb	-0.00000	13645	0.00002	-0.0020	0.00201		
Se	0.01029	138.0	0.02196	-0.0055	0.01446		
Si	4.38727	1.024	4.34193	4.38803	4.43185		
Sn	0.00243	160.8	0.00328	0.00584	-0.0018		
Sr	0.06293	1.224	0.06207	0.06314	0.06357		
Ti	0.00236	36.69	0.00208	0.00333	0.00166		
Tl	-0.07621	36.94	-0.0599	-0.1087	-0.0599		
V	-0.00477	34.87	-0.0066	-0.0038	-0.0038		
Zn	0.01632	3.291	0.01663	0.01663	0.01570		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 34 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

34 1908985 C RTFB 12/20/92 00:39

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00130	86.57	-0.0000	-0.0019	-0.0019		
X Al	0.01905	41.22	0.02528	0.02165	0.01023		
As	-0.00151	379.3	0.00100	0.00254	-0.0081		
B	0.01291	9.597	0.01362	0.01363	0.01148		
X Ba	0.00240	0.001	0.00240	0.00240	0.00240		
X Be	0.00000	301.3	-0.0000	0.00000	0.00001		
X Ca	0.17065	6.451	0.18299	0.16183	0.16712		
Cd	0.00131	57.66	0.00175	0.00175	0.00043	Δ	
X Co	-0.00021	895.2	0.00124	0.00047	-0.0023		
X Cr	-0.00154	105.6	0.00008	-0.0015	-0.0031		
X Cu	-0.00083	172.9	-0.0000	0.00000	-0.0024		
X Fe	0.00029	620.6	0.00176	0.00088	-0.0017		
X K	0.20506	0.001	0.20505	0.20506	0.20506		
Li	0.00202	173.2	0.00000	0.00606	-0.0000		
X Mg	0.00432	313.3	0.01952	-0.0065	-0.0000		
X Mn	-0.00122	57.75	-0.0004	-0.0016	-0.0016		
Mo	0.00047	458.5	0.00282	-0.0000	-0.0014		
X Na	0.41743	2.325	0.42864	0.41183	0.41182		
X Ni	-0.00244	112.7	0.00031	-0.0024	-0.0051		
P	-0.02439	224.7	0.03692	-0.0687	-0.0413		
Pb	-0.00176	211.9	0.00214	-0.0053	-0.0021		
X Sb	0.00535	78.06	0.00401	0.00200	0.01003		
Se	0.00457	177.5	0.01291	0.00416	-0.0033		
Si	0.85246	0.540	0.85706	0.84784	0.85247		
Sn	-0.01025	76.37	-0.0017	-0.0170	-0.0119		
Sr	0.00096	22.20	0.00117	0.00096	0.00074		
Ti	-0.00076	253.3	0.00083	-0.0002	-0.0029		
Tl	0.04893	226.8	0.13593	-0.0760	0.08694		
X V	-0.00096	300.5	0.00192	-0.0009	-0.0038		
X Zn	0.00375	0.402	0.00373	0.00374	0.00376		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 35 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

35 1908986 5 SBMT 12/20/92 00:44
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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00212	53.01	-0.0034	-0.0014	-0.0014		
X Al	1.49836	1.435	1.51074	1.51082	1.47352		
As	-0.00071	735.1	0.00233	-0.0067	0.00229		
B	0.03072	0.166	0.03070	0.03068	0.03078		
X Ba	0.06089	2.276	0.06009	0.06249	0.06009		
X Be	-0.00000	4.424	-0.0000	-0.0000	-0.0000		
X Ca	17.75707	0.668	17.8027	17.8461	17.6223		
Cd	0.00109	69.21	0.00153	0.00153	0.00021	Δ	
X Co	-0.00114	164.5	0.00096	-0.0017	-0.0026		
X Cr	0.00007	1885.	0.00055	-0.0014	0.00116		
X Cu	0.00327	29.21	0.00431	0.00243	0.00306		
X Fe	2.60573	0.785	2.60867	2.62456	2.58396		
X K	6.04959	1.271	5.97269	6.12649	6.04958		
Li	0.00202	86.60	0.00303	-0.0000	0.00303		
X Mg	6.97566	0.782	7.00388	7.01036	6.91273		
X Mn	0.12637	0.965	0.12760	0.12637	0.12516		
Mo	-0.00266	30.78	-0.0031	-0.0031	-0.0017		
X Na	9.23745	0.945	9.18701	9.33831	9.18701		
X Ni	0.00160	87.55	0.00190	0.00007	0.00283		
P	0.10887	30.68	0.14132	0.11072	0.07458		
Pb	-0.01488	29.83	-0.0184	-0.0099	-0.0163		
X Sb	-0.00137	935.4	0.00165	0.00968	-0.0154		
Se	-0.00361	369.5	-0.0107	0.01180	-0.0119		
Si	6.54029	0.788	6.57386	6.56615	6.48088		
Sn	-0.00725	70.77	-0.0021	-0.0072	-0.0123		
Sr	0.08718	0.765	0.08697	0.08793	0.08665		
Ti	0.10118	1.559	0.10007	0.10299	0.10049		
Tl	-0.10730	96.14	0.00949	-0.1859	-0.1454		
X V	0.00197	0.251	0.00197	0.00197	0.00198		
X Zn	0.08980	1.196	0.09104	0.08918	0.08918		

AR318942

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 36 INSTRUMENT ID: 02360

36 CCV . 12/20/92 00:49

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.46376	0.727	0.46570	0.46571	0.45986		
Al	23.82816	0.630	23.9607	23.8585	23.6651		
As	2.05655	0.740	2.06212	2.06822	2.03931		
B	0.97392	0.711	0.98182	0.97107	0.96887		
Ba	4.94625	0.718	4.97986	4.94985	4.90904		
Be	0.47744	0.686	0.48048	0.47786	0.47397		
Ca	48.78312	0.543	48.9872	48.8788	48.4832		
Cd	4.84404	0.574	4.86417	4.85564	4.81231		
Co	0.97061	0.349	0.96874	0.97453	0.96856		
Cr	5.11786	0.650	5.14119	5.13267	5.07972		
Cu	4.83376	0.642	4.86219	4.83850	4.80058		
Fe	4.90580	0.559	4.92925	4.91253	4.87562		
K	49.00690	0.353	49.1860	48.9939	48.8406		
Li	25.65088	0.827	25.8589	25.6589	25.4346		
Mg	23.76412	0.557	23.8901	23.7761	23.6260		
Mn	4.91667	0.560	4.93666	4.92809	4.88526		
Mo	24.58717	0.788	24.7428	24.6487	24.3698		
Na	50.17737	0.486	50.4154	50.1885	49.9280		
Ni	9.70111	0.668	9.74828	9.72785	9.62718		
P	25.25264	0.598	25.2130	25.4195	25.1252		
Pb	24.76503	0.596	24.9096	24.7710	24.6143		
Sb	5.03710	0.447	5.05093	5.04931	5.01106		
Se	2.02966	2.191	2.00586	2.00214	2.08096		
Si	1.99582	0.654	1.99675	2.00839	1.98231		
Sn	4.90283	0.988	4.95048	4.90445	4.85355		
Sr	1.01406	0.743	1.02072	1.01559	1.00587		
Ti	2.04779	0.806	2.06288	2.05037	2.03013		
Tl	5.28792	0.907	5.33623	5.28728	5.24025		
V	2.02085	0.708	2.03197	2.02589	2.00470		
Zn	4.93608	0.442	4.94952	4.94783	4.91088		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 37

INSTRUMENT ID: 02360

37 CCB 12/20/92 00:53

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.00000	84.24	0.00000	0.00000	0.00000		
__Al	0.02021	59.54	0.03394	0.01520	0.01149		
__As	-0.00506	159.3	0.00408	-0.0081	-0.0111		
__B	-0.00070	175.0	0.00072	-0.0014	-0.0014		
__Ba	0.00400	69.28	0.00720	0.00240	0.00240		
__Be	0.00044	168.8	0.00132	0.00001	0.00001		
__Ca	0.04054	94.86	0.08462	0.02327	0.01375		
__Cd	0.00481	87.64	0.00963	0.00175	0.00306		
__Co	0.00159	124.8	0.00047	0.00390	0.00041		
__Cr	0.00042	751.1	0.00409	-0.0011	-0.0016		
__Cu	0.00040	991.7	0.00496	-0.0012	-0.0025		
__Fe	0.00234	188.7	0.00703	0.00176	-0.0017		
__K	0.12813	60.01	0.12808	0.20504	0.05125		
__Li	0.01818	101.3	0.03940	0.00606	0.00909		
__Mg	0.00218	690.5	0.01957	-0.0065	-0.0065		
__Mn	0.00122	321.0	0.00571	-0.0004	-0.0016		
__Mo	0.01863	114.5	0.04317	0.00849	0.00424		
__Na	0.07003	90.86	0.14287	0.02521	0.04201		
__Ni	0.00279	275.0	0.01142	-0.0033	0.00031		
__P	-0.04942	194.3	-0.0411	-0.1493	0.04225		
__Pb	0.01191	178.6	0.03625	-0.0031	0.00266		
__Sb	-0.00040	1201.	0.00285	0.00196	-0.0060		
__Se	0.00832	34.71	0.00665	0.00666	0.01166		
__Si	0.00588	58.76	0.00561	0.00947	0.00257		
__Sn	-0.00344	563.1	0.01874	-0.0171	-0.0119		
__Sr	0.00035	261.9	0.00138	0.00010	-0.0004		
__Ti	0.00041	624.0	0.00333	-0.0004	-0.0016		
__Tl	-0.05173	152.0	0.00525	-0.0189	-0.1414		
__V	-0.00281	61.81	-0.0008	-0.0038	-0.0038		
__Zn	0.00497	77.44	0.00929	0.00375	0.00188		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 38

INSTRUMENT ID: 02360

A1 Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

38 1908987 J BMTF 12/20/92 00:58

****,100-100,DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00062	180.1	-0.0019	0.00002	0.00002		
X Al	0.01374	29.76	0.01791	0.01360	0.00973		
As	-0.00884	155.4	0.00437	-0.0078	-0.0230		
B	0.03614	4.597	0.03650	0.03759	0.03433		
X Ba	0.03841	0.000	0.03841	0.03841	0.03841		
X Be	0.00000	57.95	0.00001	0.00000	0.00001		
X Ca	17.00306	0.585	16.9158	16.9819	17.1114		
Cd	-0.00007	2856.	0.00036	0.00168	-0.0022		
X Co	0.00123	138.1	-0.0004	0.00300	0.00112		
X Cr	-0.00170	38.76	-0.0015	-0.0011	-0.0024		
X Cu	0.00062	439.0	-0.0006	0.00374	-0.0012		
X Fe	0.17020	1.584	0.17079	0.16726	0.17255		
X K	5.76762	1.539	5.81889	5.66509	5.81890		
Li	-0.00101	173.2	0.00000	0.00000	-0.0030		
X Mg	6.46044	0.485	6.44742	6.43766	6.49623		
X Mn	0.07028	0.002	0.07028	0.07028	0.07028		
Mo	0.00022	1198.	0.00211	0.00140	-0.0028		
X Na	9.02403	0.918	8.96800	8.98480	9.11928		
X Ni	-0.00215	48.82	-0.0033	-0.0015	-0.0015		
P	-0.03197	171.6	-0.0949	0.00590	-0.0069		
Pb	-0.01384	27.73	-0.0149	-0.0170	-0.0095		
X Sb	0.01506	17.54	0.01604	0.01708	0.01207		
Se	0.00049	1464.	0.00465	-0.0078	0.00466		
Si	4.71109	0.602	4.67932	4.72005	4.73390		
Sn	0.01181	45.21	0.00582	0.01352	0.01609		
Sr	0.08252	0.651	0.08195	0.08259	0.08302		
Ti	0.00069	138.3	0.00125	0.00125	-0.0004		
Tl	-0.06269	146.2	0.03793	-0.1414	-0.0845		
X V	-0.00286	58.17	-0.0038	-0.0009	-0.0038		
X Zn	0.03342	0.038	0.03343	0.03340	0.03342		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 39 INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00046	624.4	-0.0004	0.00245	-0.0033		
X Al	2.32638	0.907	2.30923	2.34997	2.31994		
As	-0.01013	114.9	-0.0203	-0.0126	0.00257		
B	0.02876	4.322	0.02733	0.02949	0.02947		
X Ba	0.05049	0.000	0.05049	0.05049	0.05049		
X Be	0.00000	1339.	-0.0000	-0.0000	0.00000		
X Ca	10.35077	0.389	10.3629	10.3058	10.3835		
Cd	0.00026	496.1	-0.0010	0.00158	0.00026	D	
X Co	0.00113	63.05	0.00031	0.00160	0.00148		
X Cr	0.00265	134.9	0.00066	0.00679	0.00051		
X Cu	0.00451	57.57	0.00243	0.00743	0.00368		
X Fe	2.74065	0.098	2.74301	2.73771	2.74124		
X K	2.84537	3.120	2.89664	2.89664	2.74284		
Li	-0.00000	17938	-0.0030	0.00303	-0.0000		
X Mg	4.36122	0.258	4.35471	4.35473	4.37424		
X Mn	0.11707	0.603	0.11667	0.11789	0.11666		
Mo	-0.00177	79.56	-0.0017	-0.0003	-0.0031		
X Na	9.72492	0.792	9.74175	9.64087	9.79213		
X Ni	0.00311	90.19	0.00372	0.00555	0.00004		
P	0.06270	196.4	0.18787	0.05861	-0.0583		
Pb	-0.00345	251.1	-0.0095	-0.0073	0.00648		
X Sb	0.00263	311.8	0.00968	0.00463	-0.0064		
Se	0.01807	62.35	0.00641	0.01890	0.02890		
Si	7.44640	0.840	7.37823	7.50121	7.45977		
Sn	0.00324	158.2	0.00323	0.00837	-0.0018		
Sr	0.06186	0.690	0.06229	0.06143	0.06186		
Ti	0.09854	5.324	0.10341	0.09924	0.09298		
Tl	-0.05089	80.38	-0.0886	-0.0074	-0.0565		
X V	0.00103	161.7	0.00198	0.00200	-0.0008		
X Zn	0.02934	0.060	0.02935	0.02932	0.02935		

AR318946

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 40 INSTRUMENT ID: 02360

AlSbBaBeCdCaCrCoCuFeMgMnNiKAgNaVZn

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00126	177.5	-0.0038	0.00003	0.00003		
X Al	0.01920	0.037	0.01921	0.01919	0.01920		
As	-0.00572	110.6	0.00137	-0.0077	-0.0107		
B	0.02658	4.054	0.02659	0.02549	0.02765		
X Ba	0.03121	0.000	0.03121	0.03121	0.03121		
X Be	0.00000	84.42	0.00000	0.00000	0.00001		
X Ca	10.55016	0.472	10.4941	10.5670	10.5892		
Cd	0.00038	342.0	-0.0009	0.00038	0.00169	D	
X Co	0.00051	281.8	-0.0006	0.00009	0.00214		
X Cr	-0.00400	40.41	-0.0041	-0.0055	-0.0022		
X Cu	0.00291	65.57	0.00124	0.00249	0.00499		
X Fe	0.21474	0.855	0.21327	0.21415	0.21680		
X K	2.84539	1.560	2.81976	2.81976	2.89666		
Li	-0.00101	173.2	-0.0000	-0.0000	-0.0030		
X Mg	4.31051	0.455	4.28990	4.32895	4.31269		
X Mn	0.08745	0.000	0.08745	0.08745	0.08745		
Mo	-0.00143	0.039	-0.0014	-0.0014	-0.0014		
X Na	10.18951	0.381	10.2119	10.1446	10.2119		
X Ni	0.00028	1296.	-0.0033	0.00028	0.00394		
P	-0.02320	93.58	-0.0204	-0.0461	-0.0030		
Pb	-0.00958	38.52	-0.0138	-0.0074	-0.0074		
X Sb	0.01275	80.92	0.01004	0.00405	0.02416		
Se	0.00781	359.7	-0.0230	0.03197	0.01447		
Si	4.60090	0.072	4.59860	4.59938	4.60473		
Sn	-0.00353	83.69	-0.0018	-0.0069	-0.0018		
Sr	0.06165	0.600	0.06122	0.06186	0.06186		
Ti	0.00208	59.97	0.00208	0.00083	0.00333		
Tl	0.08143	144.9	0.21743	0.02157	0.00529		
X V	-0.00190	86.79	-0.0009	-0.0009	-0.0038		
X Zn	0.01601	3.470	0.01665	0.01570	0.01568		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 41 INSTRUMENT ID: 02360

Al Ba Sb Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00273	41.19	-0.0014	-0.0033	-0.0033		
X Al	2.28620	3.139	2.32003	2.33483	2.20376		
As	-0.00438	106.4	-0.0003	-0.0033	-0.0094		
B	0.04430	2.805	0.04360	0.04356	0.04573		
X Ba	0.06010	0.000	0.06010	0.06010	0.06010		
X Be	0.00000	193.1	0.00000	0.00000	-0.0000		
X Ca	15.92290	0.425	15.8448	15.9659	15.9579		
Cd	0.00132	28.72	0.00154	0.00154	0.00088	D	
X Co	0.00242	29.92	0.00203	0.00198	0.00326		
X Cr	0.00275	48.15	0.00163	0.00421	0.00240		
X Cu	0.01242	10.06	0.01117	0.01242	0.01368		
X Fe	2.83761	0.342	2.83114	2.84879	2.83290		
X K	5.44718	1.078	5.39592	5.43436	5.51126		
Li	0.00404	114.5	-0.0000	0.00303	0.00909		
X Mg	6.33783	0.330	6.31396	6.34652	6.35301		
X Mn	0.13453	1.048	0.13372	0.13616	0.13372		
Mo	0.00010	803.4	-0.0003	0.00104	-0.0003		
X Na	9.51207	0.620	9.45605	9.57369	9.50646		
X Ni	0.00127	166.3	0.00371	0.00005	0.00004		
P	0.08999	63.32	0.03584	0.14945	0.08468		
Pb	-0.00240	423.6	-0.0041	-0.0116	0.00856		
X Sb	0.00024	1239.	-0.0003	-0.0024	0.00360		
Se	0.00490	561.6	0.00656	0.03155	-0.0234		
Si	7.70292	1.788	7.76079	7.80232	7.54567		
Sn	-0.00368	383.0	0.00572	-0.0199	0.00313		
Sr	0.08081	0.807	0.08024	0.08152	0.08067		
Ti	0.11355	1.811	0.11591	0.11216	0.11258		
Tl	-0.09354	116.1	-0.0228	-0.2186	-0.0391		
X V	0.00055	257.2	-0.0008	0.00056	0.00199		
X Zn	0.08162	0.000	0.08162	0.08162	0.08162		

