

**Appendix A.  
Abbreviations**

ACFB	atmospheric circulating fluidized bed
AF	as fed
B&W	Babcock & Wilcox
CFD Model	computer fluid dynamic model
COE	cost of energy
DB	dry basis
DOE	Department of Energy
EPC	engineer, procure, and construct
FIR	Forced Internal Recirculation
GTI	Gas Technology Institute
HHV	high heating value
LCV	low caloric value gas
LHV	low heating value
MF	moisture free
NEPA	National Environmental Protection Act
NOx	nitrogen oxides
O&M	operating and maintenance
psia	pounds per square inch - absolute
VOC	volatile organic carbon

**Appendix B**  
**List of Specifications Developed**

Gasifier (including cyclone and startup heater)  
Fuel Feed  
Limestone Feed  
Gasifier Ash Removal  
Process Air  
Product Gas Ducting  
Product Gas Cooling  
Product Gas Filtering  
Filter Ash Removal  
Flare  
High Pressure Cooling Water  
Nitrogen Distribution  
Biomass Dryer Specification  
Biomass Feed Prep  
Inert Gas Generation System  
Waste Heat Boiler  
Startup Gas Supply

**Appendix C**  
**Analytical Data**



## **Analytical Report**

Batch # : 011049

March 16, 2001

### **Prepared for:**

**Dave Stopek, Process Engineering**

**Phone: (847) 768-0853**

**Kentucky Co-gasification Project**

**Project #: 61129-01**

**Received Date: 2/5/2001**

### Disclaimer:

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In April 2000, the Institute of Gas Technology (IGT) and the Gas Research Institute (GRI) combined to form the Gas Technology Institute (GTI).

Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report: \_\_\_\_\_

Alan G. Janos

(847) 768-0603



Sample Login No: 011049-002  
 Sample Description: Estill County Saw Mill Dust

Date: February 19, 2001

Air-Dry Moisture, %	48.93		
<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u> w/SO <sub>3</sub> correction	<u>(Dry basis)</u>
Moisture, %	50.52	50.52	
Ash, % (550° C)	0.46	0.46	

**Ultimate Analysis** (Dry basis)

Ash, % (550° C)	0.92
Carbon, %	49.07
Hydrogen, %	6.00
Nitrogen, %	0.27
Sulfur, %	0.13
Oxygen, % (by difference)	43.61

*Fusion Temperature of Ash (ASTM D1857), °F*  
*Reducing    Oxidizing*

*Initial Deformation (IT)*  
*Softening (ST)*  
*Hemispherical (HT)*  
*Fluid (FT)*

**Heating Value** (Dry basis)

BTU/lb.	8,990
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Analyst: NJP

**Bulk Density:** 20.8 lbs/ft<sup>3</sup>

Analyst: AGJ

(Per submitter's request, this analysis was performed on the "as-received" sample, which was NOT air-dried, chopped, or ground into any smaller pieces than were originally received. The graduated cylinder used for the volume measurement was approximately 2¼" in diameter. No significant void spaces were noted in the packed sample at the end of the analysis.)



Login No.: 01-1049-002

Sample ID: Estill County Saw Mill Dust

**MAJOR / MINOR OXIDES IN ASH (ASTM D-2795)**

% ash at 750 degrees C = 0.74

<u>% Element</u>		<u>% Oxide</u>		<u>% Oxide, Normalized</u>
Na	0.17	Na <sub>2</sub> O	0.23	0.24
Mg	2.35	MgO	3.90	4.12
Al	1.86	Al <sub>2</sub> O <sub>3</sub>	3.52	3.72
Si	10.3	SiO <sub>2</sub>	22.0	23.3
P	1.06	P <sub>2</sub> O <sub>5</sub>	2.43	2.57
S	0.41	SO <sub>3</sub>	1.02	1.08
K	12.6	K <sub>2</sub> O	15.2	16.1
Ca	31.5	CaO	44.1	46.6
Ti	0.15	TiO <sub>2</sub>	0.25	0.26
Fe	1.31	Fe <sub>2</sub> O <sub>3</sub>	1.87	1.98
		TOTAL	94.53	100.0

Values reported on an ash basis.

Ash concentration calculated on a dry basis with SO<sub>3</sub> correction and no VM cycle.

Chloride and/or carbonate may be present, resulting in a total % less than 100.

Analyst: KC



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FAX: (708) 333-3060

March 1, 2001

INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mt. Prospect Road  
Des Plaines, IL 60018  
Attn: Norman J. Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample  
reported to us Wood Ash

Sample ID: 011049-002  
Estill County Saw Dust

Sample taken at G.T.I.

Sample taken by G.T.I.

Date sampled February 16, 2001

Date received February 27, 2001

P.O No. PF8847

Analysis Report No. 71-144427

Page 1 of 1

FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2540	2565
Softening (ST)	2555	2580
Hemispherical (HT)	2570	2600
Fluid (FT)	2585	2620

METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory



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Sample Login No: 011049-003

Date: February 19, 2001

Sample Description: Estill County Saw Mill Slabs/Bark

Air-Dry Moisture, %	38.33		
<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u> <u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	40.55	40.55	
Ash, % (550° C)	6.49	6.46	

**Ultimate Analysis** (Dry basis)

Ash, % (550° C)	10.87
Carbon, %	46.34
Hydrogen, %	5.38
Nitrogen, %	0.40
Sulfur, %	0.03
Oxygen, % (by difference)	36.97

*Fusion Temperature of Ash (ASTM D1857), °F*  
*Reducing    Oxidizing*

*Initial Deformation (IT)*  
*Softening (ST)*  
*Hemispherical (HT)*  
*Fluid (FT)*

**Heating Value** (Dry basis)

BTU/lb.	7,630
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Analyst: NJP

**Bulk Density:** 16.5 lbs/ft<sup>3</sup>

Analyst: AGJ

(Per submitter's request, this analysis was performed on the "as-received" sample, which was NOT air-dried but was chopped into pieces of < 1½" in length. The graduated cylinder used for the volume measurement was approximately 3" in diameter. Void spaces were noted which were not removed by additional tapping. Total sample volume measured includes these voids.)





Login No.: 01-1049-003

Sample ID: Estill County Saw Mill Slabs/Bark

**MAJOR / MINOR OXIDES IN ASH (ASTM D-2795)**

% ash at 750 degrees C = 7.93

<u>% Element</u>		<u>% Oxide</u>		<u>% Oxide, Normalized</u>
Na	0.09	Na <sub>2</sub> O	0.12	0.12
Mg	0.68	MgO	1.13	1.15
Al	2.43	Al <sub>2</sub> O <sub>3</sub>	4.59	4.68
Si	15.6	SiO <sub>2</sub>	33.4	34.0
P	0.30	P <sub>2</sub> O <sub>5</sub>	0.69	0.70
S	0.16	SO <sub>3</sub>	0.41	0.42
K	2.66	K <sub>2</sub> O	3.21	3.3
Ca	37.4	CaO	52.3	53.3
Ti	0.22	TiO <sub>2</sub>	0.37	0.37
Fe	1.32	Fe <sub>2</sub> O <sub>3</sub>	1.89	1.92
		TOTAL	98.09	100.0

Values reported on an ash basis.