AR318948

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 42

INSTRUMENT ID: 02360

Al Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00257	115.6	-0.0058	0.00002	-0.0019		
X Al	0.01720	43.91	0.00912	0.01840	0.02409		
As	0.00185	494.9	0.01050	-0.0077	0.00285		
B	0.03521	3.557	0.03666	0.03450	0.03448		
X Ba	0.03841	0.000	0.03841	0.03841	0.03841		
X Be	0.00001	74.52	0.00002	0.00001	0.00000		
X Ca	15.62144	0.416	15.5496	15.6380	15.6766		
Cd	-0.00006	4065.	-0.0022	0.00299	-0.0009	D	
X Co	0.00138	170.9	-0.0011	0.00344	0.00189		
X Cr	0.00037	259.1	0.00009	0.00146	-0.0004		
X Cu	0.00166	114.8	0.00124	0.00374	-0.0000		
X Fe	0.18521	1.672	0.18227	0.18845	0.18491		
X K	5.17806	0.857	5.12679	5.20369	5.20369		
Li	0.00101	173.2	-0.0000	-0.0000	0.00303		
X Mg	6.05489	0.543	6.03752	6.03429	6.09287		
X Mn	0.09148	1.542	0.08985	0.09229	0.09229		
Mo	-0.00190	43.00	-0.0014	-0.0014	-0.0028		
X Na	9.64039	0.174	9.64039	9.62358	9.65720		
X Ni	0.00089	427.6	0.00211	0.00394	-0.0033		
P	-0.03038	194.7	-0.0930	0.02454	-0.0226		
Pb	-0.00691	87.42	-0.0026	-0.0042	-0.0138		
X Sb	0.00034	617.6	-0.0020	0.00203	0.00099		
Se	0.01045	151.9	0.01962	-0.0078	0.01962		
Si	4.81584	0.217	4.80379	4.82300	4.82071		
Sn	-0.00014	7405.	0.00327	0.00840	-0.0121		
Sr	0.07864	0.407	0.07832	0.07864	0.07896		
Ti	0.00146	14.27	0.00125	0.00167	0.00146		
Tl	0.07045	166.7	0.03778	0.20085	-0.0272		
X V	-0.00381	75.23	-0.0066	-0.0038	-0.0009		
X Zn	0.02781	0.060	0.02781	0.02779	0.02782		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 43 INSTRUMENT ID: 02360

A 1 Sb Ba Be Cd Ca Cr Co Cu Fe Mg Mn Ni K Ag Na V Zn

43 1908992 CRHLT 12/20/92 01:21

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00191	102.0	-0.0038	-0.0019	0.00003		
X Al	0.06511	24.26	0.04745	0.06998	0.07790		
__As	-0.01633	35.24	-0.0229	-0.0138	-0.0122		
__B	0.02981	8.359	0.02693	0.03125	0.03125		
X Ba	0.02441	2.838	0.02401	0.02401	0.02521		
X Be	0.00000	85.67	0.00001	0.00000	0.00000		
X Ca	8.38826	0.359	8.39496	8.41453	8.35529		
__Cd	-0.00004	1673.	-0.0009	0.00039	0.00039	D	
X Co	0.00030	1079.	-0.0011	0.00408	-0.0019		
X Cr	-0.00424	47.64	-0.0057	-0.0019	-0.0050		
X Cu	0.00082	174.0	-0.0000	-0.0000	0.00249		
X Fe	0.24983	0.815	0.25218	0.24865	0.24865		
X K	2.56340	1.732	2.58904	2.51213	2.58903		
__Li	-0.00101	173.2	-0.0030	-0.0000	-0.0000		
X Mg	3.32215	0.299	3.33301	3.31999	3.31346		
X Mn	0.07363	1.915	0.07526	0.07281	0.07281		
__Mo	-0.00143	98.44	-0.0028	-0.0014	-0.0000		
X Na	6.84717	0.464	6.82475	6.88358	6.83317		
X Ni	-0.00094	112.9	-0.0015	-0.0015	0.00028		
__P	-0.07955	48.45	-0.0993	-0.0351	-0.1041		
__Pb	-0.00440	104.0	0.00055	-0.0052	-0.0084		
X Sb	0.00135	172.0	0.00403	0.00000	0.00000		
__Se	0.00522	280.3	-0.0056	-0.0006	0.02188		
__Si	3.41187	0.496	3.40061	3.43135	3.40365		
__Sn	-0.00522	150.0	-0.0120	-0.0069	0.00332		
__Sr	0.04754	0.000	0.04754	0.04754	0.04754		
__Ti	0.00055	302.5	-0.0004	-0.0004	0.00250		
__Tl	-0.10342	39.73	-0.1415	-0.1088	-0.0598		
X V	-0.00191	87.21	-0.0038	-0.0009	-0.0009		
X Zn	0.01107	0.085	0.01107	0.01107	0.01106		

132

AR318950

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 44

INSTRUMENT ID: 02360

A/SbBaBeCdCaCrCoCuFeMgMnNi'KAgNaVZn

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ELEM	AVERAGE (ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
X Ag	-0.00265	85.04	-0.0013	-0.0013	-0.0052		
X Al	3.00210	0.786	3.01548	3.01598	2.97484		
As	-0.00879	0.029	-0.0087	-0.0087	-0.0088		
B	0.02911	0.033	0.02912	0.02911	0.02910		
X Ba	0.04571	0.000	0.04571	0.04571	0.04571		
X Be	-0.00000	77.60	-0.0000	-0.0000	-0.0000		
X Ca	8.65006	0.184	8.64512	8.63719	8.66786		
Cd	-0.00062	122.3	-0.0010	0.00025	-0.0010	D	
X Co	0.00105	289.4	-0.0021	0.00388	0.00143		
X Cr	0.00299	53.62	0.00177	0.00239	0.00480		
X Cu	0.00533	48.69	0.00242	0.00742	0.00617		
X Fe	3.23942	0.041	3.23913	3.23824	3.24089		
K	2.92224	1.519	2.89662	2.97351	2.89661		
Li	0.00101	173.2	0.00303	-0.0000	-0.0000		
X Mg	3.57565	0.689	3.54744	3.59301	3.58650		
X Mn	0.10689	0.000	0.10689	0.10689	0.10689		
Mo	-0.00185	76.09	-0.0032	-0.0004	-0.0018		
X Na	6.77773	0.715	6.83374	6.74972	6.74973		
X Ni	0.00428	24.71	0.00550	0.00367	0.00367		
P	0.04343	34.69	0.03810	0.06044	0.03175		
Pb	0.00079	1597.	0.01536	-0.0059	-0.0070		
X Sb	-0.00045	1784.	-0.0004	-0.0084	0.00757		
Se	0.00785	91.82	0.00368	0.01618	0.00369		
Si	7.40199	0.950	7.46556	7.41401	7.32640		
Sn	-0.01294	97.69	-0.0043	-0.0274	-0.0069		
Sr	0.05018	0.245	0.05011	0.05011	0.05032		
Ti	0.12300	3.439	0.11842	0.12384	0.12676		
Tl	-0.12069	94.78	-0.2514	-0.0391	-0.0715		
X V	0.00247	33.65	0.00198	0.00199	0.00343		
X Zn	0.03302	0.028	0.03303	0.03301	0.03302		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 45

INSTRUMENT ID: 02360

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	-0.03458	6.498	-0.0321	-0.0350	-0.0365		
__Al	467.04864	0.529	465.178	466.114	469.853		
__As	0.20143	16.26	0.18368	0.18138	0.23924		
__B	0.27537	2.250	0.27030	0.27353	0.28228		
__Ba	0.00565	0.663	0.00562	0.00564	0.00570		
__Be	0.00108	35.66	0.00064	0.00131	0.00130		
__Ca	461.27288	0.755	458.259	460.467	465.090		
__Cd	0.03102	2.457	0.03059	0.03190	0.03057		
__Co	0.00127	141.2	0.00036	0.00010	0.00334		
__Cr	-0.01621	6.322	-0.0168	-0.0150	-0.0167		
__Cu	0.00077	302.8	-0.0019	0.00182	0.00243		
__Fe	176.77893	0.663	175.901	176.324	178.110		
__K	0.24379	18.24	0.21810	0.21813	0.29515		
__Li	-0.00202	173.1	-0.0060	-0.0000	-0.0000		
__Mg	482.41149	0.558	480.316	481.468	485.448		
__Mn	-0.01240	1.050	-0.0123	-0.0123	-0.0125		
__Mo	-0.04901	4.816	-0.0464	-0.0494	-0.0511		
__Na	0.10373	9.347	0.10933	0.09254	0.10934		
__Ni	-0.02623	6.633	-0.0261	-0.0280	-0.0245		
__P	-1.69193	11.18	-1.5375	-1.6352	-1.9030		
__Pb	0.02358	17.19	0.02051	0.02205	0.02818		
__Sb	0.00537	289.3	0.00646	-0.0106	0.02036		
__Se	-0.05911	36.90	-0.0521	-0.0835	-0.0416		
__Si	-0.00012	4168.	0.00226	0.00359	-0.0062		
__Sn	0.09084	22.93	0.06837	0.10953	0.09462		
__Sr	0.01763	0.000	0.01762	0.01763	0.01763		
__Ti	0.01097	13.34	0.00958	0.01250	0.01083		
__Tl	-0.44057	14.22	-0.5099	-0.3881	-0.4235		
__V	0.00414	69.39	0.00699	0.00125	0.00417		
__Zn	-0.00981	2.264	-0.0098	-0.0100	-0.0095		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 46

INSTRUMENT ID: 02360

46 ICSAB 12/20/92 01:35

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.87996	1.210	0.89220	0.87491	0.87278		
Al	470.97653	0.329	469.226	472.191	471.511		
As	0.20806	9.376	0.22839	0.18949	0.20630		
B	0.28389	0.851	0.28341	0.28651	0.28175		
Ba	0.48022	0.292	0.47860	0.48104	0.48101		
Be	0.45086	0.445	0.44868	0.45262	0.45130		
Ca	464.76666	0.457	463.003	467.127	464.169		
Cd	0.88326	0.299	0.88085	0.88609	0.88285		
Co	0.42927	1.334	0.42312	0.43447	0.43021		
Cr	0.43870	1.202	0.43713	0.44459	0.43439		
Cu	0.45843	1.100	0.45302	0.46301	0.45926		
Fe	178.30604	0.388	177.691	179.055	178.171		
K	0.24419	18.18	0.21850	0.21860	0.29547		
Li	-0.00101	624.4	-0.0060	0.00606	-0.0030		
Mg	485.87545	0.361	483.946	487.386	486.293		
Mn	0.43241	0.546	0.43005	0.43478	0.43240		
Mo	-0.05085	7.273	-0.0469	-0.0542	-0.0513		
Na	0.10619	16.47	0.09218	0.10059	0.12581		
Ni	0.83005	0.254	0.83136	0.83118	0.82761		
P	-1.86569	6.756	-1.7276	-1.9746	-1.8947		
Pb	0.91539	0.967	0.92411	0.90641	0.91566		
Sb	0.01255	102.8	-0.0021	0.01785	0.02198		
Se	-0.05860	69.54	-0.0733	-0.0125	-0.0899		
Si	-0.00063	1324.	-0.0003	0.00760	-0.0091		
Sn	0.11042	4.466	0.11527	0.10541	0.11058		
Sr	0.01774	0.695	0.01781	0.01781	0.01760		
Ti	0.00993	6.289	0.00930	0.00993	0.01055		
Tl	-0.41729	10.99	-0.3645	-0.4399	-0.4474		
V	0.46639	0.713	0.46255	0.46833	0.46829		
Zn	0.89625	0.529	0.89273	0.90165	0.89436		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 47 INSTRUMENT ID: 02360

47 CRI 12/20/92 01:40

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.09825	1.135	0.09762	0.09760	0.09954		
__Al	0.57454	33.94	0.78838	0.52875	0.40650		
__As	0.00899	71.82	0.01624	0.00693	0.00381		
__B	0.00084	124.2	0.00206	0.00018	0.00029		
__Ba	0.00240	0.106	0.00241	0.00240	0.00240		
__Be	0.01004	3.745	0.01025	0.01026	0.00960		
__Ca	0.58723	34.17	0.80972	0.53205	0.41993		
__Cd	0.01095	11.99	0.01095	0.00963	0.01226		
__Co	0.10075	3.540	0.09674	0.10356	0.10196		
__Cr	0.02231	3.713	0.02268	0.02289	0.02136		
__Cu	0.04748	2.632	0.04748	0.04873	0.04623		
__Fe	0.23587	33.93	0.32471	0.21351	0.16938		
__K	0.15382	28.85	0.12819	0.12819	0.20508		
__Li	-0.00000	152.7	-0.0000	-0.0000	-0.0000		
__Mg	0.59271	34.56	0.82049	0.53414	0.42350		
__Mn	0.02891	0.030	0.02890	0.02891	0.02892		
__Mo	-0.00053	152.7	-0.0000	-0.0014	-0.0000		
__Na	0.03916	44.65	0.05876	0.02516	0.03356		
__Ni	0.07962	5.791	0.07534	0.08450	0.07901		
__P	-0.10809	48.18	-0.0624	-0.1648	-0.0969		
__Pb	0.00668	155.2	0.01851	0.00239	-0.0008		
__Sb	0.11088	5.798	0.11350	0.10355	0.11557		
__Se	0.01898	94.83	0.03895	0.01398	0.00400		
__Si	0.00464	50.05	0.00718	0.00412	0.00262		
__Sn	-0.00845	92.92	-0.0015	-0.0067	-0.0170		
__Sr	-0.00018	67.12	-0.0001	-0.0003	-0.0001		
__Ti	0.00292	37.68	0.00209	0.00418	0.00251		
__Tl	-0.01846	204.1	0.00322	-0.0620	0.00338		
__V	0.09880	3.365	0.10264	0.09688	0.09688		
__Zn	0.04064	0.059	0.04063	0.04063	0.04067		

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 48

INSTRUMENT ID: 02360

48 CCV . 12/20/92 01:45

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.46441	0.484	0.46181	0.46571	0.46571	_____	_____
__Al	23.91110	0.367	23.9111	23.9989	23.8231	_____	_____
__As	2.06928	0.348	2.07742	2.06671	2.06371	_____	_____
__B	0.97421	0.482	0.97091	0.97959	0.97212	_____	_____
__Ba	4.95586	0.452	4.95946	4.97626	4.93185	_____	_____
__Be	0.47830	0.415	0.47786	0.48047	0.47657	_____	_____
__Ca	48.79493	0.318	48.7270	48.9729	48.6847	_____	_____
__Cd	4.85434	0.501	4.82937	4.87798	4.85565	_____	_____
__Co	0.96776	0.431	0.96483	0.97254	0.96592	_____	_____
__Cr	5.11542	0.167	5.11293	5.12498	5.10835	_____	_____
__Cu	4.83562	0.390	4.84045	4.85159	4.81481	_____	_____
__Fe	4.94020	0.364	4.94788	4.95309	4.91964	_____	_____
__K	49.14785	0.413	49.0710	49.3783	48.9941	_____	_____
__Li	25.78120	0.568	25.8014	25.9165	25.6256	_____	_____
__Mg	23.84337	0.458	23.8996	23.9129	23.7174	_____	_____
__Mn	4.92522	0.287	4.91584	4.94154	4.91829	_____	_____
__Mo	24.60817	0.317	24.5538	24.6975	24.5730	_____	_____
__Na	53.75790	0.945	54.0885	54.0128	53.1723	_____	_____
__Ni	9.71369	0.284	9.70102	9.74540	9.69465	_____	_____
__P	25.41547	1.110	25.1546	25.7150	25.3767	_____	_____
__Pb	24.76223	0.259	24.7727	24.8207	24.6932	_____	_____
__Sb	5.06257	0.444	5.04766	5.08846	5.05157	_____	_____
__Se	2.02839	0.647	2.02965	2.04085	2.01466	_____	_____
__Si	2.00963	0.922	1.99214	2.02907	2.00767	_____	_____
__Sn	4.88400	0.456	4.89433	4.89927	4.85840	_____	_____
__Sr	1.01441	0.395	1.01431	1.01847	1.01046	_____	_____
__Ti	2.05064	0.377	2.05286	2.05704	2.04202	_____	_____
__Tl	5.31213	1.189	5.28001	5.38491	5.27147	_____	_____
__V	2.02667	0.472	2.02696	2.03610	2.01695	_____	_____
__Zn	4.94132	0.329	4.92543	4.95799	4.94055	_____	_____

Run Name: 9235402I11

LANCASTER LABORATORIES

PAGE: 49

INSTRUMENT ID: 02360

49 CCB 12/20/92 01:49

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	0.00000	24065	0.00390	0.00000	-0.0039		
__Al	0.07335	36.18	0.07578	0.09859	0.04568		
__As	-0.01114	72.34	-0.0080	-0.0050	-0.0202		
__B	0.00069	307.7	0.00284	0.00069	-0.0014		
__Ba	0.01040	48.03	0.01440	0.01200	0.00480		
__Be	0.00066	98.39	0.00066	0.00131	0.00001		
__Ca	0.11177	40.77	0.13539	0.14068	0.05923		
__Cd	0.00875	43.30	0.01094	0.01094	0.00437		
__Co	0.00307	67.30	0.00429	0.00068	0.00424		
__Cr	0.00708	77.09	0.01167	0.00855	0.00104		
__Cu	0.00788	74.73	0.01245	0.00995	0.00123		
__Fe	0.01615	50.68	0.01850	0.02291	0.00705		
__K	0.25625	17.31	0.28186	0.28187	0.20502		
__Li	0.04445	51.62	0.06061	0.05455	0.01818		
__Mg	0.05864	40.07	0.07819	0.06516	0.03256		
__Mn	0.00693	61.12	0.00937	0.00937	0.00203		
__Mo	0.04482	45.60	0.05732	0.05591	0.02123		
__Na	0.09244	48.10	0.12606	0.10925	0.04202		
__Ni	0.01692	84.96	0.02611	0.02428	0.00035		
__P	0.02083	81.03	0.02070	0.00401	0.03778		
__Pb	0.04159	38.28	0.04799	0.05333	0.02346		
__Sb	0.01393	87.52	0.01994	0.02196	-0.0001		
__Se	-0.00085	774.6	0.00664	-0.0033	-0.0058		
__Si	0.01046	51.97	0.00943	0.01635	0.00561		
__Sn	-0.00689	148.4	0.00334	-0.0069	-0.0171		
__Sr	0.00160	69.27	0.00224	0.00224	0.00032		
__Ti	0.00361	56.94	0.00459	0.00500	0.00125		
__Tl	0.00508	1599.	0.00486	0.08647	-0.0760		
__V	-0.00079	371.2	-0.0007	0.00212	-0.0037		
__Zn	0.00926	52.68	0.01296	0.01110	0.00373		