Ash concentration calculated on a dry basis with SO<sub>3</sub> correction and no VM cycle.

Chloride and/or carbonate may be present, resulting in a total % less than 100.

Analyst: KC



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March 1, 2001

INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mt. Prospect Road  
Des Plaines, IL 60018  
Attn: Norman J. Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Wood Ash

Sample taken at G.T.I.

Sample ID: 011049-003  
Estill County Bark

Sample taken by G.T.I.

Date sampled February 16, 2001

Date received February 27, 2001

P.O No. PF8847

Analysis Report No. 71-144428

Page 1 of 1

FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2550	2385
Softening (ST)	2560	2410
Hemispherical (HT)	2575	2425
Fluid (FT)	2600	2440

METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
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Sample Login No: 011049-001

Date: February 19, 2001

Sample Description: Estill County Prep. Plant coal sample

Air-Dry Moisture, % 6.63

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u>	
		<u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	7.62	7.62	
Volatile Matter, %	32.94	32.94	35.66
Ash, % (750° C)	5.56	5.40	
Fixed Carbon, % (by difference)	53.87	54.04	

**Ultimate Analysis** (Dry basis)

Ash, % (750° C)	5.85
Carbon, %	78.83
Hydrogen, %	5.30
Nitrogen, %	1.83
Sulfur, %	0.96
Oxygen, % (by difference)	7.24

*Fusion Temperature of Ash (ASTM D1857), °F*  
*Reducing    Oxidizing*

*Initial Deformation (IT)*  
*Softening (ST)*  
*Hemispherical (HT)*  
*Fluid (FT)*

**Heating Value** (Dry basis)

BTU/lb. 14,140

Analyst: NJP

**Bulk Density:** 45.5 lbs/ft<sup>3</sup>

Analyst: AGJ

(Per submitter's request, this analysis was performed on the "as-received" sample, which was NOT air-dried, crushed, or ground into any smaller pieces than were originally received. The graduated cylinder used for the volume measurement was approximately 1 1/8" in diameter. No visible void spaces were noted in the packed sample at the end of the analysis.)



Login No.: 01-1049-001

Sample ID: Estill County Prep. Plant coal sample

**MAJOR / MINOR OXIDES IN ASH (ASTM D-2795)**

% ash at 750 degrees C = 6.29

<u>% Element</u>		<u>% Oxide</u>		<u>% Oxide, Normalized</u>
Na	0.41	Na <sub>2</sub> O	0.55	0.55
Mg	0.85	MgO	1.41	1.40
Al	14.4	Al <sub>2</sub> O <sub>3</sub>	27.2	27.0
Si	24.6	SiO <sub>2</sub>	52.6	52.2
P	0.21	P <sub>2</sub> O <sub>5</sub>	0.48	0.48
S	1.31	SO <sub>3</sub>	3.26	3.24
K	2.13	K <sub>2</sub> O	2.57	2.55
Ca	2.59	CaO	3.62	3.60
Ti	1.01	TiO <sub>2</sub>	1.68	1.67
Fe	5.14	Fe <sub>2</sub> O <sub>3</sub>	7.35	7.29
		TOTAL	100.8	100.0

Values reported on an ash basis.

Ash concentration calculated on a dry basis with SO<sub>3</sub> correction and no VM cycle.

Analyst: KC



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March 1, 2001

INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mt. Prospect Road  
Des Plaines, IL 60018  
Attn: Norman J. Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Coal Ash

Sample taken at G.T.I.

Sample ID: 011049-001  
Estill County Coal

Sample taken by G.T.I.

Date sampled February 16, 2001

Date received February 27, 2001

P.O No. PF8847

Analysis Report No. 71-144426

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2485	2640
Softening (ST)	2515	2675
Hemispherical (HT)	2570	2700+
Fluid (FT)	2610	2700+

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
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## Analytical Report

Batch # : 021478

December 18, 2002

**Prepared for:**

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 12/4/2002**

**Disclaimer:**

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Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report:

A handwritten signature in black ink that reads "Alan G. Janos". The signature is written in a cursive style and is positioned above a horizontal line.

Alan G. Janos

(847) 768-0603



Sample Login No: 021478-001

Date: December 18, 2002

Sample Description: Activated coconut charcoal, -4+8 mesh, Springfield Scientific #TC4716

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u>	
		<u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	3.28	3.28	
Volatile Matter, %	4.98	4.98	5.15
Ash (750°F), %	2.06	2.02	
Fixed Carbon, % (by difference)	89.68	89.72	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>	<b>Screen Analysis</b>	
		<u>Retained on</u>	<u>Wt. %</u>
Ash (750°F), %	2.09		
Carbon, %	94.62		
Hydrogen, %	0.23	3	0.0%
Nitrogen, %	0.61	4	5.8%
Sulfur, %	0.04	6	47.6%
Oxygen, %	2.40	8	34.1%
(by difference)		PAN	<u>12.5%</u>
		Total	100.0%

<b>Heating Value</b>	<u>(Dry basis)</u>
BTU/lb.	14,050

Analyst: NJP



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INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mount Prospect Road  
Des Plaines, IL 60018  
Attn: Norman Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Activated Coconut Charcoal

Sample ID: 021478-001  
Activated Coconut Charcoal

Sample taken at -----

Sample taken by -----

Date sampled -----

Date received February 3, 2003

P.O No. PF00020103

Analysis Report No. 71-198925

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2400	2445
Softening (ST)	2415	2460
Hemispherical (HT)	2430	2485
Fluid (FT)	2450	2505

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory







Sample Login No: 021478-002

Date: December 18, 2002

Sample Description: Calla Energy, Mill #1 sawdust, 8/29/02; (air-dried and shredded to <8mm, then  
 fines <1mm, amounting to 38.3wt% of total, were removed before analysis)

Air-Dry Moisture, % 33.42

Proximate Analysis	(As received)	(As received)	
		w/SO3 correction	(Dry basis)
Moisture, %	39.23	39.23	
Volatile Matter, %	49.35	49.35	81.20
Ash (750°F), %	0.39	0.38	
Fixed Carbon, % (by difference)	11.03	11.04	

Ultimate Analysis	(Dry basis)	Screen Analysis	
		Retained on	Wt. %
Ash (750°F), %	0.63		
Carbon, %	49.51		
Hydrogen, %	5.88	3	0.0%
Nitrogen, %	0.11	4	0.2%
Sulfur, %	0.02	6	4.0%
Oxygen, %	43.85	8	16.3%
(by difference)		10	20.2%
		12	19.4%
		14	25.4%
Heating Value	(Dry basis)	16	8.6%
		PAN	<u>5.9%</u>
BTU/lb.	8,440	Total	100.0%

Analyst: NJP



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INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mount Prospect Road  
Des Plaines, IL 60018  
Attn: Norman Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Sawdust

Sample ID: 021478-002  
Wood Mill #1 Sawdust

Sample taken at -----

Sample taken by -----

Date sampled -----

Date received February 3, 2003

P.O No. PF00020103

Analysis Report No. 71-198921

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2700+	2655
Softening (ST)	2700+	2670
Hemispherical (HT)	2700+	2680
Fluid (FT)	2700+	2695

*looks backwards*

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory





Sample Login No: 021478-003

Date: December 18, 2002

Sample Description: Calla Energy, Mill #2 sawdust, 8/29/02; (air-dried and shredded to <8mm, then  
fines <1mm, amounting to 18.1wt% of total, were removed before analysis)

Air-Dry Moisture, % 21.83

Proximate Analysis	(As received)	(As received)	
		w/SO3 correction	(Dry basis)
Moisture, %	35.96	35.96	
Volatile Matter, %	53.21	53.21	83.09
Ash (750°F), %	0.27	0.26	
Fixed Carbon, % (by difference)	10.56	10.57	

Ultimate Analysis	(Dry basis)	Screen Analysis	
		Retained on	Wt. %
Ash (750°F), %	0.41		
Carbon, %	49.82		
Hydrogen, %	5.84	3	0.0%
Nitrogen, %	0.09	4	0.9%
Sulfur, %	0.01	6	7.2%
Oxygen, %	43.82	8	33.5%
(by difference)		10	19.5%
		12	14.4%
		14	15.9%
Heating Value	(Dry basis)	16	5.4%
		PAN	3.2%
BTU/lb.	8,430	Total	100.0%

Analyst: NJP



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www.comteco.com

INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mount Prospect Road  
Des Plaines, IL 60018  
Attn: Norman Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Sawdust

Sample ID: 021478-003  
Mill #2 Sawdust

Sample taken at -----

Sample taken by -----

Date sampled -----

Date received February 3, 2003

P.O No. PF00020103

Analysis Report No. 71-198922

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2655	2560
Softening (ST)	2670	2575
Hemispherical (HT)	2680	2590
Fluid (FT)	2690	2625

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory





Sample Login No: 021478-004

Date: December 18, 2002

Sample Description: Bush Ind. chipboard waste, composite of 2 containers; (shredded to <8mm, then fines <1mm, amounting to 25.4wt% of total, were removed before analysis)

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u>	
		<u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	5.53	5.53	
Volatile Matter, %	74.85	74.85	79.23
Ash (750°F), %	0.73	0.67	
Fixed Carbon, % (by difference)	18.89	18.96	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>	<b>Screen Analysis</b>	
		<u>Retained on</u>	<u>Wt. %</u>
Ash (750°F), %	0.71		
Carbon, %	48.90		
Hydrogen, %	6.04	3	0.0%
Nitrogen, %	3.02	4	3.5%
Sulfur, %	0.07	6	19.5%
Oxygen, %	41.27	8	28.5%
(by difference)		10	14.0%
		12	10.1%
		14	13.1%
		16	6.3%
		PAN	<u>5.0%</u>
		Total	100.0%

<b>Heating Value</b>	<u>(Dry basis)</u>
BTU/lb.	8,480

Analyst: NJP



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Des Plaines, IL 60018  
Attn: Norman Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Chipboard Waste

Sample ID: 021478-004  
Chipboard Waste

Sample taken at -----

Sample taken by -----

Date sampled -----

Date received February 3, 2003

P.O No. PF00020103

Analysis Report No. 71-198923

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2460	2560
Softening (ST)	2470	2570
Hemispherical (HT)	2485	2585
Fluid (FT)	2500	2595

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory





Sample Login No: 021478-005

Date: December 18, 2002

Sample Description: Bush Ind. wood pellets, 8/21/02

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received) w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	6.89	6.89	
Volatile Matter, %	73.29	73.29	78.71
Ash (750°F), %	0.84	0.76	
Fixed Carbon, % (by difference)	18.98	19.06	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>	<b>Screen Analysis</b>	
		<u>Retained on</u>	<u>Wt. %</u>
Ash (750°F), %	0.82		
Carbon, %	48.65		
Hydrogen, %	6.03	3/8"	0.0%
Nitrogen, %	3.75	3	89.0%
Sulfur, %	0.08	PAN	<u>11.0%</u>
Oxygen, % (by difference)	40.66	Total	100.0%

<b>Heating Value</b>	<u>(Dry basis)</u>
BTU/lb.	8,540

Analyst: NJP



# COMMERCIAL TESTING & ENGINEERING CO.

GENERAL OFFICES: 1919 SOUTH HIGHLAND AVE., SUITE 210-B, LOMBARD, ILLINOIS 60148 • TEL: 630-953-9300 FAX: 630-953-9306

SINCE 1908°



Member of the SGS Group (Société Générale de Surveillance)

February 6, 2003

ADDRESS ALL CORRESPONDENCE TO:  
16130 VAN DRUNEN ROAD  
SOUTH HOLLAND, IL 60473  
TEL: (708) 331-2900  
FAX: (708) 333-3060  
www.comteco.com

INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mount Prospect Road  
Des Plaines, IL 60018  
Attn: Norman Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Wood Pellets

Sample ID: 021478-005  
Wood Pellets  
Bush Industries

Sample taken at -----

Sample taken by -----

Date sampled -----

Date received February 3, 2003

P.O No. PF00020103

Analysis Report No. 71-198924

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2400	2380
Softening (ST)	2415	2405
Hemispherical (HT)	2430	2430
Fluid (FT)	2445	2465

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory

MEMBER  
**ACIL**





Sample Login No: 021506-001

Date: December 31, 2002

Sample Description: Wood pellets (from Menard's)

*Menard's*

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received) w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	<0.05	<0.05	
Volatile Matter, %	80.61	80.61	80.61
Ash (750°F), %	0.46	0.44	
Fixed Carbon, % (by difference)	18.93	18.95	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash (750°F), %	0.44
Carbon, %	47.84
Hydrogen, %	6.14
Nitrogen, %	0.10
Sulfur, %	0.01
Oxygen, % (by difference)	45.47

<b>Heating Value</b>	<u>(Dry basis)</u>
BTU/lb.	8,060

Analyst: NJP

*NJP binder*



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GENERAL OFFICES: 1919 SOUTH HIGHLAND AVE., SUITE 210-B, LOMBARD, ILLINOIS 60148 • TEL: 630-953-9300 FAX: 630-953-9306

SINCE 1908\*



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February 6, 2003

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16130 VAN DRUNEN ROAD  
SOUTH HOLLAND, IL 60473  
TEL: (708) 331-2900  
FAX: (708) 333-3060  
www.comteco.com

INSTITUTE OF GAS TECHNOLOGY  
1700 S. Mount Prospect Road  
Des Plaines, IL 60018  
Attn: Norman Petrulis

Sample identification by  
Institute of Gas Technology

Kind of sample reported to us Wood Pellets

Sample taken at -----

Sample ID: 021506-001  
Wood Pellets - Menard's

Sample taken by -----

Date sampled -----

Date received February 3, 2003

P.O No. PF00020103

Analysis Report No. 71-198920

Page 1 of 1

### FUSION TEMPERATURE OF ASH, (°F)

	<u>Reducing</u>	<u>Oxidizing</u>
Initial Deformation (IT)	2700+	2700+
Softening (ST)	2700+	2700+
Hemispherical (HT)	2700+	2700+
Fluid (FT)	2700+	2700+

#### METHOD

Fusion Temperature of Ash: ASTM D 1857

Respectfully submitted,  
COMMERCIAL TESTING & ENGINEERING CO.