Data Reviewed By: [illegible] 12-20-92

Data Verified By: [illegible] 12-21-92

137

Name: 9235601I11
Data Reviewed By: [illegible] 12/21/92
Data Verified By: Mary Snavely 12-22-92

LANCASTER LABORATORIES

PAGE: 1 INSTRUMENT ID: 02360

1 SO 5Q 12/21/92 14:33

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	RE[illegible]	REDIG
__Ag	22.66666	5.094	22.0000	22.0000	24.0000		
__Al	577.33331	0.529	578.000	574.000	580.000		
__As	23.00000	15.67	26.0000	24.0000	19.0000		
__B	6.66666	17.32	6.00000	8.00000	6.00000		
__Ba	52.00000	0.000	52.0000	52.0000	52.0000		
__Be	22.00000	0.000	22.0000	22.0000	22.0000		
__Ca	600.66668	1.017	594.000	606.000	602.000		
__Cd	0.00000	0.000	0.00000	0.00000	0.00000		
__Co	3.18518	96.75	-0.2222	4.00000	5.77777		
__Cr	10.29628	13.18	9.11108	11.7777	10.0000		
__Cu	8.33333	24.97	9.00000	10.0000	6.00000		
__Fe	9.33333	32.73	12.0000	10.0000	6.00000		
__K	258.33334	1.117	255.000	260.000	260.000		
__Li	0.00000	0.000	0.00000	0.00000	0.00000		
__Mg	7.00000	14.28	7.00000	6.00000	8.00000		
__Mn	3.66666	15.74	3.00000	4.00000	4.00000		
__Mo	1.33333	86.60	2.00000	0.00000	2.00000		
__Na	152.66667	0.756	152.000	154.000	152.000		
__Ni	5.33333	86.60	0.00000	8.00000	8.00000		
__P	-0.87356	174.8	0.20689	-2.6206	-0.2068		
__Pb	2.66666	60.62	3.60000	3.60000	0.79998		
__Sb	6.66666	217.0	14.0000	16.0000	-10.000		
__Se	6.00000	145.2	0.00000	16.0000	2.00000		
__Si	41.00000	24.02	33.0000	52.0000	38.0000		
__Sn	-1.33333	229.1	2.00000	-4.0000	-2.0000		
__Sr	0.00000	0.000	-2.0000	0.00000	2.00000		
__Ti	67.33333	14.03	60.0000	78.0000	64.0000		
__Tl	21.66666	37.01	21.0000	30.0000	14.0000		
__V	0.00000	0.000	0.00000	0.00000	0.00000		
__Zn	43.66666	1.322	44.0000	43.0000	44.0000		

[illegible]

AR318957

Run Name: 1235601I11

LANCASTER LABORATORIES

PAGE:

INSTRUMENT ID: 02360

? 51 S 12/21/92 14:37

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ba	8620.33300	0.736	8622.00	8556.00	8683.00	_____	_____
__Cd	16200.00000	0.407	16191.0	16139.0	16270.0	_____	_____
__Cr	22612.96289	0.653	22587.5	22479.4	22771.8	_____	_____
__Cu	16429.00000	0.712	16425.0	16314.0	16548.0	_____	_____
__Fe	11747.66699	0.719	11744.0	11665.0	11834.0	_____	_____
__Mn	8524.33300	0.694	8514.00	8471.00	8588.00	_____	_____
__Mo	74180.33593	0.658	74146.0	73710.0	74685.0	_____	_____
__Ni	22877.00000	0.592	22874.0	22743.0	23014.0	_____	_____
__Pb	19857.80078	0.413	19858.8	19775.1	19939.4	_____	_____
__Zn	11370.33300	0.557	11358.0	11314.0	11439.0	_____	_____

139

AR318958

n Name: 9235601I11

LANCASTER LABORATORIES

PAGE: 3

INSTRUMENT [illegible]: 02360

3 S2 S 12/21/92 14:41

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	RELOAD	REDIG
__As	1806.33337	1.474	1777.00	1813.00	1829.00		
__Be	1622.33337	1.560	1595.00	1627.00	1645.00		
__Co	3663.51855	1.664	3596.66	3677.77	3716.11		
__Li	16614.00000	1.676	16326.0	16634.0	16882.0		
__P	5985.39062	1.558	5892.48	5984.68	6079.00		
__Sb	10583.33300	1.168	10442.0	10636.0	10672.0		
__Se	4340.33349	0.837	4329.00	4311.00	4381.00		
__Si	6792.66650	1.723	6665.00	6818.00	6895.00		
__Sn	4484.66650	1.046	4432.00	4500.00	4522.00		
__Sr	9759.00000	1.708	9572.00	9813.00	9892.00		
__Ti	25267.33398	1.596	24844.0	25311.0	25647.0		
__Tl	1323.33337	1.513	1304.00	1322.00	1344.00		
__V	3675.00000	1.558	3610.00	3697.00	3718.00		

120

AR318959

Run Name 1235601I11

LANCASTER LABORATORIES

PAGE: 4 INSTRUMENT ID: 02360

4 S 12/21/92 14:46

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	1068.00000	0.374	1072.00	1068.00	1064.00		
B	1914.33337	0.030	1914.00	1914.00	1915.00		

AR318960

Run Name: 935601I11

LANCASTER LABORATORIES

PAGE: 5

INSTRUMENT ID: 02360

S 91 S

12/21/92 14:49

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__K	13548.00000	0.472	13480.0	13557.0	13607.0	_____	_____
__Na	58359.00000	0.736	57865.0	58566.0	58646.0	_____	_____

Run Name: 235601I11

LANCASTER LABORATORIES

PAGE: 6

INSTRUMENT ID: 02360

6 55 S 12/21/92 14:51

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Al	7183.66601	0.531	27072.0	27347.0	27132.0		
Ca	97571.32812	0.409	196943.	198484.	197287.		
Mg	1405.66601	0.437	31316.0	31564.0	31337.0		

Name: [illegible]5601I11

LANCASTER LABORATORIES

PAGE: 7 INSTRUMENT ID: 02360

7 ICV 12/21/92 14:54

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible]8267	4.118	0.50562	0.47120	0.47119		
Al	.86079	0.501	23.9986	23.7846	23.7990		
As	[illegible]7305	0.912	1.99356	1.96566	1.95991		
B	[illegible]7181	0.004	0.97176	0.97184	0.97183		
Ba	[illegible]7207	0.691	4.91097	4.85262	4.85262		
Be	[illegible]6108	0.459	0.46353	0.45989	0.45982		
Ca	[illegible].00273	0.474	49.2709	48.8566	48.8805		
Cd	[illegible]5882	0.339	4.77733	4.75142	4.74772		
Co	[illegible]4843	0.572	0.95443	0.94385	0.94700		
Cr	[illegible]4182	0.490	5.07033	5.02576	5.02938		
Cu	[illegible]4641	0.643	4.78146	4.72552	4.73225		
Fe	[illegible]5225	0.472	4.87858	4.84121	4.83695		
[illegible]	.24823	1.087	50.7748	49.6839	50.2859		
Li	[illegible].64915	0.634	24.8287	24.5428	24.5759		
Mg	[illegible].66769	0.536	23.8133	23.6091	23.5805		
Mn	[illegible]4966	0.497	4.87745	4.83400	4.83752		
Mo	[illegible].22146	0.305	24.3039	24.1994	24.1610		
Na	[illegible].97985	0.734	51.4093	50.7221	50.8080		
Ni	[illegible]5005	0.398	9.59282	9.53746	9.51987		
P	[illegible].26931	0.992	24.4224	24.3937	23.9917		
Pb	[illegible].52565	0.516	24.6719	24.4517	24.4532		
Sb	[illegible].90474	0.095	4.90450	4.90955	4.90018		
Se	[illegible]6019	0.579	1.94707	1.96676	1.96675		
Si	[illegible]4644	0.852	2.05079	2.02723	2.06130		
Sn	[illegible]9816	0.819	4.84338	4.77217	4.77892		
Sr	[illegible]8035	0.722	0.98851	0.97581	0.97673		
Ti	[illegible]7015	0.448	1.98027	1.96399	1.96617		
Tl	[illegible]0465	0.459	5.02953	4.98416	5.00026		
V	[illegible]6055	0.963	1.97722	1.94003	1.96440		
Zn	[illegible]3694	0.281	4.85225	4.82619	4.83239		

Run Name: 9235601I11

LANCASTER LABORATORIES

PAGE: 8 INSTRUMENT ID: 02360

8 ILP 12/21/92 14:58

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.00255	86.60	0.00510	0.00127	0.00127		
Al	0.02624	29.49	0.03491	0.02382	0.02000		
As	0.00373	229.0	-0.0112	0.00560	-0.0056		
B	0.00139	173.6	0.00140	-0.0027	-0.0028		
Ba	0.00233	0.002	0.00233	0.00233	0.00233		
Be	0.00041	173.3	0.00124	-0.0000	-0.0000		
Ca	0.02504	28.08	0.03316	0.02098	0.02098		
Cd	0.00226	41.66	0.00308	0.00246	0.00123		
Co	0.00008	1757.	0.00101	0.00046	-0.0017		
Cr	0.00211	148.0	0.00461	-0.0013	0.00312		
Cu	0.00222	98.50	0.00465	0.00040	0.00161		
Fe	0.00000	18415	0.00055	-0.0011	0.00056		
K	0.13793	0.006	0.13792	0.13793	0.13793		
Li	0.00702	65.46	0.01203	0.00300	0.00601		
Mg	0.00743	275.2	0.01594	-0.0159	0.02229		
Mn	0.00078	86.49	0.00156	0.00039	0.00039		
Mo	0.01078	47.32	0.01662	0.00853	0.00718		
Na	0.00572	458.7	0.02862	-0.0229	0.01145		
Ni	0.00055	483.8	0.00237	-0.0028	-0.0011		
P	0.00385	1489.	-0.0190	-0.0520	0.05959		
Pb	0.01159	125.8	0.01596	-0.0046	0.02351		
Sb	0.01204	128.3	0.00495	-0.0158	-0.0252		
Se	0.01691	138.0	0.00229	0.00460	0.04383		
Si	0.00271	191.1	0.00221	-0.0081	-0.0022		
Sn	0.00742	103.9	0.01633	0.00296	0.00296		
Sr	0.00037	110.2	0.00081	-0.0000	0.00030		
Ti	0.00006	1549.	0.00033	-0.0010	0.00092		
Tl	0.01792	301.1	-0.0128	-0.0742	0.03336		
V	0.00004	53.58	0.00006	0.00002	0.00003		
Zn	0.00234	21.01	0.00291	0.00206	0.00205		

Name [illegible]5601I11

LANCASTER LABORATORIES

PAGE: 9

INSTRUMENT ID: 02360

[illegible] ICSA 12/21/02 15:03

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].03073	5.649	-0.0325	-0.0305	-0.0290		
Al	[illegible]4.34030	0.856	465.827	467.361	459.832		
As	[illegible]16836	31.15	0.22428	0.16052	0.12026		
B	[illegible].28098	1.998	0.28746	0.27789	0.27758		
Ba	[illegible]00637	21.36	0.00704	0.00562	0.00554		
Be	[illegible].00122	0.203	0.00122	0.00122	0.00122		
Ca	[illegible]1.39697	0.609	462.702	463.319	458.168		
Cd	[illegible]02888	6.605	0.03095	0.02850	0.02719		
Co	[illegible].00050	172.0	0.00007	-0.0014	-0.0000		
Cr	[illegible].01461	16.18	-0.0156	-0.0162	-0.0118		
Cu	[illegible]00214	85.67	0.00214	0.00031	0.00399		
Fe	[illegible]4.78601	0.690	175.252	175.690	173.415		
K	[illegible]17600	24.63	0.15095	0.15098	0.22607		
Li	[illegible].00000	9.569	-0.0000	-0.0000	-0.0000		
Mg	[illegible]7.95496	0.740	479.296	480.646	473.952		
Mn	[illegible].01255	11.63	-0.0118	-0.0142	-0.0115		
Mo	[illegible].04557	9.569	-0.0429	-0.0431	-0.0506		
Na	[illegible]07738	27.94	0.05447	0.08024	0.09744		
Ni	[illegible].02658	10.51	-0.0272	-0.0290	-0.0235		
P	[illegible].60866	2.696	-1.5651	-1.6518	-1.6090		
Pb	[illegible]04145	19.88	0.03884	0.05068	0.03483		
Sb	[illegible].02595	53.30	-0.0408	-0.0134	-0.0236		
Se	[illegible].06640	44.19	-0.0331	-0.0887	-0.0772		
Si	[illegible].00247	259.4	0.00403	-0.0064	-0.0059		
Sn	[illegible]08650	18.61	0.09566	0.09593	0.06791		
Sr	[illegible]01694	0.698	0.01700	0.01680	0.01700		
Ti	[illegible].01176	19.78	0.01084	0.01004	0.01441		
Tl	[illegible].35012	27.61	-0.3716	-0.4342	-0.2444		
V	[illegible]00822	9.891	0.00779	0.00916	0.00772		
Zn	[illegible].00952	8.608	-0.0102	-0.0086	-0.0096		

Run Name: 235601I11

LANCASTER LABORATORIES

PAGE: 1

INSTRUMENT ID: 02360

10 [illegible] 12/21/92 15:07

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.89022	2.403	0.91269	0.87009	0.88788		
Al	466.80548	1.010	468.691	461.438	470.286		
As	0.25473	26.38	0.33073	0.20312	0.23035		
B	0.28801	0.588	0.28619	0.28954	0.28831		
Ba	0.46896	1.303	0.47248	0.46190	0.47249		
Be	0.43229	0.997	0.43478	0.42732	0.43478		
Ca	463.77499	0.795	465.357	459.556	466.411		
Cd	0.86925	0.822	0.87172	0.86119	0.87483		
Co	0.42134	1.099	0.42373	0.41599	0.42428		
Cr	0.43286	1.465	0.43757	0.42565	0.43535		
Cu	0.45164	0.618	0.45326	0.44842	0.45325		
Fe	175.92738	0.900	176.671	174.107	177.002		
K	0.07603	0.153	0.07608	0.07590	0.07611		
Li	0.00100	173.2	-0.0000	0.00300	-0.0000		
Mg	480.51742	0.925	482.393	475.437	483.721		
Mn	0.42713	0.906	0.42939	0.42266	0.42934		
Mo	0.04138	5.850	-0.0406	-0.0440	-0.0394		
Na	0.05129	19.34	0.04556	0.06275	0.04556		
Ni	0.81582	0.790	0.82040	0.80844	0.81861		
P	1.60671	3.170	-1.5480	-1.6393	-1.6328		
Pb	0.93145	3.063	0.96295	0.92407	0.90734		
Sb	0.00132	955.6	-0.0023	-0.0134	0.01183		
Se	-0.04922	51.75	-0.0249	-0.0757	-0.0470		
Si	0.00918	41.91	-0.0055	-0.0088	-0.0132		
Sn	0.09294	1.975	0.09173	0.09505	0.09203		
Sr	0.01725	0.684	0.01739	0.01718	0.01718		
Ti	0.01216	3.267	0.01217	0.01177	0.01256		
Tl	0.45116	21.02	-0.5576	-0.3755	-0.4203		
V	0.45424	1.398	0.45791	0.44691	0.45792		
Zn	0.87687	0.815	0.88016	0.86867	0.88178		

[illegible]

AR318966

Name: [illegible]35601111

LANCASTER LABORATORIES

PAGE: 11 INSTRUMENT ID: 02360

11 [illegible] 12/21/92 15:12

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.01815	8.027	0.01847	0.01656	0.01942		
Al	.12425	27.53	0.11940	0.16064	0.09272		
As	.01449	10.92	0.01357	0.01358	0.01632		
B	-0.00082	5.425	-0.0008	-0.0008	-0.0007		
Ba	.00039	172.4	0.00000	0.00000	0.00116		
Be	.00949	3.662	0.00909	0.00970	0.00969		
Ca	.12692	37.09	0.11643	0.17836	0.08596		
Cd	.00987	12.48	0.00864	0.00987	0.01111		
Co	.09533	2.087	0.09373	0.09756	0.09470		
Cr	.02249	5.960	0.02244	0.02117	0.02385		
Cu	.04890	3.800	0.04728	0.04849	0.05093		
Fe	.04735	39.45	0.03968	0.06864	0.03372		
K	.11288	38.48	0.06272	0.13796	0.13796		
Li	.00000	315.4	-0.0000	0.00000	0.00000		
Mg	.12678	35.64	0.11828	0.17561	0.08644		
Mn	.02857	0.008	0.02857	0.02856	0.02857		
Mo	.00042	315.4	-0.0009	0.00042	0.00177		
Na	-0.02868	34.60	-0.0229	-0.0401	-0.0229		
Ni	.07626	3.510	0.07392	0.07567	0.07918		
P	-0.04217	157.5	-0.1185	0.00204	-0.0099		
Pb	-0.00128	685.8	-0.0036	-0.0086	0.00844		
Sb	.10745	10.32	0.09482	0.11188	0.11565		
Se	.00219	557.4	-0.0024	-0.0070	0.01603		
Si	.00034	2879.	-0.0021	-0.0080	0.01118		
Sn	-0.00295	174.2	-0.0059	-0.0059	0.00298		
Sr	.00021	97.36	-0.0002	-0.0004	-0.0000		
Ti	.00384	59.57	0.00252	0.00252	0.00649		
Tl	-0.00937	499.8	0.03157	0.00079	-0.0604		
V	.09756	4.487	0.09256	0.09937	0.10074		
Zn	.03926	2.566	0.03869	0.03867	0.04043		

[illegible] 118

AR318967

[illegible] 25501I11

LANCASTER LABORATORIES

[illegible]

INSTRUMENT ID: 02360

[illegible] 12/21/92 15:17

ITEM	AVERAGE (ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Al	.47820	1.282	0.47121	0.48074	0.48265		
As	.99098	1.818	23.4016	24.1937	24.1076		
Au	01552	1.786	1.97686	2.04813	2.02156		
B	.97397	2.243	0.94885	0.98861	0.98444		
Ba	.88763	2.144	4.76742	4.95999	4.93548		
Be	.46270	1.921	0.45245	0.46842	0.46722		
Ca	.07770	1.791	48.0626	49.5745	49.5958		
Cd	79399	1.894	4.68912	4.84766	4.84519		
Co	.94942	2.280	0.92464	0.95893	0.96468		
Cr	95539	1.986	4.94933	5.12804	5.11882		
Cu	75705	1.911	4.65271	4.81956	4.79887		
Fe	87431	1.797	4.77329	4.92950	4.92014		
K	.11000	2.048	48.9329	50.5860	50.8118		
[illegible]	1.72941	2.148	24.1775	25.1326	24.9280		
Mg	.68695	1.896	23.1689	23.9698	23.9220		
Mn	86808	1.839	4.76470	4.92211	4.91741		
Mo	1.38794	1.897	23.9537	24.6631	24.6469		
Na	0.54456	1.950	49.1226	51.3147	50.8852		
Ni	62307	1.985	9.10278	9.74331	9.72314		
P	1.57641	1.196	24.2292	24.7104	24.7795		
Pb	4.68483	1.720	24.1962	24.9641	24.8941		
Sb	91661	2.101	4.70914	4.95738	4.99330		
Se	.96747	2.267	1.91612	1.98967	1.99662		
Si	03513	2.240	1.98795	2.05448	2.06786		
Sn	80534	1.482	4.72159	4.83841	4.85402		
Sr	98325	2.140	0.95910	0.99773	0.99291		
Ti	97942	2.231	1.93846	2.00688	2.00291		
Tl	01585	1.807	5.00218	4.91133	5.07303		
V	97658	2.405	1.92238	2.01120	1.99616		
Zn	86081	1.713	4.76514	4.90017	4.91711		

149

AR318968

N [illegible] 601111

LANCASTER LABORATORIES

[illegible] INSTRUMENT ID: 02360

[illegible] 12/21/92 15:21

ITEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	PREREAD	REDIG
Ag	.0318	59.98	0.00510	0.00127	0.00318	_____	_____
Al	.3860	51.08	0.06096	0.02366	0.03118	_____	_____
As	.00701	121.6	-0.0112	0.00280	-0.0126	_____	_____
B	.00211	57.46	-0.0028	-0.0028	-0.0007	_____	_____
Ba	.0466	49.99	0.00700	0.00466	0.00233	_____	_____
Be	.00124	0.591	0.00123	0.00124	0.00124	_____	_____
Ca	.5076	51.58	0.07987	0.04332	0.02910	_____	_____
Cd	.00328	86.64	0.00493	0.00493	-0.0000	_____	_____
Co	.00057	250.4	-0.0019	-0.0007	0.00093	_____	_____
Cr	.00332	19.47	0.00365	0.00258	0.00374	_____	_____
Cu	.0526	41.54	0.00768	0.00465	0.00344	_____	_____
Fe	.0736	61.05	0.01246	0.00566	0.00396	_____	_____
K	.3807	46.20	0.13783	0.13791	0.28842	_____	_____
Li	.01805	66.66	0.03009	0.01805	0.00601	_____	_____
Mg	.4355	47.06	0.05693	0.03505	0.02867	_____	_____
Mn	.0313	94.34	0.00626	0.00274	0.00039	_____	_____
Mo	.2201	62.14	0.03684	0.01932	0.00988	_____	_____
Na	.00286	964.4	0.03867	-0.0143	-0.0228	_____	_____
Ni	.0822	32.96	0.01117	0.00763	0.00585	_____	_____
P	.3555	60.23	0.03282	0.02356	0.06028	_____	_____
Pb	.2587	63.09	0.04468	0.01545	0.01747	_____	_____
Sb	.0434	204.7	0.01437	0.00120	-0.0025	_____	_____
Se	.1306	71.35	0.01844	0.00230	0.01845	_____	_____
Si	.00323	114.6	-0.0027	-0.0066	0.00070	_____	_____
Sn	.0888	104.2	0.01629	0.01186	-0.0014	_____	_____
Sr	.0038	58.07	0.00143	0.00081	0.00040	_____	_____
Ti	.0304	83.83	0.00181	0.00013	0.00410	_____	_____
Tl	.06923	46.38	-0.0589	-0.1052	-0.0434	_____	_____
V	.0280	85.66	0.00557	0.00143	0.00140	_____	_____
Zn	.0495	73.82	0.00904	0.00378	0.00202	_____	_____

AR318969

Cd

[illegible] 25601111

LANCASTER LABORATORIES

[illegible] INSTRUMENT ID: 02360

PBW

[illegible] 12/21/92 15:26

[illegible].DF1,92?541848001.

ELEM	IRAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
[illegible]Ag	00000	44598	0.00128	-0.0025	0.00128		
Al	04134	24.07	0.03759	0.03380	0.05262		
As	.00658	49.15	0.00845	0.00284	0.00844		
B	.00869	31.87	0.01183	0.00660	0.00764		
Ba	00233	0.002	0.00233	0.00233	0.00233		
Be	0.00000	168.4	-0.0000	-0.0000	-0.0000		
Ca	.01421	4.123	0.01489	0.01387	0.01387		
✓Cd	00082	86.70	0.00123	-0.0000	0.00123		
[illegible]Co	0.00286	9.704	-0.0027	-0.0031	-0.0027		
Cr	00040	495.9	-0.0002	-0.0024	0.00153		
[illegible]Cu	.00020	349.9	-0.0002	0.00101	-0.0002		
Fe	.02697	5.468	0.02782	0.02527	0.02782		
K	.13794	0.001	0.13794	0.13794	0.13794		
Li	00200	173.2	0.00000	0.00000	0.00601		
Mg	.01167	83.27	0.00956	0.00318	0.02228		
Mn	00156	75.03	0.00273	0.00156	0.00039		
Mo	00314	42.92	0.00179	0.00449	0.00313		
Na	00000	64168	0.01144	-0.0229	0.01145		
Ni	0.00115	151.5	0.0011	0.00059	-0.0029		
P	05656	79.01	-0.0154	-0.0501	-0.1041		
Pb	.00971	112.1	0.0127	-0.0187	0.00237		
Sb	0.00129	553.6	0.00190	-0.0063	-0.0044		
Se	01498	78.11	0.02767	0.01268	0.00460		
Si	.11601	3.149	0.11774	0.11848	0.11181		
Sn	00297	259.8	0.00743	0.00742	-0.0059		
Sr	00003	756.1	0.00030	-0.0002	-0.0000		
Ti	0.00013	457.0	-0.0006	-0.0002	0.00052		
Tl	03841	197.5	0.0896	-0.0743	0.04874		
V	00091	171.5	0.00090	0.00001	0.00273		
Zr	00235	21.76	0.00194	0.00204	0.00206		

152

AR318970

cd

Name: [illegible]601I11

LANCASTER LABORATORIES

Acct: 15 INSTRUMENT ID: 02360

[illegible] LCSW 12/21/92 15:30

[illegible] 023541848001

ITEM	AGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
[illegible]	4884	3.914	0.05076	0.04885	0.04693		
[illegible]	1084	0.202	1.91307	1.91307	1.90638		
[illegible]	9342	0.773	2.00929	1.99249	1.97848		
[illegible]	3142	0.595	2.03841	2.03841	2.01745		
[illegible]	5254	0.431	1.95955	1.95488	1.94321		
[illegible]	5041	1.214	0.05102	0.05041	0.04980		
[illegible]	0601	0.400	4.01650	4.01396	3.98756		
✓ [illegible]	4779	3.251	0.04635	0.04944	0.04759		
[illegible]	3541	1.109	0.49043	0.48608	0.47973		
[illegible]	1064	1.280	0.21319	0.21090	0.20782		
[illegible]	5174	0.738	0.25376	0.25011	0.25134		
[illegible]	1403	0.629	1.01771	1.01771	1.00666		
[illegible]	2220	1.824	4.12719	4.19743	4.04697		
[illegible]	6021	0.177	1.96772	1.96222	1.95620		
[illegible]	9159	0.389	1.90011	1.88737	1.88730		
[illegible]	0458	0.711	0.50533	0.50771	0.50066		
[illegible]	0537	0.611	1.92098	1.99368	1.97144		
[illegible]	2572	0.480	4.13819	4.13815	4.10381		
[illegible]	0948	0.719	0.51066	0.51242	0.50537		
[illegible]	1354	1.783	2.04972	1.97789	2.01302		
[illegible]	9318	2.780	0.49783	0.50393	0.47774		
[illegible]	8671	1.166	0.49269	0.48607	0.48139		
[illegible]	4646	0.860	2.06181	2.05030	2.02724		
[illegible]	3094	0.518	4.94200	4.94826	4.90156		
[illegible]	0990	0.484	3.89131	3.92921	3.90918		
[illegible]	6540	0.467	0.98892	0.96697	0.96031		
[illegible]	0977	0.421	1.00456	0.99681	0.99781		
[illegible]	2086	4.207	1.96182	1.99207	1.83868		
[illegible]	9491	0.865	0.48856	0.48584	0.48032		
[illegible]	1228	0.523	0.51462	0.51286	0.50935		

154

AR318971

cd

Run Name: [illegible]35601I11

LANCASTER LABORATORIES

PAGE: 15 INSTRUMENT ID: 02360

16 [illegible] JAMMT 12/21/92 15:35
U***,10[illegible].DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	[illegible].00039	280.4	-0.0002	-0.0002	0.00167		
__Al	[illegible]82650	0.961	1.80676	1.84036	1.83238		
__As	[illegible].00013	5126.	-0.0076	0.00078	0.00641		
__B	[illegible].05483	2.225	0.05624	0.05414	0.05411		
__Ba	[illegible]06620	2.035	0.06542	0.06776	0.06542		
__Be	[illegible].00000	113.6	0.00000	-0.0000	-0.0000		
__Ca	[illegible].41562	0.424	17.3636	17.3829	17.5002		
X Cd	[illegible].00017	362.4	0.00044	-0.0007	-0.0001		
__Co	[illegible].00185	148.5	0.00106	-0.0044	-0.0022		
__Cr	[illegible].00048	433.3	0.00021	0.00118	-0.0028		
__Cu	[illegible].00420	16.73	0.00461	0.00339	0.00461		
__Fe	[illegible].15535	0.474	2.14513	2.15535	2.16557		
__K	[illegible].90701	1.470	5.85685	5.85684	6.00733		
__Li	[illegible].00000	146.4	0.00000	0.00000	0.00000		
__Mg	[illegible].04638	0.333	7.05488	7.01984	7.06441		
__Mn	[illegible].10923	1.074	0.10923	0.11041	0.10806		
__Mo	[illegible]00106	146.4	0.00016	0.00016	0.00285		
__Na	[illegible].83610	0.220	9.81317	9.85617	9.83896		
__Ni	[illegible].00076	132.0	-0.0013	-0.0013	0.00040		
__P	[illegible].16364	12.81	0.14898	0.18766	0.15429		
__Pb	[illegible].00091	414.2	-0.0012	-0.0012	0.00528		
__Sb	[illegible].00946	55.40	-0.0103	-0.0038	-0.0141		
__Se	[illegible].00337	978.2	0.00008	-0.0379	0.02777		
__Si	[illegible].26504	0.379	7.24459	7.25412	7.29640		
__Sn	[illegible].00022	4183.	0.00275	0.00720	-0.0106		
__Sr	[illegible].08511	0.367	0.08484	0.08504	0.08545		
__Ti	[illegible]08293	6.971	0.07751	0.08902	0.08227		
__Tl	[illegible].01491	312.7	0.03602	-0.0555	-0.0252		
__V	[illegible]00368	85.17	0.00006	0.00550	0.00550		
__Zn	[illegible]07297	1.201	0.07297	0.07209	0.07384		

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AR318972

Name: [redacted] 601I11

LANCASTER LABORATORIES

PAGE: 17 INSTRUMENT ID: 02360

17 1908[redacted] JAMMT A 12/21/92 15:39
[redacted].100-[redacted] DF1.923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[redacted]5321	6.229	0.05512	0.05512	0.04938		
Al	[redacted]4554	1.113	4.06783	4.07511	3.99367		
As	[redacted]0578	11.10	0.22166	0.21607	0.17960		
B	[redacted]5511	1.336	0.15511	0.15718	0.15304		
Ba	[redacted]4935	0.649	0.54858	0.55324	0.54624		
Be	[redacted]4933	0.009	0.04933	0.04933	0.04934		
Ca	[redacted]8396	0.484	20.6910	20.7804	20.5803		
Cd	[redacted]9049	0.645	0.48967	0.49399	0.48782		
Co	[redacted]9547	1.259	0.09478	0.09686	0.09478		
Cr	[redacted]1608	0.522	0.51507	0.51914	0.51404		
Cu	[redacted]9034	1.032	0.48872	0.49602	0.48629		
Fe	[redacted]4313	0.475	2.44512	2.45362	2.43064		
[redacted]	[redacted].29215	0.730	10.2921	10.3673	10.2169		
[redacted]	[redacted]7384	0.643	2.46782	2.49190	2.46180		
Mg	[redacted]4947	0.373	8.74097	8.78560	8.72184		
Mn	[redacted]9283	0.604	0.59205	0.59674	0.58970		
Mo	[redacted]6933	0.397	2.46619	2.48034	2.46147		
Na	[redacted].13629	0.530	14.1162	14.2193	14.0733		
Ni	[redacted]8273	0.778	0.97923	0.99150	0.97747		
P	[redacted]3229	1.610	2.68102	2.60391	2.61195		
Pb	[redacted]9054	0.471	2.48535	2.50399	2.48229		
Sb	[redacted]8817	1.231	0.48470	0.48470	0.49511		
Se	[redacted]1842	12.18	0.19304	0.21611	0.24612		
Si	[redacted]0324	0.775	6.93040	6.93777	6.84155		
Sn	[redacted]9136	1.383	0.48542	0.49878	0.48988		
Sr	[redacted]7589	0.597	0.17562	0.17705	0.17500		
Ti	[redacted]7335	0.866	0.27229	0.27606	0.27170		
Tl	[redacted]4469	5.356	0.53192	0.52408	0.57808		
V	[redacted]0450	0.781	0.20539	0.20545	0.20265		
Zn	[redacted]6070	0.176	0.56131	0.56124	0.55956		

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Run Name: [illegible]35601I11

LANCASTER LABORATORIES

PAGE: 18

INSTRUMENT ID: 02360

18 [illegible]: JAMMT D 12/21/92 15:44

D***.100 [illegible].DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].00021	1.446	-0.0002	-0.0002	-0.0002		
Al	[illegible]94023	1.335	1.95357	1.95676	1.91038		
As	[illegible].00435	145:4	-0.0017	0.01092	0.00386		
B	[illegible].04538	0.111	0.04535	0.04535	0.04544		
Ba	[illegible].06698	2.012	0.06776	0.06776	0.06543		
Be	[illegible].00081	88.90	-0.0000	0.00123	0.00124		
Ca	[illegible].10709	0.812	18.1851	18.1988	17.9373		
X Cd	[illegible]00084	41.94	0.00105	0.00104	0.00043		
Co	[illegible].00105	164.0	-0.0029	-0.0006	0.00044		
Cr	[illegible]00338	22.05	0.00310	0.00423	0.00281		
Cu	[illegible]00521	34.99	0.00521	0.00704	0.00339		
Fe	33252	0.959	2.34189	2.34871	2.30697		
K	.08258	1.237	6.08258	6.00734	6.15783		
Li	00501	34.64	0.00300	0.00601	0.00601		
Mg	31711	0.676	7.35746	7.33197	7.26191		
Mn	11587	1.169	0.11743	0.11508	0.11509		
Mo	[illegible]00373	75.11	0.00688	0.00148	0.00284		
Na	[illegible].34867	1.506	10.3715	10.4918	10.1825		
Ni	[illegible]00097	103.5	0.00040	0.00038	0.00213		
P	[illegible]11465	63.58	0.19846	0.06599	0.07948		
Pb	[illegible].00876	100.2	-0.0148	0.00131	-0.0128		
Sb	[illegible]00089	369.0	-0.0029	0.00277	0.00280		
Se	[illegible].00261	226.2	0.00238	-0.0010	-0.0091		
Si	[illegible]51897	1.571	7.55333	7.61616	7.38743		
Sn	[illegible].00024	2088.	0.00272	-0.0061	0.00272		
Sr	08805	0.940	0.08853	0.08853	0.08709		
Ti	[illegible]08941	6.974	0.08505	0.09656	0.08664		
Tl	[illegible].00444	1212.	-0.0558	-0.0091	0.05171		
V	00552	49.39	0.00825	0.00551	0.00279		
Zn	07793	1.305	0.07910	0.07733	0.07735		