South Holland Laboratory

MEMBER  
**ACIL**

# Customer Sample Request Certificate of Analysis

## CARBO CERAMICS, INC.

McIntyre Operations  
2295 Wriley Rd.  
McIntyre, GA 31054 U.S.A.  
ph. (478) 943-6500  
fax. (478) 943-6570  
972 401 0090  
972 401 0100

<b>Request To:</b>	Alan Conger
<b>Request Via:</b>	Steve Canova/Brandi Thayer
<b>Date Rec'd:</b>	12/3/2002
<b>Date Shipped:</b>	12/3/2002
<b>Product:</b>	CARBOPROP <span style="float: right;"><b>Size:</b> 30/60</span>
<b>Quantity:</b>	10 - 20 lbs
<b>Ship From Loc. :</b>	McIntyre
<b>Ship To :</b>	Andy Kramer
	Gas Technology Institute
	1700 South Mount Prospect Rd
	Desplaines, IL 60018

<b>Method of Shipment:</b>	Ground	Next day	2nd day
<b>Phone No.</b>	847-768-0612		
<b>Fax No.</b>			

**Purpose of use :** Fluidization Bed Media

\_\_\_\_\_

\_\_\_\_\_

\_\_\_\_\_

**SIEVE DISTRIBUTION**

S-1	20 mesh	0.0
S-2	30 mesh	0.0
S-3	40 mesh	80.9
S-4	50 mesh	18.6
S-5	60 mesh	0.5
S-6	70 mesh	0.0
S-7	100 mesh	0.0
	<b>Pan</b>	0.0
	<b>Total</b>	100.0

<input checked="" type="checkbox"/> <b>ASG</b>	3.23
<input checked="" type="checkbox"/> <b>BD (g/cc)</b>	1.84
<input type="checkbox"/> <b>API CRUSH @ 10 Kpsi</b>	_____
<input type="checkbox"/> <b>OTHER</b>	_____

**APPROVAL** Steve Canova



## Analytical Report

Batch # : 031306

July 11, 2003

**Prepared for:**

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 6/26/2003**

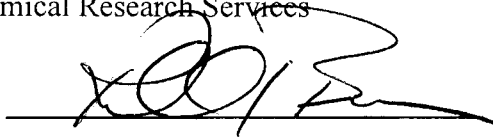
**Disclaimer:**

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Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report:

  
Russell Bora

(847) 768-0693



### Major Component Gas Analysis By Gas Chromatography

Report Date: 11-Jul-03

Client Name: 61129-08

GTI Sample Number: 031306-001

Sample Description: Gas Sample 1 6/20/2003

Date Analyzed: 27-Jun-03

Analyst: MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	12.7%	0.1%	0.94%
Carbon Dioxide	17.4%	0.03%	28.2%
Oxygen/Argon	0.88%	0.03%	1.05%
Nitrogen	52.6%	0.03%	54.2%
Carbon Monoxide	11.6%	0.03%	11.9%
Methane	3.82%	0.002%	2.25%
Ethane	0.080%	0.002%	0.088%
Ethene	0.689%	0.002%	0.711%
Ethyne		0.002%	
Propane		0.002%	
Propene	0.006%	0.002%	0.009%
Propadiene		0.002%	
Propyne		0.002%	
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene		0.002%	
i-Butene		0.002%	
trans-2-Butene		0.002%	
cis-2-Butene		0.002%	
1,3-Butadiene	0.003%	0.002%	0.006%
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
1-Pentene		0.002%	
Hexane Plus	0.173%	0.002%	0.549%
Hydrogen Sulfide	0.0164%	0.0001%	0.0206%
Carbonyl Sulfide	0.0013%	0.0001%	0.0029%
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

#### Calculated Real Gas Properties per ASTM D3588-98

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99924	0.99924
Compressibility Factor [z] (Sat.) =	0.99903	0.99903
Relative Density (Dry) =	0.9389	0.9389
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	138.0	138.3
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	135.6	135.9
Wobbe Index =	142.4	142.7
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	126.3	126.6
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	124.1	124.4

Notes: All blank values are below detection limit

N.A. - Not Analyzed



**TRACE SULFUR DETERMINATION BY ASTM D6228-98**

**Report Date:** 11-Jul-03  
**Client Name:** 61129-08  
**GTI Sample Number:** 031306-001  
**Sample Description:** Gas Sample 1 6/20/2003  
**Date Analyzed:** 27-Jun-03  
**Analyst:** MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	164	Thiophene	4.40
Sulfur Dioxide		C1-Thiophenes	0.07
Carbonyl Sulfide	13.2	C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan		Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide		Individual Unidentified	
Methyl Ethyl Sulfide		Sulfur Compounds	
Diethyl Sulfide		(all as monosulfides)	
Di-t-Butyl Sulfide			
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.00
i-Propyl n-Propyl Disulfide		Total Identified:	182
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	182
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	11.4
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	10.8
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

**Notes:** Component Detection Limit:  
 1 ppmv for H2S, COS, and SO2  
 0.05 ppmv for all other compounds per sulfur  
 All blank values are below detection limit.  
 STP= 14.696psia, 0°C




---

**Major Component Gas Analysis By Gas Chromatography**


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**Report Date:** 11-Jul-03**Client Name:** 61129-08**GTI Sample Number:** 031306-002**Sample Description:** Gas Sample 2 6/20/2003**Date Analyzed:** 27-Jun-03**Analyst:** MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	11.8%	0.1%	0.86%
Carbon Dioxide	17.9%	0.03%	28.7%
Oxygen/Argon	1.07%	0.03%	1.25%
Nitrogen	55.9%	0.03%	56.8%
Carbon Monoxide	9.37%	0.03%	9.53%
Methane	3.21%	0.002%	1.87%
Ethane	0.065%	0.002%	0.071%
Ethene	0.503%	0.002%	0.512%
Ethyne		0.002%	
Propane		0.002%	
Propene	0.004%	0.002%	0.006%
Propadiene		0.002%	
Propyne		0.002%	
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene		0.002%	
i-Butene		0.002%	
trans-2-Butene		0.002%	
cis-2-Butene		0.002%	
1,3-Butadiene		0.002%	
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
1-Pentene		0.002%	
Hexane Plus	0.122%	0.002%	0.383%
Hydrogen Sulfide	0.0137%	0.0001%	0.0169%
Carbonyl Sulfide	0.0012%	0.0001%	0.0026%
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

---

**Calculated Real Gas Properties per ASTM D3588-98**


---

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99925	0.99924
Compressibility Factor [z] (Sat.) =	0.99903	0.99903
Relative Density (Dry) =	0.9515	0.9515
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	116.0	116.3
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	114.0	114.3
Wobbe Index =	118.9	119.2
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	105.8	106.0
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	104.0	104.2