AR318974

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Name: 9 [illegible]01I11

LANCASTER LABORATORIES

PAGE: 19

INSTRUMENT ID: 02360

ID 190827 JAMMTS 12/21/92 15:48

[illegible].100-1[illegible] 1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible]568	3.198	0.04536	0.04728	0.04441		
Al	[illegible]295	0.360	3.34911	3.37166	3.36808		
As	[illegible]631	0.624	1.94230	1.96193	1.96471		
B	[illegible]789	0.186	1.96406	1.96824	1.97138		
Ba	[illegible]077	0.000	1.96077	1.96077	1.96077		
Be	[illegible]814	1.488	0.04732	0.04857	0.04855		
Ca	[illegible]45162	0.219	19.4040	19.4639	19.4868		
X Cd	[illegible]4705	1.515	0.04746	0.04746	0.04622		
Co	[illegible]7232	1.514	0.47562	0.46411	0.47723		
Cr	[illegible]0620	0.820	0.20426	0.20703	0.20732		
Cu	[illegible]4890	0.002	0.24890	0.24890	0.24889		
Fe	[illegible]851	0.136	2.81482	2.81822	2.82248		
[illegible]	[illegible]329	1.747	8.78774	9.08870	9.01345		
[illegible]	[illegible]3582	0.327	1.83582	1.82980	1.84184		
Mg	8.[illegible]2566	0.239	8.10442	8.14264	8.12993		
Mn	[illegible]8646	0.231	0.58489	0.58724	0.58724		
Mo	[illegible]819	0.423	1.92876	1.94223	1.94358		
Na	[illegible]0863	0.387	12.7800	12.7800	12.8658		
Ni	[illegible]9406	0.412	0.49286	0.49641	0.49290		
P	[illegible]9186	4.045	1.89990	2.02551	2.05016		
Pb	[illegible].7746	0.920	0.47443	0.47545	0.48250		
Sb	[illegible]5295	1.423	0.45670	0.46984	0.46230		
Se	[illegible]8313	1.629	1.98236	2.01582	1.95121		
Si	[illegible]9793	0.198	10.6151	10.6040	10.5745		
Sn	[illegible]3401	0.146	3.83997	3.82881	3.83327		
Sr	[illegible]1903	0.078	1.01862	1.01852	1.01995		
Ti	[illegible].7419	0.232	1.02658	1.02419	1.02181		
Tl	[illegible]4432	5.022	1.85473	1.93130	1.74692		
V	[illegible]7664	0.663	0.47479	0.47484	0.48029		
Zn	[illegible]5893	0.181	0.55834	0.56010	0.55834		

156

AR318975

Run Name: [illegible]35601I11

LANCASTER LABORATORIES

PAGE: 2[illegible]

INSTRUMENT ID: 02360

20 [illegible] JAMMT L 12/21/92 15:53

[illegible]**,100 : [illegible].DF1,923541848001,1,,

ELEM	[illegible]ERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].00055	0.101	-0.0005	-0.0005	-0.0005		
Al	[illegible]48814	1.951	0.48723	0.49809	0.47910		
As	[illegible]00029	3247.	0.00355	-0.0104	0.00777		
B	[illegible]01379	7.638	0.01484	0.01274	0.01377		
Ba	[illegible]01479	9.106	0.01635	0.01401	0.01401		
Be	[illegible]00000	18.54	0.00000	0.00000	0.00000		
Ca	[illegible]55484	0.934	3.51693	3.56872	3.57887		
Cd	[illegible]00078	180.5	-0.0000	-0.0000	0.00243		
Co	[illegible].00023	342.3	0.00017	-0.0011	0.00028		
Cr	[illegible]00191	51.26	0.00185	0.00096	0.00292		
Cu	[illegible]00099	121.9	-0.0002	0.00099	0.00221		
Fe	[illegible]45310	0.868	0.44855	0.45537	0.45537		
K	[illegible]21651	9.448	1.11617	1.19142	1.34193		
Li	[illegible]00000	118.5	0.00000	0.00000	-0.0000		
Mg	[illegible]43858	1.022	1.43009	1.43008	1.45557		
Mn	[illegible]02341	2.892	0.02380	0.02262	0.02380		
Mo	[illegible].00262	118.5	0.00442	0.00442	-0.0009		
Na	[illegible].02166	0.850	2.00447	2.02166	2.03885		
Ni	[illegible].00054	317.3	-0.0011	0.00055	0.00228		
P	[illegible].00337	1047.	-0.0228	-0.0247	0.03747		
Pb	[illegible].00460	236.4	0.00763	0.01367	-0.0074		
Sb	[illegible]00305	215.0	-0.0007	-0.0007	0.01064		
Se	[illegible].00883	208.0	0.01383	0.02421	-0.0115		
Si	[illegible]68643	1.848	1.65263	1.69260	1.71406		
Sn	.00151	294.0	0.00294	-0.0059	-0.0015		
Sr	[illegible]01796	0.658	0.01803	0.01803	0.01782		
Ti	[illegible]02447	7.313	0.02275	0.02433	0.02632		
Tl	[illegible]00348	440.9	-0.0119	0.01883	0.00354		
V	[illegible]00002	38.55	0.00002	0.00002	0.00001		
Zn	[illegible]01605	0.066	0.01606	0.01605	0.01604		

152

AR318976

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ame: [illegible]

LANCASTER LABORATORIES

[illegible]: 21

INSTRUMENT ID: 02360

[illegible] 190835 JAMMF 12/21/92 15:58

[illegible].100-10[illegible] F1.923541848001.1.,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.[illegible]0003	7132.	-0.0025	0.00130	0.00130		
Al	0.[illegible]0606	59.45	0.00265	0.00984	0.00570		
As	-0.[illegible]0198	334.7	0.00036	0.00315	-0.0094		
B	0.[illegible]1692	2.568	0.04625	0.04832	0.04621		
Ba	0.[illegible]1392	1.823	0.07236	0.07470	0.07470		
Be	-0.[illegible]0000	86.99	0.00000	-0.0000	-0.0000		
Ca	18.[illegible]3126	0.596	18.3754	18.4729	18.5953		
X Cd	0.[illegible]1032	216.6	-0.0000	-0.0000	0.00115		
Co	-0.[illegible]0031	542.0	0.00155	-0.0007	-0.0017		
Cr	0.[illegible]0117	43.22	-0.0017	-0.0009	-0.0008		
Cu	0.[illegible]0445	20.90	0.00344	0.00527	0.00466		
Fe	0.[illegible]1813	0.001	0.21814	0.21813	0.21813		
[illegible]	[illegible]1277	1.874	6.00736	6.15786	6.23310		
[illegible]	[illegible]0200	173.2	0.00000	-0.0000	0.00601		
[illegible]	[illegible]0127	0.680	7.29713	7.36082	7.39587		
Mn	0.[illegible]8545	1.585	0.08467	0.08467	0.08702		
Mo	0.[illegible]0133	233.8	0.00312	-0.0022	0.00312		
Na	11.[illegible]8436	2.386	10.8865	11.2645	11.4019		
Ni	0.[illegible]0001	15823	-0.0011	-0.0011	0.00232		
P	0.[illegible]2191	334.4	-0.0288	0.10591	-0.0113		
Pb	-0.[illegible]0805	151.2	-0.0218	-0.0036	0.00134		
Sb	-0.[illegible]0063	1478.	0.00880	-0.0100	-0.0006		
Se	0.[illegible]0171	793.9	-0.0017	-0.0098	0.01671		
Si	5.[illegible]2963	0.584	5.04024	5.06912	5.09951		
Sn	0.[illegible]0129	524.4	0.00723	0.00278	-0.0061		
Sr	0.[illegible]8989	0.732	0.08914	0.09017	0.09037		
Ti	0.[illegible]0277	45.88	0.00370	0.00330	0.00132		
Tl	0.[illegible]3853	184.0	0.00251	-0.1204	0.00228		
V	0.[illegible]0183	85.32	0.00002	0.00273	0.00275		
Zn	0.[illegible]4823	2.096	0.04765	0.04764	0.04940		

158

AR318977

Run Name: [illegible]35601I11

LANCASTER LABORATORIES

PAGE: 22

INSTRUMENT ID: 02360

22 [illegible] JAMFM 12/21/92 16:02
****,103 [illegible].DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].04434	7.778	0.04721	0.04530	0.04051		
Al	[illegible].00681	142.6	0.00220	-0.0171	-0.0055		
As	[illegible].01095	76.81	-0.0109	-0.0193	-0.0025		
B	[illegible].05069	0.009	0.05068	0.05068	0.05069		
Ba	[illegible].04902	0.000	0.04902	0.04902	0.04902		
Be	[illegible].00000	270.7	-0.0000	0.00000	0.00000		
Ca	[illegible].34606	0.222	17.3159	17.3895	17.3327		
X Cd	[illegible].00048	146.6	-0.0000	-0.0000	-0.0013		
Co	[illegible].00124	164.8	-0.0005	0.00034	-0.0035		
Cr	[illegible].00212	144.8	0.00128	-0.0047	-0.0029		
Cu	[illegible].00303	46.33	0.00466	0.00222	0.00222		
Fe	[illegible].15626	0.629	0.15683	0.15512	0.15683		
K	63112	2.672	5.78161	5.48062	5.63112		
Li	[illegible]00000	75.51	0.00000	0.00000	0.00000		
Mg	[illegible]92447	0.200	6.93084	6.93402	6.90854		
Mn	.07999	0.000	0.07999	0.07999	0.07999		
Mo	[illegible]00178	75.51	0.00043	0.00313	0.00178		
Na	[illegible].69185	0.334	10.7319	10.6632	10.6804		
Ni	[illegible]00116	174.4	-0.0011	0.00232	0.00232		
P	[illegible]01383	211.3	0.02560	-0.0194	0.03535		
Pb	[illegible].00235	150.3	0.00134	-0.0026	-0.0057		
Sb	[illegible].00062	1885.	-0.0101	-0.0043	0.01261		
Se	[illegible]00976	108.3	0.01206	0.01899	-0.0017		
Si	[illegible]76008	0.490	4.73737	4.78403	4.75883		
Sn	[illegible]01023	25.16	0.01172	0.01171	0.00725		
Sr	[illegible]08395	0.508	0.08361	0.08443	0.08382		
Ti	[illegible]00238	63.07	0.00132	0.00172	0.00410		
Tl	[illegible].06418	136.1	0.03313	-0.0898	-0.1358		
V	[illegible]00001	16.38	0.00002	0.00001	0.00001		
Zn	[illegible]04414	0.002	0.04414	0.04414	0.04414		

159

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Name: [redacted]601111

LANCASTER LABORATORIES

PAGE: 23

INSTRUMENT ID: 02360

[redacted] ZZZZZZ

[redacted] 12/21/92 16:07

****.100-[redacted] F1.923541848001,1,,

ELEM	AVERAGE (ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[redacted]0318	0.020	0.00318	0.00318	0.00318		
Al	[redacted]3078	18.61	0.03197	0.03583	0.02455		
As	[redacted]0232	363.3	0.00561	-0.0013	-0.0112		
B	[redacted]1257	9.570	0.01395	0.01187	0.01187		
Ba	[redacted]0077	173.1	0.00000	0.00000	0.00233		
Be	[redacted]0000	115.1	-0.0000	0.00001	0.00000		
Ca	[redacted]7735	10.60	0.19867	0.17024	0.16313		
X Cd	[redacted]0000	30076	0.00123	0.00000	-0.0012		
Co	[redacted]0376	31.89	-0.0029	-0.0031	-0.0051		
Cr	[redacted]0085	325.3	0.00380	0.00049	-0.0017		
Cu	[redacted]0020	917.1	-0.0002	-0.0014	0.00223		
Fe	[redacted]2056	300.6	0.00226	0.00056	-0.0011		
K	[redacted]6336	16.49	0.21319	0.28844	0.28844		
[redacted]	[redacted]0000	301.3	0.00000	0.00000	-0.0000		
[redacted]	[redacted]441	60.25	0.04140	0.01592	0.01591		
Mn	[redacted]0078	86.68	0.00039	0.00156	0.00039		
Mo	[redacted]0044	301.3	0.00179	0.00044	-0.0009		
Na	[redacted]7797	5.248	0.38942	0.35506	0.38942		
Ni	[redacted]0407	24.84	-0.0029	-0.0046	-0.0046		
P	[redacted]3214	183.8	-0.0763	0.03498	-0.0550		
Pb	[redacted]1072	46.96	-0.0107	-0.0056	-0.0157		
Sb	[redacted]0951	46.05	-0.0120	-0.0120	-0.0044		
Se	[redacted]230	200.1	-0.0023	0.00230	0.00692		
Si	[redacted]1895	0.433	0.65019	0.65093	0.64573		
Sn	[redacted]0593	198.6	-0.0148	-0.0103	0.00743		
Sr	[redacted]0099	21.52	0.00122	0.00081	0.00092		
Ti	[redacted]105	57.30	0.00092	0.00052	0.00171		
Tl	[redacted]2033	99.79	-0.0280	0.00268	-0.0356		
V	[redacted]0181	114.9	0.00001	-0.0040	-0.0013		
Zn	[redacted]148	69.25	0.00206	0.00207	0.00029		

160

AR318979

File Name: 25601I11

LANCASTER LABORATORIES

PAGE: 21 INSTRUMENT ID: 02360

24 OF 12/21/92 16:11

ELEM	AVERAGE (ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.47437	0.930	0.46977	0.47692	0.47692		
Al	23.78264	0.296	23.7271	23.8620	23.7587		
As	2.01223	0.928	2.02442	2.02154	1.99072		
B	.97192	0.215	0.96983	0.97402	0.97189		
Ba	4.86312	0.386	4.85729	4.88413	4.84795		
Be	.46146	0.406	0.45982	0.46350	0.46104		
Ca	48.88037	0.382	48.6855	49.0582	48.8973		
Cd	4.80262	0.610	4.77299	4.83162	4.80325		
Co	.93978	0.923	0.93685	0.93294	0.94954		
Cr	5.05818	0.360	5.03894	5.07522	5.06039		
Cu	4.74220	0.440	4.73330	4.76607	4.72723		
Fe	4.85386	0.323	4.84197	4.87169	4.84793		
K	49.49541	0.075	49.4579	49.5328	49.4955		
Li	24.84477	0.430	24.8497	24.9491	24.7354		
Mg	23.70277	0.252	23.6665	23.7719	23.6698		
Mn	.85595	0.328	4.84108	4.87280	4.85398		
Mo	24.42637	0.390	24.3875	24.5351	24.3565		
Na	51.14303	0.197	51.1682	51.2289	51.0313		
Ni	9.60133	0.324	9.56771	9.62932	9.60696		
P	24.66125	0.575	24.4983	24.7271	24.7583		
Pb	24.66882	0.237	24.6305	24.7363	24.6396		
Sb	4.90491	0.639	4.86904	4.92714	4.91856		
Se	.96751	0.339	1.97137	1.95979	1.97136		
Si	2.11806	1.049	2.09229	2.13084	2.13094		
Sn	4.79336	0.907	4.75179	4.83853	4.78975		
Sr	.97761	0.346	0.97642	0.98144	0.97498		
Ti	1.98788	0.430	1.98344	1.99774	1.98246		
Tl	4.96069	0.799	4.93876	5.00649	4.93720		
V	1.97351	0.566	1.96245	1.98480	1.97328		
Zn	.85879	0.326	4.84171	4.87516	4.85599		

161

AR318980

Name: [redacted] 501111

LANCASTER LABORATORIES

PAGE: 25 INSTRUMENT ID: 02360

[redacted] CCB 12/21/92 16:16

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible]0318	60.04	0.00510	0.00127	0.00318		
Al	[illegible]2846	54.24	0.04575	0.02354	0.01608		
As	-0.00468	105.2	-0.0098	-0.0042	0.00000		
B	[illegible]0279	43.01	0.00349	0.00140	0.00348		
Ba	[illegible]0427	78.73	0.00816	0.00233	0.00233		
Be	[illegible]0082	86.62	0.00123	0.00123	-0.0000		
Ca	[illegible]4704	79.48	0.09003	0.02910	0.02199		
Cd	[illegible]0452	103.2	0.00987	0.00123	0.00246		
Co	[illegible]00000	30270	0.00112	-0.0020	0.00093		
Cr	[illegible]0485	71.57	0.00819	0.00126	0.00509		
Cu	[illegible]0546	33.90	0.00707	0.00343	0.00587		
Fe	[illegible]0339	153.1	0.00906	-0.0011	0.00226		
[redacted]	[illegible]8807	23.09	0.21312	0.13792	0.21318		
[redacted]	[illegible]2006	96.43	0.04213	0.01203	0.00601		
Mg	[illegible]1808	73.35	0.02236	0.00320	0.02867		
Mn	[illegible]0352	101.8	0.00744	0.00274	0.00039		
Mo	[illegible]2336	76.32	0.04358	0.01662	0.00988		
Na	[illegible]2290	229.1	0.08015	-0.0229	0.01147		
Ni	[illegible]0880	50.29	0.01294	0.00413	0.00935		
P	[illegible]6547	61.69	0.05588	0.03074	0.10981		
Pb	[illegible]2284	70.34	0.04065	0.01847	0.00940		
Sb	[illegible]00007	2733.	0.00109	-0.0025	0.00124		
Se	[illegible]1460	134.3	0.01382	-0.0046	0.03460		
Si	[illegible]0541	100.4	0.00741	-0.0007	0.00957		
Sn	[illegible]0145	467.7	0.00291	-0.0059	0.00741		
Sr	[illegible]0068	96.44	0.00143	0.00040	0.00020		
Ti	[illegible]0244	86.66	0.00132	0.00112	0.00489		
Tl	[illegible]00781	1502.	0.01760	-0.1357	0.09473		
V	[illegible]0371	85.61	0.00560	0.00550	0.00004		
Zn	[illegible]0436	93.06	0.00904	0.00203	0.00200		

Run Name [illegible]35601I11

LANCASTER LABORATORIES

PAGE: 26 INSTRUMENT ID: 02360

26 [illegible] JAMLT 12/21/92 16:20
****,100 [illegible],DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	[illegible].00095	230.8	0.00159	-0.0022	-0.0022		
__Al	[illegible]91832	0.588	0.91777	0.92398	0.91321		
__As	[illegible].00832	58.41	-0.0111	-0.0111	-0.0027		
__B	[illegible].03726	3.357	0.03803	0.03581	0.03793		
__Ba	[illegible].05607	0.001	0.05607	0.05607	0.05607		
__Be	[illegible]00040	179.6	-0.0000	0.00124	-0.0000		
__Ca	[illegible].55151	1.014	17.3464	17.6667	17.6413		
X Cd	[illegible].00045	136.4	0.00107	-0.0001	0.00045		
__Co	[illegible].00190	131.7	0.00040	-0.0045	-0.0015		
__Cr	[illegible]00022	1299.	-0.0018	-0.0011	0.00365		
__Cu	[illegible].00361	9.699	0.00401	0.00341	0.00341		
__Fe	[illegible].73253	1.184	1.70897	1.74644	1.74218		
__K	[illegible].35520	0.811	5.33012	5.33013	5.40537		
__Li	[illegible]00000	40.41	0.00000	0.00000	0.00000		
__Mg	[illegible]02522	1.081	6.94028	7.08679	7.04858		
__Mn	[illegible]10730	1.259	0.10574	0.10808	0.10808		
__Mo	[illegible].00385	40.41	0.00565	0.00295	0.00295		
__Na	[illegible].83303	0.825	9.77004	9.92466	9.80440		
__Ni	[illegible]00161	62.33	0.00045	0.00219	0.00219		
__P	[illegible]08489	65.67	0.10693	0.12626	0.02149		
__Pb	[illegible]00264	177.1	0.00784	-0.0012	0.00129		
__Sb	[illegible]00543	80.45	0.01049	0.00292	0.00290		
__Se	[illegible]00566	338.9	0.02103	0.01182	-0.0158		
__Si	[illegible].02983	1.073	5.96170	6.09055	6.03724		
__Sn	[illegible].00859	29.92	0.00711	0.00710	0.01156		
__Sr	[illegible]08340	0.885	0.08259	0.08402	0.08361		
__Ti	[illegible]04180	0.948	0.04219	0.04140	0.04180		
__Tl	[illegible].08830	79.74	-0.1036	-0.1498	-0.0114		
__V	[illegible]00460	34.20	0.00551	0.00278	0.00551		
__Zn	[illegible]07746	0.008	0.07747	0.07746	0.07746		

AR318982 163

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Name: [redacted]601111

LANCASTER LABORATORIES

AGE: 27 INSTRUMENT ID: 02360

[redacted] JAMLF 12/21/92 16:25
[redacted].100-[redacted] F1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[redacted]1002	10532	0.00129	0.00129	-0.0025	_____	_____
Al	[redacted]2351	32.12	0.02345	0.03108	0.01598	_____	_____
As	[redacted]0066	245.8	-0.0025	0.00025	0.00030	_____	_____
B	[redacted]8325	1.254	0.08221	0.08430	0.08325	_____	_____
Ba	[redacted]3968	0.000	0.03968	0.03968	0.03968	_____	_____
Be	[redacted]0000	616.1	-0.0000	-0.0000	0.00001	_____	_____
Ca	[redacted]7648	0.561	16.9661	17.1245	17.1387	_____	_____
X Cd	[redacted]0075	94.85	0.00116	-0.0000	0.00116	_____	_____
Co	[redacted]0085	232.4	-0.0029	0.00101	-0.0006	_____	_____
Cr	[redacted]0109	293.4	0.00166	-0.0003	-0.0045	_____	_____
Cu	[redacted]547	12.83	0.00588	0.00587	0.00466	_____	_____
Fe	[redacted]251	0.505	0.16365	0.16194	0.16195	_____	_____
[redacted]	[redacted]1726	1.249	5.17964	5.17963	5.29251	_____	_____
[redacted]	[redacted]0000	108.8	-0.0000	0.00000	0.00000	_____	_____
Mg	[redacted]4932	0.249	6.73021	6.76207	6.75569	_____	_____
Mn	[redacted]8234	0.001	0.08234	0.08234	0.08234	_____	_____
Mo	[redacted]0155	108.8	-0.0002	0.00313	0.00178	_____	_____
Na	[redacted]7249	0.570	9.63241	9.73547	9.64958	_____	_____
Ni	[redacted]0232	0.100	0.00232	0.00232	0.00232	_____	_____
P	[redacted]6703	57.02	0.04790	0.11106	0.04215	_____	_____
Pb	[redacted]0569	98.52	-0.0067	0.00035	-0.0107	_____	_____
Sb	[redacted]0410	265.2	0.00314	0.01545	-0.0062	_____	_____
Se	[redacted]0973	185.0	0.03050	-0.0017	0.00051	_____	_____
Si	[redacted]5783	0.266	4.85436	4.84698	4.87215	_____	_____
Sn	[redacted]0132	968.8	-0.0061	0.01618	-0.0061	_____	_____
Sr	[redacted]7978	0.535	0.07931	0.07992	0.08013	_____	_____
Ti	[redacted]0105	184.7	0.00330	0.00013	-0.0002	_____	_____
Tl	[redacted]1038	62.87	-0.0436	-0.1053	-0.1821	_____	_____
V	[redacted]0092	610.1	0.00546	0.00275	-0.0054	_____	_____
Zn	[redacted]1060	0.011	0.04060	0.04059	0.04060	_____	_____

Run Name: [illegible]25601I11

LANCASTER LABORATORIES

PAGE: 26 INSTRUMENT ID: 02360

28 [illegible] CHULT 12/21/92 16:30
****,[illegible],DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].00018	1022.	-0.0020	0.00172	-0.0001		
Al	[illegible] 49558	0.398	1.48898	1.50058	1.49718		
As	[illegible] 00217	260.2	0.00125	-0.0029	0.00826		
B	[illegible] 11457	0.922	0.11350	0.11458	0.11562		
Ba	[illegible].05104	1.319	0.05026	0.05142	0.05142		
Be	[illegible].00001	26.64	-0.0000	-0.0000	-0.0000		
Ca	[illegible].28190	0.300	10.3022	10.2463	10.2971		
Cd	[illegible].00037	339.8	0.00106	-0.0014	-0.0007		
Co	[illegible].00014	197.5	0.00000	-0.0000	0.00048		
Cr	[illegible].00162	121.9	0.00070	0.00027	0.00390		
Cu	[illegible].00238	64.21	0.00218	0.00400	0.00096		
Fe	[illegible].43121	0.368	2.44001	2.42212	2.43149		
K	[illegible].04758	1.425	3.07266	2.99741	3.07267		
Li	00401	86.60	0.00000	0.00601	0.00601		
Mg	[illegible] 50408	0.186	4.51043	4.50726	4.49453		
Mn	[illegible].14920	0.786	0.14802	0.14920	0.15037		
Mo	[illegible].00052	330.6	0.00017	0.00287	-0.0011		
Na	[illegible].46086	1.271	9.58112	9.46085	9.34060		
Ni	[illegible].00079	255.6	-0.0031	0.00038	0.00036		
P	[illegible].12808	32.15	0.16572	0.13441	0.08409		
Pb	[illegible] 00242	145.8	0.00208	0.00612	-0.0009		
Sb	[illegible].00788	113.3	-0.0047	-0.0009	-0.0179		
Se	[illegible].00145	1078.	0.01376	0.00683	-0.0162		
Si	[illegible].72424	0.429	6.75238	6.72573	6.69461		
Sn	[illegible].00556	120.1	0.01160	-0.0017	0.00715		
Sr	[illegible].05953	0.620	0.05994	0.05922	0.05943		
Ti	[illegible] 05449	3.173	0.05568	0.05251	0.05529		
Tl	[illegible].03152	225.1	-0.0723	-0.0726	0.05043		
V	[illegible].00643	24.35	0.00824	0.00553	0.00552		
Zn	[illegible].03450	2.946	0.03509	0.03508	0.03333		

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ame: [illegible] 601111

LANCASTER LABORATORIES

[illegible] E: 29 INSTRUMENT ID: 02360

[illegible] 1908358 CHULF 12/21/92 16:34
[illegible].100-100 [illegible]1.923541848001,1,.

[illegible] EM	A[illegible] AGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible]0090	265.1	0.00132	-0.0034	-0.0005		
Al	0.[illegible]329	23.11	0.03965	0.03548	0.02474		
As	0.00225	186.3	0.00194	-0.0022	-0.0064		
B	0.[illegible]805	7.319	0.03910	0.03490	0.04017		
Ba	0.[illegible]035	0.000	0.03035	0.03035	0.03035		
Be	[illegible]00000	64.73	-0.0000	-0.0000	-0.0000		
Ca	0.[illegible]052	0.789	9.30206	9.38786	9.24164		
X Cd	0.[illegible]014	1050.	0.00179	-0.0000	-0.0012		
Co	[illegible]0019	1352.	0.00251	-0.0026	-0.0003		
Cr	[illegible]0137	104.4	-0.0009	-0.0029	-0.0002		
Cu	[illegible]141	99.46	0.00222	-0.0002	0.00222		
Fe	[illegible]508	1.090	0.32394	0.32905	0.32224		
[illegible]	[illegible]43	2.790	2.77168	2.69643	2.62119		
[illegible]	[illegible]0000	317.1	0.00000	-0.0000	0.00000		
[illegible]a	[illegible]769	0.533	3.98831	3.99151	3.95326		
Mn	0.[illegible]762	1.994	0.11762	0.11997	0.11527		
Mo	0.[illegible]042	317.1	0.00177	-0.0009	0.00042		
Na	[illegible].8748	1.370	8.75621	8.75619	8.55004		
Ni	0.[illegible]055	829.4	0.00405	-0.0046	0.00230		
P	[illegible]4517	110.0	0.00529	-0.0466	-0.0941		
Pb	0.[illegible]538	178.3	0.01445	-0.0046	0.00639		
Sb	[illegible]0252	389.5	0.00884	-0.0082	-0.0081		
Se	[illegible]0126	1566.	0.01641	-0.0228	0.00257		
Si	4.[illegible]886	0.675	4.09748	4.14048	4.08862		
Sn	[illegible]0012	1991.	-0.0016	-0.0016	0.00284		
Sr	[illegible]335	0.221	0.05328	0.05348	0.05328		
Ti	[illegible]053	448.3	0.00291	-0.0018	0.00053		
Tl	[illegible]6418	49.84	-0.0283	-0.0897	-0.0744		
V	[illegible]002	20.79	0.00003	0.00002	0.00003		
Zn	[illegible]598	0.113	0.01597	0.01601	0.01598		

166

AR318985

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Run Name: 9235601I11

LANCASTER LABORATORIES

PAGE: 20

INSTRUMENT ID: 02360

ID 1008081 JSBHT 12/21/92 16:39

****.100-100.DF1.923541848001,1,,

ITEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	-0.00087	335.3	-0.0002	-0.0040	0.00168		
Al	1.41060	0.139	1.40833	1.41181	1.41166		
As	0.00083	1346.	0.01203	-0.0103	0.00083		
B	0.04570	0.060	0.04573	0.04567	0.04571		
Ba	0.05842	0.000	0.05842	0.05842	0.05842		
Be	0.00040	180.1	0.00123	-0.0000	-0.0000		
Ca	17.96731	0.202	17.9287	18.0008	17.9723		
X Cd	-0.00100	70.80	-0.0014	-0.0014	-0.0001		
Co	-0.00187	147.6	-0.0035	0.00131	-0.0034		
Cr	0.00115	69.64	0.00128	0.00188	0.00029		
Cu	0.00603	25.36	0.00766	0.00461	0.00583		
Fe	2.18376	0.390	2.17495	2.19199	2.18433		
K	5.27995	0.822	5.25487	5.25488	5.33012		
Li	0.00000	103.6	0.00000	0.00000	0.00000		
Mg	7.04632	0.514	7.01660	7.08666	7.03569		
Mn	0.10219	1.149	0.10337	0.10219	0.10102		
Mo	0.00198	103.6	0.00423	0.00018	0.00153		
Na	9.95343	1.395	9.80453	10.0794	9.97632		
Ni	0.00098	271.3	0.00041	-0.0013	0.00390		
P	0.15714	9.629	0.17437	0.15102	0.14603		
Pb	-0.00634	92.99	-0.0019	-0.0130	-0.0039		
Sb	0.00283	176.6	-0.0009	-0.0085	0.00097		
Se	0.00632	242.9	0.00247	-0.0067	0.02324		
Si	6.52764	0.330	6.50493	6.53010	6.54790		
Sn	0.00413	164.5	0.01156	-0.0018	0.00265		
Sr	0.08641	0.362	0.08607	0.08668	0.08648		
Ti	0.06878	2.642	0.06918	0.07037	0.06680		
Tl	0.04874	78.67	-0.0102	-0.0870	-0.0489		
V	0.00596	13.09	0.00552	0.00686	0.00551		
Zn	0.07738	0.011	0.07738	0.07739	0.07737		

102

AR318986

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Name: [redacted]5601I11

LANCASTER LABORATORIES

PAGE: 31 INSTRUMENT ID: 02360

[redacted] JSBFL 12/21/92 16:43
****,100-[redacted] DF1.923541848001,1,,

ELEM	A	RAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag		1001	6328.	0.00129	-0.0006	-0.0006		
Al		00876	67.52	-0.0037	-0.0071	-0.0153		
As		00351	121.6	0.00020	-0.0081	-0.0025		
B		3678	3.305	0.03818	0.03610	0.03606		
Ba		4201	0.000	0.04201	0.04201	0.04201		
Be		00000	102.2	-0.0000	-0.0000	0.00000		
Ca		54865	0.418	17.6601	17.5697	17.7160		
X Cd		00007	0.448	-0.0000	-0.0000	-0.0000		
Co		00072	267.5	0.00045	0.00032	-0.0029		
Cr		00114	285.2	0.00076	0.00072	-0.0049		
Cu		0324	75.80	0.00466	0.00466	0.00040		
Fe		1009	1.265	0.14151	0.13810	0.14066		
[redacted]		2948	0.846	5.10439	5.17964	5.10440		
[redacted]i		00200	173.2	-0.0000	0.00000	-0.0060		
Mg		0017	0.655	6.78107	6.76833	6.85112		
Mn		6865	0.988	0.06943	0.06826	0.06825		
Mo		00135	151.4	-0.0009	0.00043	-0.0036		
Na		2403	0.713	9.73548	9.64960	9.78702		
Ni		00030	864.2	-0.0003	0.00232	-0.0029		
P		1681	338.5	0.07867	0.00508	-0.0333		
Pb		00101	1905.	0.01947	-0.0036	-0.0188		
Sb		00314	308.4	0.00504	-0.0006	-0.0138		
Se		0591	340.9	0.00052	0.02821	-0.0110		
Si		4571	0.667	4.94399	4.91360	4.97955		
Sn		00016	1597.	-0.0016	0.00280	-0.0016		
Sr		8429	0.853	0.08422	0.08361	0.08504		
Ti		0238	100.4	0.00211	0.00489	0.00013		
Tl		04107	167.7	-0.0052	0.00251	-0.1205		
V		0272	100.0	0.00545	0.00273	-0.0000		
Zn		3591	2.809	0.03531	0.03707	0.03534		

168

AR318987

Rd

[illegible] 35601I11

LANCASTER LABORATORIES

[illegible] INSTRUMENT ID: 02360

[illegible] CRCH- 12/21/92 16:48

[illegible] ,0F1,923541848001,1,,

ITEM	RAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
[illegible]	00004	3131.	0.00164	-0.0012	-0.0002		
[illegible]	5893	0.689	0.95134	0.96202	0.96343		
[illegible]	00048	1765.	0.00805	-0.0023	-0.0079		
[illegible]	3244	6.417	0.03455	0.03240	0.03038		
[illegible]	4596	2.932	0.04674	0.04674	0.04441		
[illegible]	00001	16.73	-0.0000	-0.0000	-0.0000		
[illegible]	45451	0.737	10.4601	10.5286	10.3748		
X [illegible]	0108	113.7	0.00108	0.00231	-0.0001		
[illegible]	00315	37.00	-0.0042	-0.0033	-0.0018		
[illegible]	00097	84.12	-0.0009	-0.0018	-0.0001		
[illegible]	00808	8.692	0.00889	0.00767	0.00767		
[illegible]	5276	1.093	1.95106	1.97490	1.93231		
[illegible]	2184	1.539	2.84692	2.84693	2.77168		
[illegible]	0000	126.7	0.00000	0.00000	-0.0000		
[illegible]	3597	0.529	4.43747	4.46252	4.41793		
[illegible]	1832	1.144	0.11753	0.11988	0.11753		
[illegible]	0092	126.7	0.00159	0.00159	-0.0004		
[illegible]	95256	0.993	10.9134	11.0756	10.8695		
[illegible]	0042	0.970	0.00041	0.00042	0.00042		
[illegible]	8828	72.81	0.11193	0.01667	0.10715		
[illegible]	20119	1211.	-0.0052	0.01491	-0.0132		
[illegible]	01284	22.53	-0.0160	-0.0103	-0.0122		
[illegible]	00610	121.4	-0.0114	0.00235	-0.0091		
[illegible]	2932	1.204	5.32449	5.39079	5.26267		
[illegible]	9416	61.64	0.00267	0.00713	0.00269		
[illegible]	6318	1.138	0.06312	0.06394	0.06250		
[illegible]	1062	2.973	0.02187	0.03029	0.03029		
[illegible]	04273	36.06	-0.0273	-0.0581	-0.0427		
[illegible]	0505	15.58	0.00561	0.00551	0.00414		
[illegible]	2984	0.014	0.02984	0.02984	0.02985		

169

AR318988

Name: 5601I11

LANCASTER LABORATORIES

AGE: 33

INSTRUMENT ID: 02360

33 1908 CRCHF 12/21/92 16:52
****.100-1 .DF1,923541848001,1,,

ELEM	RAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0961	22.97	0.00706	0.01089	0.01088		
Al	04974	21.02	0.03838	0.05897	0.05186		
As	.00228	212.7	0.00052	-0.0078	0.00051		
B	02709	4.493	0.02638	0.02639	0.02849		
Ba	03268	0.000	0.03268	0.03268	0.03268		
Be	00000	577.9	0.00000	0.00000	-0.0000		
Ca	31134	0.706	10.3656	10.3397	10.2285		
X Cd	00014	645.0	-0.0006	-0.0000	0.00117		
Co	.00426	59.47	-0.0071	-0.0033	-0.0023		
Cr	00222	183.8	-0.0053	-0.0037	0.00240		
Cu	00060	421.2	-0.0014	-0.0002	0.00344		
Fe	31144	0.835	0.31201	0.31371	0.30860		
K	29677	1.553	2.77169	2.77169	2.84693		
Li	00000	164.0	-0.0000	-0.0000	0.00000		
Mg	13960	0.235	4.35127	4.33536	4.33216		
Mn	10509	1.288	0.10587	0.10587	0.10353		
Mo	00047	164.0	-0.0009	-0.0009	0.00042		
Na	99537	0.786	11.0068	11.0755	10.9037		
Ni	.00118	0.267	-0.0011	-0.0011	-0.0011		
P	06487	67.18	0.06134	0.11011	0.02315		
Pb	00702	125.3	-0.0167	-0.0046	0.00036		
Sb	.00443	112.9	-0.0082	0.00125	-0.0063		
Se	02106	47.74	0.02568	0.02798	0.00952		
Si	18583	0.469	4.39300	4.40187	4.36262		
Sn	01324	19.43	0.01175	0.01176	0.01621		
Sr	06165	0.583	0.06168	0.06199	0.06127		
Ti	00092	74.08	0.00013	0.00132	0.00132		
Tl	.13073	52.97	-0.0590	-0.1972	-0.1358		
V	00002	54.03	0.00001	0.00001	0.00003		
Zn	01598	0.068	0.01599	0.01598	0.01597		