Notes: All blank values are below detection limit

N.A. - Not Analyzed



## TRACE SULFUR DETERMINATION BY ASTM D6228-98

Report Date: 11-Jul-03

Client Name: 61129-08

GTI Sample Number: 031306-002

Sample Description: Gas Sample 2 6/20/2003

Date Analyzed: 27-Jun-03

Analyst: MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	137	Thiophene	3.18
Sulfur Dioxide		C1-Thiophenes	0.05
Carbonyl Sulfide	11.7	C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan		Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide			
Methyl Ethyl Sulfide		Individual Unidentified	
Diethyl Sulfide		Sulfur Compounds	
Di-t-Butyl Sulfide		(all as monosulfides)	
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.00
i-Propyl n-Propyl Disulfide		Total Identified:	152
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	152
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	9.50
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	9.01
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

## Notes:

Component Detection Limit:

1 ppmv for H<sub>2</sub>S, COS, and SO<sub>2</sub>

0.05 ppmv for all other compounds per sulfur

All blank values are below detection limit.

STP= 14.696psia, 0°C





## Analytical Report

Batch #: 031307.doc

Date: July 18, 2003

**Prepared for:**

**Andrew Kramer  
X5612**

**Gas Technology Institute  
1700 S. Mt Prospect Rd.  
Des Plaines, IL 60018**

**Project # 61129-08**

**Received Date: 6/26/2003**

**Disclaimer:**

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In April 2000, the Institute of Gas Technology (IGT) and the Gas Research Institute (GRI) combined to form the Gas Technology Institute (GTI).

Submitted by: Karen Crippen, (847) 768-0604  
Chemical Research Services

Technical Contact:

A handwritten signature in black ink that reads "Alan G. Janos". The signature is written over a horizontal line.

Alan G. Janos, (847) 768-0603

## Analytical Report

Batch #: 031307.doc

Date: July 18, 2003

### Liquid Sample #1, 6/20/03 (031307-001)

#### Phase separation:

Organic phase (oil) recovered:	0.3 g	0.07 wt%
Aqueous phase recovered:	414.4 g	98.41 wt%
Solid phase recovered:	0.3 g	0.07 wt%
Loss (by difference):	<u>6.1 g</u>	1.45 wt%
Total sample received:	421.1 g	

Analyst: AGJ

#### Oil phase analyses (031307-003): (GC work only per A.K.)

#### Water phase analyses (031307-002):

Total Carbon:	2.41 wt%
Kjeldahl Nitrogen:	4.82 wt%
Ammonia:	5.52 wt%

Analyst: NJP, JS, AGJ



## Hydrocarbon Analysis in Oil Phase

**GTI Sample Number:** 031307-003

**Sample Description:** Liquid #1 6/20/03 Oil Phase

<u>Tentative Component ID</u>	<u>Estimated mg recovered</u>	<u>Tentative Component ID</u>	<u>Estimated mg recovered</u>
Hydroxymethylpentanone	0.548	Methoxypropenylphenol/Dimethylnaphthalenes	0.670
Ethylbenzene	0.083	Ethenylnaphthalene	0.836
m,p-Xylene	0.338	Acenaphthylene	5.39
o-Xylene	0.221	Methyl-1,1'-biphenyl	0.309
Phenol	0.815	Naphthalenecarbonitriles	1.03
Benzeneamine	0.077	Acenaphthene	0.402
Benzonitrile	0.357	Dibenzofuran	0.867
Methyl phenols/Indene	0.977	Phenylene	0.430
Methylbenzonitriles	0.161	Fluorene	1.73
Tetramethylpiperidinone	<0.020	Methyldibenzofuran	0.212
Methylbenzofuran	0.112	Methylfluorenes	0.535
Dimethylphenols	0.178	Fluorenone	0.370
Methylindenes	0.246	Phenanthrene	3.47
Methylbenzaldehydes	0.228	Anthracene	0.804
Naphthalene	5.18	Carbazole	0.130
Quinoline	2.01	Benzoquinoline	0.212
Isoquinoline	1.04	Phenylnaphthalenes	0.542
Indole	1.16	Benzo[c]cinnoline	0.198
2-Methylnaphthalene	1.40	Methylphenanthrene or Methylantracene	0.192
1-Methylnaphthalene	1.07	Methylphenanthrene or Methylantracene	0.198
Methylquinolines	0.200	4H-Cyclopenta[d,e,f]phenanthrene	0.517
Methylindoles	0.075	Fluoranthene	0.568
1,1'-Biphenyl	1.39	Pyrene	0.592
Ethyl-naphthalenes	0.221	Methylpyrenes	0.282
Dihydroacenaphthylene	0.147		
Indazole	<0.020	<b>Additional Unidentified Components</b>	<b>6.29</b>
		<b>Total :</b>	<b>45.0</b>

Notes: Detection Limit = 0.020 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.

## Analytical Report

Batch #: 031307.doc

Date: July 18, 2003

### Liquid Sample #2, 6/20/03 (031307-004)

#### Phase separation:

Organic phase (oil) recovered:	0.2 g	0.04 wt%
Aqueous phase recovered:	548.9 g	98.74 wt%
Solid phase recovered:	0.3 g	0.05 wt%
Loss (by difference):	<u>6.5 g</u>	1.17 wt%
Total sample received:	555.9 g	

Analyst: AGJ

#### Oil phase analyses (031307-006): (GC work only per A.K.)

#### Water phase analyses (031307-005):

Total Carbon:	2.18 wt%
Kjeldahl Nitrogen:	4.24 wt%
Ammonia:	4.93 wt%

Analyst: NJP, JS, AGJ



## Hydrocarbon Analysis in Oil Phase

**GTI Sample Number:** 031307-006

**Sample Description:** Liquid #2 6/20/03 Oil Phase

<b>Tentative Component ID</b>	<b>Estimated mg recovered</b>	<b>Tentative Component ID</b>	<b>Estimated mg recovered</b>
Hydroxymethylpentanone	0.213	Methoxypropenylphenol/Dimethylnaphthalene	1.14
Ethylbenzene	0.231	Ethenylnaphthalene	1.26
m,p-Xylene	0.923	Acenaphthylene	7.69
o-Xylene	0.507	Methyl-1,1'-biphenyl	0.483
Phenol	0.501	Naphthalenecarbonitriles	1.70
Benzeneamine	0.076	Acenaphthene	0.617
Benzonitrile	0.575	Dibenzofuran	1.28
Methyl phenols/Indene	1.08	Phenylene	0.686
Methylbenzotriles	0.191	Fluorene	2.47
Tetramethylpiperidinone	<0.020	Methylidibenzofuran	0.329
Methylbenzofuran	0.098	Methylfluorenes	1.14
Dimethylphenols	0.163	Fluorenone	0.531
Methylindenes	0.410	Phenanthrene	5.02
Methylbenzaldehydes	0.194	Anthracene	1.19
Naphthalene	9.02	Carbazole	0.215
Quinoline	3.46	Benzoquinoline	0.310
Isoquinoline	1.63	Phenylnaphthalenes	1.02
Indole	1.74	Benzo[c]cinnoline	0.302
2-Methylnaphthalene	2.40	Methylphenanthrene or Methylanthracene	0.297
1-Methylnaphthalene	1.78	Methylphenanthrene or Methylanthracene	0.324
Methylquinolines	0.296	4H-Cyclopenta[d,e,f]phenanthrene	0.836
Methylindoles	0.110	Fluoranthene	0.737
1,1'-Biphenyl	2.06	Pyrene	0.711
Ethyl-naphthalenes	0.350	Methylpyrenes	0.346
Dihydroacenaphthylene	0.185		
Indazole	<0.020	<b>Additional Unidentified Components</b>	<b>10.9</b>
		<b>Total :</b>	<b>59.1</b>

Notes: Detection Limit = 0.020 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.



Sample Login No: 031307-007  
Sample Description: Solids, 6/20/03

Date: July 15, 2003

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u> <u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	5.58	5.58	
Volatile Matter, %	16.82	16.82	17.81
Ash, %	45.91	45.84	
Fixed Carbon, % (by difference)	31.71	31.77	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash, %	48.54
Carbon, %	43.76
Hydrogen, %	0.46
Nitrogen, %	0.74
Sulfur, %	0.05
Oxygen, % (by difference)	6.44

Analyst: NJP



## Analytical Report

Batch # : 031308

August 06, 2003

**Prepared for:**

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 6/26/2003**

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Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report:

A handwritten signature in black ink, appearing to read "R. Bora", is written over a horizontal line.

Russell Bora

(847) 768-0693



### Major Component Gas Analysis By Gas Chromatography

Report Date: 11-Jul-03

Client Name: **61129-08**

GTI Sample Number: **031308-001**

Sample Description: **Gas Sample 1 6/24/2003**

Date Analyzed: 27-Jun-03

Analyst: MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	<b>9.5%</b>	0.1%	<b>0.68%</b>
Carbon Dioxide	<b>17.9%</b>	0.03%	<b>28.0%</b>
Oxygen/Argon	<b>0.80%</b>	0.03%	<b>0.92%</b>
Nitrogen	<b>57.4%</b>	0.03%	<b>57.1%</b>
Carbon Monoxide	<b>9.47%</b>	0.03%	<b>9.41%</b>
Methane	<b>3.49%</b>	0.002%	<b>1.99%</b>
Ethane	<b>0.185%</b>	0.002%	<b>0.197%</b>
Ethene	<b>0.895%</b>	0.002%	<b>0.891%</b>
Ethyne	<b>0.021%</b>	0.002%	<b>0.019%</b>
Propane	<b>0.009%</b>	0.002%	<b>0.014%</b>
Propene	<b>0.105%</b>	0.002%	<b>0.157%</b>
Propadiene		0.002%	
Propyne	<b>0.008%</b>	0.002%	<b>0.011%</b>
i-Butane		0.002%	
n-Butane		0.002%	
l-Butene		0.002%	
i-Butene	<b>0.005%</b>	0.002%	<b>0.011%</b>
trans-2-Butene	<b>0.004%</b>	0.002%	<b>0.007%</b>
cis-2-Butene	<b>0.002%</b>	0.002%	<b>0.004%</b>
1,3-Butadiene	<b>0.031%</b>	0.002%	<b>0.059%</b>
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
Pentenes	<b>0.004%</b>	0.002%	<b>0.011%</b>
Hexane Plus	<b>0.160%</b>	0.002%	<b>0.490%</b>
Hydrogen Sulfide	<b>0.0082%</b>	0.0001%	<b>0.0099%</b>
Carbonyl Sulfide	<b>0.0009%</b>	0.0001%	<b>0.0020%</b>
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

#### Calculated Real Gas Properties per ASTM D3588-98

Temp. (°F) =	<b>60.0</b>	<b>60.0</b>
Press. (psia) =	<b>14.696</b>	<b>14.73</b>
Compressibility Factor [z] (Dry) =	0.99919	0.99918
Compressibility Factor [z] (Sat.) =	0.99896	0.99896
Relative Density (Dry) =	0.9730	0.9730
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	126.2	126.5
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	124.0	124.3
Wobbe Index =	127.9	128.2
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	115.9	116.1
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	113.9	114.1

Notes: All blank values are below detection limit

N.A. - Not Analyzed





## TRACE SULFUR DETERMINATION BY ASTM D6228-98

Report Date: 11-Jul-03

Client Name: 61129-08

GTI Sample Number: 031308-001

Sample Description: Gas Sample 1 6/24/2003

Date Analyzed: 27-Jun-03

Analyst: MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	81.6	Thiophene	3.20
Sulfur Dioxide		C1-Thiophenes	0.40
Carbonyl Sulfide	9.34	C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan	1.11	Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide			
Methyl Ethyl Sulfide		Individual Unidentified	
Diethyl Sulfide		Sulfur Compounds	
Di-t-Butyl Sulfide		(all as monosulfides)	
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.00
i-Propyl n-Propyl Disulfide		Total Identified:	95.7
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	95.7
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	5.98
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	5.67
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

**Notes:** Component Detection Limit:  
 1 ppmv for H<sub>2</sub>S, COS, and SO<sub>2</sub>  
 0.05 ppmv for all other compounds per sulfur  
 All blank values are below detection limit.  
 STP= 14.696psia, 0°C



**Major Component Gas Analysis By Gas Chromatography**

**Report Date:** 11-Jul-03

**Client Name:** 61129-08

**GTI Sample Number:** 031308-002

**Sample Description:** Gas Sample 2 6/24/2003

**Date Analyzed:** 27-Jun-03

**Analyst:** MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	10.5%	0.1%	0.76%
Carbon Dioxide	18.5%	0.03%	29.1%
Oxygen/Argon	0.79%	0.03%	0.91%
Nitrogen	54.9%	0.03%	55.1%
Carbon Monoxide	9.72%	0.03%	9.74%
Methane	3.98%	0.002%	2.28%
Ethane	0.232%	0.002%	0.249%
Ethene	0.945%	0.002%	0.949%
Ethyne	0.016%	0.002%	0.015%
Propane	0.011%	0.002%	0.017%
Propene	0.128%	0.002%	0.192%
Propadiene	0.002%	0.002%	0.003%
Propyne	0.008%	0.002%	0.012%
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene	0.002%	0.002%	0.004%
i-Butene	0.007%	0.002%	0.013%
trans-2-Butene	0.004%	0.002%	0.008%
cis-2-Butene	0.003%	0.002%	0.005%
1,3-Butadiene	0.037%	0.002%	0.071%
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
Pentenes	0.005%	0.002%	0.013%
Hexane Plus	0.180%	0.002%	0.554%
Hydrogen Sulfide	0.0119%	0.0001%	0.0145%
Carbonyl Sulfide	0.0010%	0.0001%	0.0022%
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

**Calculated Real Gas Properties per ASTM D3588-98**

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99916	0.99916
Compressibility Factor [z] (Sat.) =	0.99893	0.99893
Relative Density (Dry) =	0.9654	0.9654
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	138.7	139.1
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	136.3	136.7
Wobbe Index =	141.2	141.5
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	127.2	127.5
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	125.0	125.3