170

AR318989

Run Name: [illegible]35601I11

LANCASTER LABORATORIES

PAGE: 34

INSTRUMENT ID: 02360

34 [illegible] CRTFB 12/21/92 16:57
[illegible].DF1,923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].00255	43.26	0.00318	0.00127	0.00318		
Al	[illegible].03118	30.63	0.04130	0.02232	0.02993		
As	[illegible].00187	228.7	0.00561	-0.0027	0.00280		
B	[illegible].01187	17.65	0.01396	0.00977	0.01187		
Ba	[illegible].00077	173.1	0.00233	0.00000	0.00000		
Be	[illegible].00000	517.8	0.00000	-0.0000	0.00000		
Ca	[illegible].16855	1.840	0.16922	0.16516	0.17126		
X Cd	[illegible].00000	51.79	0.00000	0.00000	0.00000		
Co	[illegible].00133	336.3	0.00262	-0.0062	-0.0004		
Cr	[illegible].00013	537.2	-0.0007	0.00065	-0.0003		
Cu	00162	135.2	0.00223	-0.0008	0.00345		
Fe	00198	24.77	0.00227	0.00227	0.00141		
K	.18811	23.09	0.21319	0.21319	0.13795		
Li	[illegible].00000	456.1	-0.0000	0.00000	-0.0000		
Mg	01485	24.82	0.01272	0.01272	0.01911		
Mn	[illegible].00078	149.8	-0.0019	-0.0007	0.00039		
Mo	[illegible].00045	456.1	-0.0009	0.00179	-0.0022		
Na	[illegible]37224	4.616	0.38943	0.37222	0.35506		
Ni	00058	601.2	0.00057	0.00409	-0.0029		
P	[illegible].01704	83.35	-0.0253	-0.0006	-0.0250		
Pb	[illegible].01207	91.88	-0.0228	-0.0127	-0.0006		
Sb	[illegible].01008	67.71	-0.0043	-0.0176	-0.0082		
Se	01537	97.59	0.02998	-0.0000	0.01614		
Si	84619	0.971	0.83752	0.85388	0.84717		
Sn	00075	1361.	-0.0014	-0.0081	0.01189		
Sr	00119	13.09	0.00133	0.00102	0.00122		
Ti	[illegible]00039	650.5	0.00290	-0.0022	0.00052		
Tl	[illegible].02050	209.5	-0.0126	-0.0668	0.01801		
V	[illegible].00000	446.8	-0.0000	0.00000	-0.0000		
Zn	[illegible]00293	30.11	0.00292	0.00205	0.00381		

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Name: [illegible]601I11

LANCASTER LABORATORIES

PAGE: 35 INSTRUMENT ID: 02360

[illegible] JSBMT 12/21/92 17:02

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ELEM	A	RAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag		00207	0.179	-0.0020	-0.0020	-0.0020		
Al		34292	0.270	1.34033	1.34136	1.34707		
As		00971	76.72	-0.0153	-0.0012	-0.0125		
B		02941	3.569	0.02947	0.03044	0.02834		
Ba		06038	1.115	0.06077	0.05960	0.06077		
Be		.00001	61.61	-0.0000	-0.0000	-0.0000		
Ca		.90358	0.757	17.7566	18.0236	17.9307		
X Cd		00039	155.7	-0.0002	0.00039	0.00101		
Co		.00183	115.6	-0.0000	-0.0012	-0.0041		
Cr		00111	335.8	-0.0030	0.00233	0.00409		
Cu		00338	35.91	0.00217	0.00460	0.00338		
Fe		28072	0.781	2.55744	2.59321	2.59151		
		03242	0.720	6.08259	6.00734	6.00733		
Li		.00000	278.1	-0.0000	-0.0000	0.00000		
Mg		75811	0.709	7.00077	7.08041	7.09314		
Mn		12642	1.416	0.12447	0.12799	0.12682		
Mo		.00073	278.1	-0.0025	-0.0011	0.00150		
Na		58131	0.948	9.47822	9.65003	9.61567		
Ni		.00051	452.4	-0.0013	0.00210	-0.0022		
P		16435	20.44	0.16279	0.19871	0.13156		
Pb		.00738	77.52	-0.0090	-0.0010	-0.0120		
Sb		00794	83.39	-0.0142	-0.0010	-0.0086		
Se		00555	145.8	0.00478	-0.0021	0.01401		
Si		72430	0.693	6.17592	6.25885	6.23812		
Sn		00550	123.6	-0.0019	0.01144	0.00698		
Sr		08532	0.909	0.08443	0.08586	0.08566		
Ti		09206	1.316	0.09100	0.09180	0.09338		
Tl		06019	102.1	-0.0551	-0.0013	-0.1240		
V		00551	49.46	0.00549	0.00279	0.00824		
Zn		08971	0.015	0.08972	0.08970	0.08971		

Run Name: 935601I11

LANCASTER LABORATORIES

PAGE: 36

INSTRUMENT ID: 02360

36 CCV 12/21/92 17:06

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
__Ag	.47310	1.212	0.46736	0.47311	0.47884		
__Al	3.71615	0.265	23.6445	23.7414	23.7624		
__As	.01046	0.738	1.99363	2.02170	2.01606		
__B	0.96074	0.763	0.95307	0.96144	0.96770		
__Ba	1.83317	0.310	4.82227	4.82694	4.85028		
__Be	.45384	0.340	0.45240	0.45364	0.45547		
__Ca	48.53108	0.522	48.2540	48.5881	48.7510		
__Cd	.78864	0.492	4.76807	4.78350	4.81434		
__Co	.92786	0.492	0.92265	0.92975	0.93118		
__Cr	.02548	0.471	5.00083	5.02752	5.04808		
__Cu	.72260	0.301	4.71698	4.71205	4.73877		
__Fe	4.82045	0.556	4.79268	4.82244	4.84622		
__K	49.09427	0.711	48.7056	49.1946	49.3825		
__Li	1.65116	0.325	24.6421	24.5759	24.7354		
__Mg	23.65675	0.411	23.5769	23.6280	23.7652		
__Mn	.81603	0.426	4.79528	4.81642	4.83639		
__Mo	1.31562	0.371	24.2304	24.3059	24.4104		
__Na	1.81883	0.528	51.7444	51.5897	52.1223		
__Ni	.52001	0.437	9.47287	9.53514	9.55203		
__P	1.69293	0.475	24.6292	24.6212	24.8283		
__Pb	1.60784	0.373	24.5030	24.6451	24.6753		
__Sb	.88735	0.990	4.83903	4.88718	4.93583		
__Se	.94217	1.431	1.91605	1.93909	1.97138		
__Si	.11062	0.848	2.09677	2.10425	2.13084		
__Sn	.84701	0.775	4.80551	4.85670	4.87882		
__Sr	.96972	0.405	0.96638	0.96873	0.97406		
__Ti	.97960	0.350	1.97569	1.97550	1.98760		
__Tl	.99840	0.946	5.04824	4.95416	4.99278		
__V	.94449	0.442	1.93871	1.94038	1.95438		
__Zn	.85744	0.507	4.83171	4.85980	4.88081		

Name: 35601I11

LANCASTER LABORATORIES

PAGE: 37

INSTRUMENT ID: 02360

37 CCB 12/21/92 17:11

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.00254	43.35	0.00318	0.00127	0.00318		
Al	.02741	35.90	0.03486	0.03111	0.01625		
As	.00512	16.14	0.00560	0.00417	0.00560		
B	.00105	207.2	0.00350	0.00035	-0.0006		
Ba	.00155	86.60	0.00233	0.00233	-0.0000		
Be	.00041	176.7	0.00124	-0.0000	-0.0000		
Ca	.01827	41.69	0.02707	0.01387	0.01387		
Cd	.00123	100.0	0.00246	0.00123	-0.0000		
Co	.00160	138.0	-0.0026	-0.0030	0.00094		
Cr	.00240	39.91	0.00343	0.00153	0.00224		
Cu	.00060	153.8	0.00161	-0.0002	0.00040		
Fe	.00000	13461	0.00055	-0.0011	0.00056		
K	.21317	93.38	0.43890	0.13793	0.06269		
Li	.00601	86.60	0.01203	0.00300	0.00300		
Mg	.00108	338.4	0.00321	0.00320	-0.0031		
Mn	.00313	21.66	0.00391	0.00274	0.00273		
Mo	.01010	56.70	0.01662	0.00583	0.00786		
Na	.04008	24.73	0.04580	0.04580	0.02863		
Ni	.00026	1460.	0.00237	-0.0046	0.00147		
P	.02906	167.9	0.03010	0.07734	-0.0202		
Pb	.00471	238.1	0.00035	-0.0036	0.01746		
Sb	.00780	11.74	0.00872	0.00780	0.00689		
Se	.00153	1170.	0.01614	0.00691	-0.0184		
Si	.00222	240.1	0.00666	0.00372	-0.0037		
Sn	.00000	60776	0.00295	0.00297	-0.0059		
Sr	.00020	173.3	0.00061	-0.0000	-0.0000		
Ti	.00026	312.7	0.00052	-0.0006	0.00092		
Tl	.00766	1362.	0.11005	-0.0895	-0.0435		
V	.00185	168.7	0.00006	0.00546	0.00003		
Zn	.00146	69.64	0.00204	0.00207	0.00028		

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Run Name: 935601I11

LANCASTER LABORATORIES

PAGE. 39

INSTRUMENT ID: 02360

38 JBMTF 12/21/92 17:16

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	00066	166.9	0.00129	0.00129	-0.0006		
Al	.01500	57.43	0.00965	0.01041	0.02495		
As	.00682	62.78	0.01149	0.00588	0.00308		
B	.03756	3.259	0.03615	0.03828	0.03825		
Ba	03735	0.000	0.03735	0.03735	0.03735		
Be	.00000	170.9	0.00000	-0.0000	-0.0000		
Ca	.98087	0.524	17.0458	16.8793	17.0174		
X Cd	00095	37.21	0.00116	0.00116	0.00054		
Co	.00174	137.5	-0.0041	-0.0017	0.00064		
Cr	.00265	19.64	-0.0030	-0.0028	-0.0020		
Cu	00283	56.88	0.00405	0.00344	0.00100		
Fe	.16765	0.585	0.16708	0.16708	0.16878		
K	76906	0.376	5.78161	5.74397	5.78160		
Li	00200	173.2	-0.0000	0.00000	0.00601		
Mg	51567	0.720	6.55176	6.46259	6.53266		
Mn	07061	0.002	0.07061	0.07061	0.07061		
Mo	00178	151.0	-0.0009	0.00178	0.00448		
Na	66676	0.640	9.71830	9.59804	9.68394		
Ni	00292	59.92	-0.0029	-0.0046	-0.0011		
P	05989	72.06	0.01032	0.08915	0.08021		
Pb	.00536	39.10	-0.0036	-0.0077	-0.0046		
Sb	00569	50.51	-0.0082	-0.0025	-0.0063		
Se	01664	63.48	0.00511	0.02587	0.01895		
Si	72083	0.815	4.75589	4.67959	4.72700		
Sn	00281	158.4	0.00281	0.00726	-0.0016		
Sr	08040	0.778	0.08095	0.07972	0.08054		
Ti	00125	102.6	-0.0000	0.00251	0.00132		
Tl	.08461	27.72	-0.1051	-0.0896	-0.0590		
V	.00047	166.6	0.00000	0.00137	0.00002		
Zn	03357	0.015	0.03356	0.03357	0.03357		

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Name: 35601I11

LANCASTER LABORATORIES

PAGE: 39

INSTRUMENT ID: 02360

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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.00047	647.7	0.00079	0.00175	-0.0039		
Al	37691	0.312	1.37196	1.37977	1.37901		
As	00147	953.5	0.01548	-0.0125	0.00150		
B	02394	5.062	0.02534	0.02328	0.02321		
Ba	04637	1.453	0.04676	0.04559	0.04676		
Be	.00001	1.001	-0.0000	-0.0000	-0.0000		
Ca	10.40070	0.593	10.4001	10.3392	10.4626		
Cd	.00038	91.53	-0.0008	-0.0001	-0.0001		
Co	.00130	184.1	0.00146	-0.0027	-0.0025		
Cr	00007	1776.	0.00111	0.00043	-0.0013		
Cu	00339	71.65	0.00583	0.00339	0.00096		
Fe	56273	0.674	2.55932	2.54740	2.58147		
K	78421	2.064	2.84692	2.73404	2.77167		
Li	.00000	116.0	0.00000	0.00000	0.00000		
Mg	34347	0.593	4.34772	4.31585	4.36683		
Mn	11516	2.036	0.11516	0.11281	0.11751		
Mo	00241	116.0	0.00017	0.00556	0.00151		
Na	.47448	0.751	10.4057	10.4572	10.5603		
Ni	00095	104.7	0.00211	0.00038	0.00037		
P	11569	26.65	0.12680	0.08083	0.13944		
Pb	.00334	304.9	0.00202	0.00304	-0.0150		
Sb	00316	168.9	-0.0075	0.00279	-0.0047		
Se	00531	275.5	0.01377	0.01377	-0.0116		
Si	74353	0.648	5.73856	5.70898	5.78303		
Sn	00263	447.9	0.01600	-0.0062	-0.0018		
Sr	05991	0.197	0.05984	0.05984	0.06004		
Ti	05275	4.450	0.05489	0.05053	0.05132		
Tl	12126	91.08	0.00433	-0.2032	-0.1648		
V	00553	0.172	0.00552	0.00554	0.00552		
Zn	02830	4.722	0.02976	0.02801	0.02714		

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Run Num 25631111

LANCASTER LABORATORIES

DATE: 4/

INSTRUMENT ID: 02360

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ITEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	00034	724.0	0.00021	-0.0015	-0.0006		
Al	02708	47.07	0.04174	0.01863	0.02086		
As	00619	113.7	0.00033	-0.0052	-0.0136		
B	02652	0.056	0.02653	0.02650	0.02652		
Ba	03035	0.000	0.03035	0.03035	0.03035		
Be	00000	85.99	-0.0000	-0.0000	-0.0000		
Ca	68568	0.638	10.6073	10.7317	10.7180		
Cd	00067	159.2	0.00056	-0.0012	-0.0012		
Co	00075	104.9	-0.0006	-0.0000	-0.0016		
Cr	00028	589.9	0.00061	0.00074	-0.0022		
Cu	00384	18.27	0.00314	0.00466	0.00344		
Fe	21717	0.453	0.21831	0.21660	0.21661		
K	03472	1.958	2.99742	2.88456	2.92218		
Li	00000	19.34	0.00000	0.00000	0.00000		
Mg	25116	0.887	4.30976	4.38621	4.35750		
Mn	08710	2.694	0.08710	0.08945	0.08475		
Mo	00200	19.34	0.00245	0.00178	0.00178		
Na	96101	0.478	10.9937	11.0068	10.9724		
Ni	00144	159.8	0.00219	0.00231	-0.0011		
P	03263	232.6	-0.0751	-0.0777	0.05500		
Pb	00569	127.6	0.00237	-0.0077	-0.0117		
Sb	00378	208.0	0.00501	-0.0063	-0.0101		
Se	00108	537.5	-0.0043	0.00723	0.00031		
Si	27576	0.648	4.59718	4.65738	4.62552		
Sn	00285	270.9	0.00711	0.00730	-0.0060		
Sr	06045	0.897	0.06031	0.06086	0.06066		
Ti	00158	62.85	0.00251	0.00053	0.00172		
Tl	07952	117.9	0.00243	-0.0591	-0.1818		
V	00274	85.83	0.00547	0.00139	0.00138		
Zn	01715	5.895	0.01773	0.01773	0.01598		

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Name: 935601I11

LANCASTER LABORATORIES

PAGE: 41 INSTRUMENT ID: 02360

41 JSBLT 12/21/92 17:29
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ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.00141	78.27	-0.0020	-0.0001	-0.0020		
Al	53250	0.383	1.53745	1.52602	1.53404		
As	.00204	439.6	0.00450	-0.0123	0.00165		
B	.03750	3.146	0.03687	0.03897	0.03697		
Ba	.05844	0.001	0.05844	0.05844	0.05843		
Be	00081	88.06	0.00123	0.00122	-0.0000		
Ca	.87055	0.607	15.9294	15.9228	15.7593		
X Cd	.00143	24.98	0.00163	0.00101	0.00164		
Co	.00126	178.5	-0.0021	-0.0029	0.00130		
Cr	.00425	62.76	0.00732	0.00282	0.00259		
Cu	.01597	8.806	0.01678	0.01435	0.01678		
Fe	69804	0.746	2.70656	2.71252	2.67504		
K	43043	0.799	5.40535	5.40534	5.48059		
Li	00401	43.30	0.00300	0.00300	0.00601		
Mg	31500	0.949	6.35428	6.34472	6.24600		
Mn	13153	0.002	0.13153	0.13153	0.13153		
Mo	00104	197.8	-0.0012	0.00149	0.00284		
Na	.01076	0.748	10.0451	10.0623	9.92485		
Ni	.00152	66.09	0.00210	0.00036	0.00210		
P	11694	35.47	0.11433	0.15967	0.07683		
Pb	.00309	73.39	-0.0009	-0.0029	-0.0054		
Sb	.00165	173.7	-0.0010	0.00085	-0.0047		
Se	.01236	149.7	0.03005	0.01390	-0.0068		
Si	38651	0.333	6.39367	6.40330	6.36257		
Sn	.00485	140.2	-0.0107	0.00257	-0.0063		
Sr	07746	0.825	0.07767	0.07797	0.07674		
Ti	08154	0.921	0.08187	0.08207	0.08068		
Tl	.00971	0.720	-0.0097	-0.0096	-0.0097		
V	00689	19.70	0.00689	0.00825	0.00553		
Zn	07965	1.280	0.07906	0.07907	0.08083		