Notes: All blank values are below detection limit  
 N.A. - Not Analyzed



**TRACE SULFUR DETERMINATION BY ASTM D6228-98**

**Report Date:** 11-Jul-03  
**Client Name:** 61129-08  
**GTI Sample Number:** 031308-002  
**Sample Description:** Gas Sample 2 6/24/2003  
**Date Analyzed:** 27-Jun-03  
**Analyst:** MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	119	Thiophene	3.23
Sulfur Dioxide		C1-Thiophenes	0.47
Carbonyl Sulfide	10.4	C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan	1.85	Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide		Individual Unidentified	
Methyl Ethyl Sulfide		Sulfur Compounds	
Diethyl Sulfide		(all as monosulfides)	
Di-t-Butyl Sulfide			
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.00
i-Propyl n-Propyl Disulfide		Total Identified:	135
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	135
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	8.44
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	8.00
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

**Notes:** Component Detection Limit:  
 1 ppmv for H2S, COS, and SO2  
 0.05 ppmv for all other compounds per sulfur  
 All blank values are below detection limit.  
 STP= 14.696psia, 0°C



### Hydrocarbon Analysis in Xylene Wash

GTI Sample Number: 031308-003

Sample Description: Xylene Cleanout 6/24/2003

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	4.83	Indazole	<2.12
i-Propylbenzene	2,820	Methoxypropenylphenol/Dimethylnaphthalenes	74.9
Phenol/n-Propylbenzene	1,750	Ethenylnaphthalene	95.9
Benzeneamine	<2.12	Acenaphthylene	323
Benzonitrile/1-Ethyl-3-methylbenzene	1,900	Methyl-1,1'-biphenyl	26.2
1-Ethyl-4-methylbenzene	501	Naphthalenecarbonitriles	91.0
1-Ethyl-2-methylbenzene	162	Acenaphthene	35.7
Trimethylbenzenes	319	Dibenzofuran	69.0
Methyl phenols/Indene	367	Phenylene	45.7
Methylbenzonitriles	27.1	Fluorene	164
Tetramethylpiperidinone	<2.12	Methyldibenzofuran	21.2
Methylbenzofuran	20.8	Methylfluorenes	101
Dimethylphenols	22.3	Fluorenone	41.1
Methylindenes	28.4	Phenanthrene	557
Methylbenzaldehydes	56.1	Anthracene	127
Naphthalene	632	Carbazole	17.3
Quinoline	57.8	Benzoquinoline	26.8
Isoquinoline	41.2	Phenylnaphthalenes	72.3
Indole	67.5	Benzo[c]cinnoline	22.1
2-Methylnaphthalene	210	Methylphenanthrene or Methylanthracene	22.0
1-Methylnaphthalene	149	Methylphenanthrene or Methylanthracene	25.4
Methylquinolines	14.4	4H-Cyclopenta[d,e,f]phenanthrene	71.6
Methylindoles	9.84	Fluoranthene	128
1,1'-Biphenyl	86.2	Pyrene	124
Ethyl-naphthalenes	28.3	Methylpyrenes	39.1
Dihydroacenaphthylene	19.3	<b>Additional Unidentified Components</b>	<b>1,200</b>
		<b>Total :</b>	<b>12,815</b>

Total Filtered Solids (Dry Weight): 1.2 g (0.13% wt.)

Notes: Detection Limit = 2.12 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Toluene and Ethylbenzene were found in large quantities but not calculated, based on the assumption that the source of these components were from the Xylene solvent.

Due to the lack of a solvent blank to identify additional solvent impurities, all additional components found have been reported.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.

Results are based on a total volume of 1060 mLs.



## Analytical Report

Batch #: 031309.doc  
Date: July 18, 2003

**Prepared for:**

**Andrew Kramer  
X5612**

**Gas Technology Institute  
1700 S. Mt Prospect Rd.  
Des Plaines, IL 60018**

**Project # 61129-08**

**Received Date: 6/26/2003**

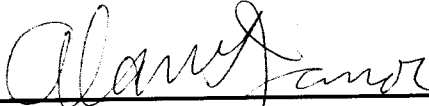
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In April 2000, the Institute of Gas Technology (IGT) and the Gas Research Institute (GRI) combined to form the Gas Technology Institute (GTI).

**Submitted by: Karen Crippen, (847) 768-0604  
Chemical Research Services**

**Technical Contact:**

  
\_\_\_\_\_  
Alan G. Janos, (847) 768-0603

## Analytical Report

Batch #: 031309.doc

Date: July 18, 2003

### Liquid Sample #1, 6/24/03 (031309-001)

#### Phase separation:

Organic phase (oil) recovered:	1.1 g	0.16 wt%
Aqueous phase recovered:	703.5 g	99.31 wt%
Solid phase recovered:	0.3 g	0.04 wt%
Loss (by difference):	<u>3.5 g</u>	0.49 wt%
Total sample received:	708.4 g	

Analyst: AGJ

#### Oil phase analyses (031309-003): (GC work only per A.K.)

#### Water phase analyses (031309-002):

Total Carbon:	2.80 wt%
Kjeldahl Nitrogen:	4.69 wt%
Ammonia:	5.67 wt%

Analyst: NJP, JS, AGJ



### Hydrocarbon Analysis in Oil Phase

GTI Sample Number: 031309-003

Sample Description: Liquid #1 6/24/03 Oil Phase

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	0.18	Methoxypropenylphenol/Dimethylnaphthalenes	8.78
Ethylbenzene	<0.09	Ethenylnaphthalene	10.7
m,p-Xylene	0.29	Acenaphthylene	34.8
o-Xylene	0.12	Methyl-1,1'-biphenyl	3.25
Phenol	27.9	Naphthalenecarbonitriles	12.7
Benzeneamine	0.82	Acenaphthene	4.08
Benzonitrile	1.26	Dibenzofuran	8.14
Methylphenols/Indene	24.3	Phenalene	5.30
Methylbenzonitriles	2.26	Fluorene	20.1
Tetramethylpiperidinone	<0.09	Methyldibenzofuran	2.40
Methylbenzofuran	1.39	Methylfluorenes	5.62
Dimethylphenols	2.80	Fluorenone	3.57
Methylindenes	3.52	Phenanthrene	56.7
Methylbenzaldehydes	6.53	Anthracene	10.6
Naphthalene	36.4	Carbazole	1.65
Quinoline	6.94	Benzoquinoline	2.39
Isoquinoline	5.20	Phenylnaphthalenes	6.98
Indole	8.16	Benzo[c]cinnoline	1.71
2-Methylnaphthalene	18.0	Methylphenanthrene or Methylanthracene	2.06
1-Methylnaphthalene	13.6	Methylphenanthrene or Methylanthracene	2.36
Methylquinolines	2.07	4H-Cyclopenta[d,e,f]phenanthrene	6.19
Methylindoles	1.36	Fluoranthene	7.84
1,1'-Biphenyl	9.12	Pyrene	7.14
Ethyl-naphthalenes	3.11	Methylpyrenes	2.32
Dihydroacenaphthylene	1.99		
Indazole	0.53	<b>Additional Unidentified Components</b>	<b>102</b>
		<b>Total :</b>	<b>507</b>

Notes: Detection Limit = 0.09 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.