128

AR318997

Cd

Run Name: [illegible]5601I11

LANCASTER LABORATORIES

RACE: 42 INSTRUMENT ID: 02360

12 [illegible] JSBLF 12/21/92 17:34
[illegible].100 1 .DF1,923541846001,1,,

ELEM	RAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	00002	4624.	0.00129	-0.0006	-0.0006		
Al	01780	48.71	0.01292	0.01266	0.02781		
As	.00670	55.21	-0.0081	-0.0095	-0.0025		
B	03980	3.049	0.03840	0.04051	0.04050		
Ba	03813	3.534	0.03735	0.03735	0.03968		
Be	00000	88.67	-0.0000	0.00000	-0.0000		
Ca	68473	0.122	15.6674	15.7055	15.6811		
X Cd	00006	0.412	-0.0000	-0.0000	-0.0000		
Co	00197	60.22	-0.0006	-0.0023	-0.0029		
Cr	.00106	217.9	0.00093	-0.0036	-0.0005		
Cu	0344	35.37	0.00466	0.00222	0.00344		
Fe	18526	0.530	0.18413	0.18583	0.18583		
K	5487	1.431	5.25487	5.17962	5.33011		
Li	00000	295.1	-0.0000	-0.0000	0.00000		
Mg	0604	0.270	6.11558	6.08694	6.11560		
Mn	9057	2.244	0.08823	0.09175	0.09175		
Mo	00091	295.1	-0.0035	-0.0009	0.00178		
Na	36543	0.834	10.4570	10.3539	10.2852		
Ni	00321	15.91	-0.0038	-0.0029	-0.0029		
P	00433	993.5	0.04321	-0.0420	0.01181		
Pb	0068	519.8	-0.0026	0.00437	0.00034		
Sb	0754	126.0	0.00880	-0.0025	0.01634		
Se	2775	120.1	0.02172	-0.0018	0.06390		
Si	1808	0.296	4.80548	4.83363	4.81513		
Sr	0579	88.84	0.01174	0.00282	0.00282		
Sr	7661	0.278	0.07654	0.07644	0.07685		
Ti	0211	56.22	0.00330	0.00211	0.00092		
Tl	07685	56.79	-0.1126	-0.0282	-0.0897		
V	0182	86.12	0.00273	0.00001	0.00274		
Zn	2740	3.207	0.02740	0.02652	0.02828		

AR318998

cd

Run Nam 9235601I11

LANCASTER LABORATORIES

PAGE: 1 INSTRUMENT ID: 02360

43 [illegible]02 CRHLT 12/21/92 17:39
****,10[illegible]0.DF1.923541848001,1,,

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible].00057	163.2	0.00131	0.00131	-0.0005		
Al	[illegible].06889	13.54	0.07047	0.07734	0.05887		
As	0.00240	364.2	0.00037	-0.0121	0.00461		
B	[illegible].02901	7.162	0.02694	0.03109	0.02901		
Ba	[illegible].02529	2.664	0.02568	0.02568	0.02451		
Be	0.00000	173.4	-0.0000	-0.0000	0.00000		
Ca	[illegible].30394	0.997	8.22271	8.38822	8.30089		
X Cd	-0.00004	2.270	-0.0000	-0.0000	-0.0000		
Co	-0.00238	63.29	-0.0006	-0.0033	-0.0031		
Cr	-0.00191	116.6	-0.0027	0.00061	-0.0036		
Cu	[illegible].00344	35.33	0.00344	0.00466	0.00222		
Fe	[illegible].24565	1.200	0.24395	0.24905	0.24394		
K	[illegible].69641	2.790	2.69641	2.62116	2.77166		
Li	[illegible].00000	315.0	-0.0000	0.00000	0.00000		
Mg	[illegible].31171	1.236	3.26500	3.34143	3.32869		
Mn	[illegible].07305	0.002	0.07305	0.07304	0.07305		
Mo	[illegible].00042	315.0	-0.0009	0.00177	0.00042		
Na	[illegible].99809	0.567	6.95227	7.02101	7.02100		
Ni	-0.00001	6777.	0.00056	-0.0011	0.00056		
P	[illegible].00708	225.5	0.02019	0.01175	-0.0107		
Pb	-0.01003	110.0	-0.0227	-0.0036	-0.0036		
Sb	[illegible].00377	252.5	0.01262	0.00498	-0.0063		
Se	-0.01054	63.21	-0.0182	-0.0066	-0.0066		
Si	[illegible].36409	1.026	3.32807	3.39690	3.36731		
Sr	[illegible].00510	226.8	0.0015	-0.0015	0.01847		
Sr	[illegible].04566	1.151	0.04508	0.04611	0.04580		
Ti	[illegible].00132	130.6	0.00053	0.00330	0.00013		
Tl	0.07181	72.80	-0.0129	-0.0896	-0.1128		
V	[illegible].00183	171.2	0.00545	0.00002	0.00001		
Zn	[illegible].01248	0.032	0.01248	0.01247	0.01248		

180

AR318999

RCI

Run Nam /35601I11

LANCASTER LABORATORIES

PAGE: A

INSTRUMENT ID: 02360

44 1 3 CRHLF 12/21/92 17:43
****.100 0.DF1,923541848001,1..

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.00070	156.2	-0.0000	-0.0019	-0.0000		
Al	1.85279	0.940	1.85778	1.86718	1.83340		
As	0.00526	201.8	-0.0174	0.00223	-0.0005		
B	.02924	7.233	0.03135	0.02712	0.02926		
Ba	.04211	0.000	0.04211	0.04211	0.04211		
Be	0.00001	27.19	-0.0000	-0.0000	-0.0000		
Ca	.60263	0.399	8.62734	8.61718	8.56337		
X Cd	.00021	865.5	0.00227	-0.0014	-0.0001		
Co	0.00057	303.3	0.00144	-0.0015	-0.0015		
Cr	.00046	191.5	0.00055	-0.0004	0.00129		
Cu	.00481	36.49	0.00582	0.00278	0.00582		
Fe	.00805	0.256	3.00919	3.01515	2.99982		
K	.79673	1.553	2.77165	2.84690	2.77165		
Li	.00000	139.0	0.00000	0.00000	0.00000		
Mg	.55343	1.050	3.52582	3.59590	3.53856		
Mn	.10186	1.328	0.10108	0.10342	0.10108		
Mo	.00055	139.0	0.00145	0.00010	0.00011		
Na	.34487	0.472	7.36493	7.36490	7.30479		
Ni	.00032	1417.	0.00557	-0.0031	-0.0014		
P	.06471	92.04	0.07544	0.11818	0.00051		
Pb	.00008	9979.	0.00630	0.00378	-0.0098		
Sb	0.00165	728.5	0.01225	-0.0086	-0.0086		
Se	.00206	193.6	-0.0002	0.00667	-0.0002		
Si	.32440	0.074	5.32141	5.32888	5.32291		
Sn	.00711	62.56	0.01157	0.00266	0.00711		
Sr	.04822	0.648	0.04857	0.04816	0.04795		
Ti	.06481	5.231	0.06799	0.06124	0.06521		
Tl	0.04659	75.99	-0.0760	-0.0262	-0.0874		
V	.00644	24.42	0.00825	0.00553	0.00553		
Zn	.03091	1.693	0.03060	0.03151	0.03062		

[illegible] 235601I11

LANCASTER LABORATORIES

PAGE [illegible]

INSTRUMENT ID: 02360

[illegible] 12/21/92 17:48

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	0.03423	12.21	-0.0364	-0.0294	-0.0368		
Al	462.89083	1.070	468.123	458.270	462.278		
As	[illegible].21061	48.25	0.32631	0.13583	0.16968		
B	[illegible].26814	1.561	0.26332	0.27087	0.27023		
Ba	[illegible].00554	0.960	0.00560	0.00549	0.00553		
Be	[illegible].00163	44.09	0.00122	0.00122	0.00247		
Ca	455.22958	0.858	459.583	452.031	454.073		
Cd	[illegible].02894	6.617	0.03102	0.02725	0.02855		
Co	0.00282	68.91	-0.0029	-0.0008	-0.0046		
Cr	0.01335	27.34	-0.0165	-0.0093	-0.0141		
Cu	[illegible].00216	101.5	0.00153	0.00461	0.00034		
Fe	[illegible]73.24012	0.959	175.079	171.841	172.799		
K	[illegible].22514	33.24	0.22629	0.30127	0.15088		
Li	0.00100	173.1	-0.0000	-0.0000	-0.0030		
Mg	477.04806	1.036	482.420	472.694	476.028		
Mn	0.01167	1.854	-0.0119	-0.0114	-0.0116		
Mo	0.04852	10.03	-0.0485	-0.0436	-0.0534		
Na	[illegible].13466	20.50	0.12320	0.16616	0.11461		
Ni	0.02702	11.76	-0.0289	-0.0233	-0.0287		
P	1.70644	10.23	-1.5251	-1.8735	-1.7207		
Pb	[illegible].02562	49.55	0.03998	0.02096	0.01591		
Sb	0.01160	317.7	0.02353	-0.0083	-0.0499		
Se	0.06802	26.00	-0.0541	-0.0619	-0.0879		
Si	[illegible].00296	257.2	0.00531	0.00914	-0.0055		
Sn	[illegible].09318	25.67	0.10068	0.11245	0.06640		
Sr	[illegible].01683	1.531	0.01711	0.01680	0.01659		
Ti	[illegible].01322	7.959	0.01282	0.01441	0.01242		
Tl	0.39504	5.305	-0.4191	-0.3815	-0.3843		
V	[illegible].01044	0.405	0.01049	0.01042	0.01041		
Zn	0.00984	5.486	-0.0104	-0.0093	-0.0097		

132

AR319001

Run Name: [illegible]3601I11

LANCASTER LABORATORIES

PAGE: 4[illegible]

INSTRUMENT ID: 02360

16 ICS [illegible] 12/21/92 17:53

ITEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	[illegible]7375	0.736	0.88118	0.86997	0.87009		
Al	[illegible].94210	0.595	460.295	462.731	465.799		
As	[illegible]23303	21.49	0.22232	0.28762	0.18916		
B	[illegible]27405	1.347	0.27512	0.27708	0.26994		
Ba	[illegible]46265	0.641	0.45950	0.46305	0.46540		
Be	[illegible]42273	0.612	0.41983	0.42354	0.42481		
Ca	[illegible]4.96691	0.613	451.773	456.204	456.923		
Cd	[illegible]86026	0.831	0.85204	0.86497	0.86379		
Co	[illegible]40978	1.789	0.40178	0.41137	0.41618		
Cr	[illegible]42635	1.569	0.41932	0.43264	0.42708		
Cu	[illegible]44924	0.154	0.44844	0.44965	0.44964		
Fe	[illegible].28697	0.582	172.178	173.531	174.150		
K	[illegible]27659	15.73	0.22634	0.30169	0.30175		
Li	[illegible].00000	1355	-0.0000	-0.0030	0.00300		
Mg	[illegible]6.77355	0.610	473.734	477.052	479.533		
Mn	[illegible]42110	0.615	0.41811	0.42265	0.42255		
Mo	[illegible].04637	4.787	-0.0438	-0.0481	-0.0470		
Na	[illegible]16297	10.96	0.14865	0.15725	0.18300		
Ni	[illegible]80878	0.532	0.80688	0.81372	0.80576		
P	[illegible].73846	5.853	-1.6383	-1.7351	-1.8418		
Pb	[illegible]92415	1.602	0.91040	0.93983	0.92222		
Sb	[illegible].01779	7.711	-0.0189	-0.0181	-0.0162		
Se	[illegible].06589	53.83	-0.0259	-0.0781	-0.0936		
Si	[illegible]00183	601.4	-0.0100	0.01173	0.00385		
Sn	[illegible]08592	7.884	0.08852	0.09101	0.07823		
Sr	[illegible]01695	1.256	0.01677	0.01718	0.01688		
Ti	[illegible]01230	16.56	0.01177	0.01455	0.01058		
Tl	[illegible]41233	17.74	-0.4360	-0.3302	-0.4706		
V	[illegible]44960	1.216	0.44411	0.45505	0.44962		
Zn	[illegible]87336	0.804	0.85547	0.87567	0.87895		

[redacted] 35601111

LANCASTER LABORATORIES

[redacted] INSTRUMENT ID: 02360

[redacted] 12/21/93 17:57

[redacted]	PPAGE (ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
[redacted]	02165	5.103	0.02229	0.02229	0.02037		
[redacted]	04909	49.32	0.06668	0.05912	0.02147		
[redacted]	01350	41.58	0.01912	0.00789	0.01348		
[redacted]	.00354	33.89	-0.0028	-0.0028	-0.0049		
[redacted]	00000	26.58	0.00000	0.00000	0.00000		
[redacted]	00866	4.192	0.00908	0.00845	0.00844		
[redacted]	04332	30.73	0.05753	0.04129	0.03113		
[redacted]	00905	15.74	0.00740	0.00987	0.00987		
[redacted]	09557	1.084	0.09506	0.09676	0.09489		
[redacted]	02155	6.690	0.02213	0.02262	0.01991		
[redacted]	04829	3.174	0.04849	0.04971	0.04667		
[redacted]	01611	26.06	0.02004	0.01327	0.01413		
[redacted]	13796	0.002	0.13796	0.13796	0.13795		
[redacted]	.00000	167.8	0.0000	-0.0000	0.00000		
[redacted]	04185	45.65	0.07696	0.04186	0.02274		
[redacted]	02857	0.003	0.02857	0.02857	0.02857		
[redacted]	.00046	167.8	0.0009	-0.0009	0.00043		
[redacted]	04576	0.016	0.04576	0.04576	0.04575		
[redacted]	07247	4.238	0.07567	0.06955	0.07218		
[redacted]	.01714	180.0	0.0251	-0.0432	0.01692		
[redacted]	00786	46.16	0.0102	-0.0097	-0.0036		
[redacted]	10555	5.742	0.09863	0.10994	0.10808		
[redacted]	02365	40.28	0.01373	0.01834	0.02989		
[redacted]	.01408	55.55	0.0065	-0.0021	-0.0035		
[redacted]	00301	21792	0.00298	-0.0014	-0.0014		
[redacted]	00011	83.13	0.0002	-0.0002	-0.0000		
[redacted]	00384	31.53	0.00490	0.00410	0.00252		
[redacted]	[redacted]44	87.81	0.00171	0.00112	0.04709		
[redacted]	09391	1.587	0.09801	0.09801	0.10073		
[redacted]	03870	0.027	0.03863	0.03870	0.03871		

File Name: 0235601I11

LANCASTER LABORATORIES

PAGE: 4

INSTRUMENT ID: 02360

AS CO 12/21/92 18:02

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.46482	0.627	0.46546	0.46737	0.46164		
Al	23.34699	0.894	23.4892	23.4443	23.1072		
As	1.95629	1.576	1.98803	1.95443	1.92642		
B	.94118	0.645	0.94465	0.94472	0.93416		
Ba	.74019	1.087	4.77559	4.76392	4.68106		
Be	.44501	0.959	0.44745	0.44749	0.44008		
Ca	47.39250	0.720	47.5590	47.6189	46.9995		
Cd	.66237	0.775	4.67676	4.68910	4.62126		
Co	.90927	1.388	0.91938	0.91331	0.89512		
Cr	.91870	1.015	4.94867	4.94641	4.86104		
Cu	.64100	0.982	4.66971	4.66484	4.58844		
Fe	.71714	0.906	4.74515	4.73835	4.66791		
K	48.31766	0.468	48.5556	48.2923	48.1050		
Li	24.46256	1.305	24.6902	24.5999	24.0974		
Mg	23.20896	0.799	23.3695	23.2515	23.0057		
Mn	.72439	0.861	4.75297	4.74241	4.67778		
Mo	23.76880	1.105	23.9197	23.9211	23.4655		
Na	53.43265	1.422	54.0896	53.7288	52.6294		
Ni	.32358	0.808	9.38110	9.35138	9.23826		
P	24.00597	1.636	24.4221	23.9542	23.6415		
Pb	24.06859	0.482	24.1640	24.1025	23.9392		
Sb	.76348	1.052	4.75513	4.81728	4.71802		
Se	.90617	2.352	1.93919	1.92420	1.85512		
Si	.15961	0.341	2.16057	2.16646	2.15181		
Sn	.72592	1.055	4.72795	4.77474	4.67507		
Sr	.95135	1.077	0.95869	0.95572	0.93964		
Ti	93982	1.024	1.94988	1.95265	1.91693		
Tl	95624	2.077	5.04957	4.85084	4.99830		
V	90659	1.000	1.92259	1.91170	1.88549		
Zn	76425	0.643	4.78101	4.78289	4.72886		

Run Name: [illegible]5601T11

LANCASTER LABORATORIES

CASE: 47 INSTRUMENT ID: 02360

10 CI[illegible] 12/21/92 18:06

ELEM	AVERAGE(ppm)	%RSD	BURN#1	BURN#2	BURN#3	REREAD	REDIG
Ag	.00063	173.0	0.00127	0.00127	-0.0006		
Al	.02121	40.41	0.03111	0.01611	0.01641		
As	.00653	24.79	-0.0055	-0.0084	-0.0055		
B	.00174	103.2	-0.0007	-0.0007	-0.0038		
Ba	.00311	86.60	0.00466	0.00466	-0.0000		
Be	00083	85.77	0.00124	0.00124	0.00000		
Ca	.03384	61.79	0.05347	0.03621	0.01184		
Cd	00370	33.30	0.00493	0.00370	0.00246		
Co	.00292	90.61	0.00011	-0.0047	-0.0041		
Cr	.00014	2493.	0.00364	-0.0002	-0.0037		
Cu	00242	210.4	0.00830	-0.0002	-0.0008		
Fe	00339	115.5	0.00566	0.00567	-0.0011		
K	.12538	45.82	0.17553	0.13792	0.06269		
Li	.01504	52.91	0.02106	0.01805	0.00601		
Mg	.01805	146.9	0.04779	0.00956	-0.0031		
Mn	00195	69.24	0.00274	0.00274	0.00039		
Mo	.01662	48.64	0.02471	0.01662	0.00853		
Na	.07444	48.04	0.11453	0.06298	0.04581		
Ni	.00471	93.79	0.00939	0.00413	0.00061		
P	.02793	45.51	-0.0419	-0.0245	-0.0172		
Pb	.00269	714.9	0.00840	0.01847	-0.0187		
Sb	.01109	112.5	-0.0016	-0.0063	-0.0252		
Se	.00845	87.68	0.01383	-0.0000	0.01153		
Si	.00421	180.4	-0.0022	0.00222	-0.0125		
Sn	00295	399.4	0.00739	0.01186	-0.0104		
Sr	.00020	199.9	0.00061	0.00020	-0.0002		
Ti	.00145	113.5	0.00330	0.00092	0.00013		
Tl	.08458	37.84	-0.1204	-0.0589	-0.0742		
V	.00085	188.0	0.00009	0.00005	-0.0027		
Zn	00291	89.61	0.00552	0.00293	0.00029		

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196