Sample Login No: 031309-004  
Sample Description: Solids, 6/24/03

Date: July 15, 2003

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u> <u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	3.44	3.44	
Volatile Matter, %	12.99	12.99	13.45
Ash, %	33.38	33.18	
Fixed Carbon, % (by difference)	50.19	50.39	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash, %	34.36
Carbon, %	60.79
Hydrogen, %	1.04
Nitrogen, %	1.28
Sulfur, %	0.14
Oxygen, % (by difference)	2.38

Analyst: NJP



## Analytical Report

Batch # : 031324

August 06, 2003

**Prepared for:**

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 7/2/2003**

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Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report:

  
Russell Bora

(847) 768-0693



### Major Component Gas Analysis By Gas Chromatography

Report Date: 17-Jul-03

Client Name: 61129-08

GTI Sample Number: 031324-001

Sample Description: Gas Sample #1 6/26/2003 15:35

Date Analyzed: 15-Jul-03

Analyst: MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	7.0%	0.1%	0.50%
Carbon Dioxide	16.2%	0.03%	25.1%
Oxygen/Argon	1.37%	0.03%	1.56%
Nitrogen	56.5%	0.03%	55.7%
Carbon Monoxide	12.5%	0.03%	12.3%
Methane	4.52%	0.002%	2.55%
Ethane	0.160%	0.002%	0.169%
Ethene	1.30%	0.002%	1.28%
Ethyne	0.057%	0.002%	0.052%
Propane		0.002%	
Propene	0.033%	0.002%	0.048%
Propadiene		0.002%	
Propyne	0.004%	0.002%	0.006%
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene		0.002%	
i-Butene		0.002%	
trans-2-Butene		0.002%	
cis-2-Butene		0.002%	
1,3-Butadiene	0.015%	0.002%	0.029%
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
1-Pentene		0.002%	
Hexane Plus	0.220%	0.002%	0.665%
Hydrogen Sulfide	0.00225%	0.00005%	0.00270%
Carbonyl Sulfide		0.00005%	
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

#### Calculated Real Gas Properties per ASTM D3588-98

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99918	0.99918
Compressibility Factor [z] (Sat.) =	0.99895	0.99895
Relative Density (Dry) =	0.9824	0.9824
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	145.0	145.3
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	142.5	142.8
Wobbe Index =	146.3	146.6
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	134.5	134.8
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	132.2	132.5

Notes: All blank values are below detection limit

N.A. - Not Analyzed



**TRACE SULFUR DETERMINATION BY ASTM D6228-98**

**Report Date:** 17-Jul-03  
**Client Name:** 61129-08  
**GTI Sample Number:** 031324-001  
**Sample Description:** Gas Sample #1 6/26/2003 15:35  
**Date Analyzed:** 2-Jul-03  
**Analyst:** MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	22.5	Thiophene	1.38
Sulfur Dioxide		C1-Thiophenes	0.05
Carbonyl Sulfide		C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan		Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide			
Methyl Ethyl Sulfide		Individual Unidentified	
Diethyl Sulfide		Sulfur Compounds	
Di-t-Butyl Sulfide		(all as monosulfides)	
		Compound #1	<u>0.04</u>
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.04
i-Propyl n-Propyl Disulfide		Total Identified:	23.9
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	24.0
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	1.50
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	1.42
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

**Notes:** Component Detection Limit:  
 0.5 ppmv for H<sub>2</sub>S, COS, and SO<sub>2</sub>  
 0.05 ppmv for all other compounds per sulfur  
 Underlined numbers are below standard detection  
 limits, and are included for information only.  
 All blank values are below detection limit.  
 STP= 14.696psia, 0°C



8210 Mosley Rd.  
Houston, TX 77075  
713 943-9776 Telephone  
713 943-3846 Facsimile

# CORE LABORATORIES

KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-001  
Sample Date: 6/26/03 3:35:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 1 (06-26-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
<b>Nitrogen Compounds, Comprehensive</b>					
Ammonia	1	ppm v / v	GC-Chemiluminesci	7/24/03	JB
Nitric Oxide	< 1	ppm v / v			
Nitrous Oxide	< 1	ppm v / v			
Hydrogen Cyanide	< 1	ppm v / v			
Methyl amine	< 1	ppm v / v			
Dimethyl amine	< 1	ppm v / v			
Acetonitrile	< 1	ppm v / v			
Acrylonitrile	< 1	ppm v / v			
Ethyl amine	< 1	ppm v / v			
Diethyl amine	< 1	ppm v / v			
Isopropyl Amine	< 1	ppm v / v			
Propionitrile	2	ppm v / v			
Tert-Butyl Amine	< 1	ppm v / v			
Isobutyl nitrile	< 1	ppm v / v			
Butyl nitrile	< 1	ppm v / v			
Pyridine	< 1	ppm v / v			
Pyrrole	< 1	ppm v / v			
Pentane nitrile	< 1	ppm v / v			
2-Methyl pyridine	< 1	ppm v / v			
2-Methyl pyrrole	< 1	ppm v / v			
3-Methyl pyrrole	< 1	ppm v / v			
3 & 4-Methyl pyridine	< 1	ppm v / v			
2,6-Dimethyl pyridine	< 1	ppm v / v			
2-Ethyl pyridine	< 1	ppm v / v			
2-Ethyl pyrrole	< 1	ppm v / v			
2,4-Dimethyl pyridine	< 1	ppm v / v			
2,4 Dimethyl pyrrole	< 1	ppm v / v			

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# CORE LABORATORIES

KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-001  
Sample Date: 6/26/03 3:35:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 1 (06-26-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
2,3-Dimethyl pyridine	< 1	ppm v / v			
4-Ethyl pyridine	< 1	ppm v / v			
3,5-Dimethyl pyridine	< 1	ppm v / v			
Aniline	< 1	ppm v / v			
Unidentified Nitrogen	< 1	ppm v / v			

Approved By:

Jean Waits  
Supervising Chemist

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**Major Component Gas Analysis By Gas Chromatography**


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**Report Date:** 17-Jul-03

**Client Name:** 61129-08

**GTI Sample Number:** 031324-002

**Sample Description:** Gas sample #2 6/26/2003 17:25

**Date Analyzed:** 15-Jul-03

**Analyst:** MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	8.1%	0.1%	0.58%
Carbon Dioxide	16.4%	0.03%	25.7%
Oxygen/Argon	1.42%	0.03%	1.63%
Nitrogen	56.2%	0.03%	55.8%
Carbon Monoxide	12.0%	0.03%	11.9%
Methane	4.57%	0.002%	2.60%
Ethane	0.11%	0.002%	0.12%
Ethene	0.92%	0.002%	0.92%
Ethyne	0.021%	0.002%	0.019%
Propane		0.002%	
Propene	0.007%	0.002%	0.010%
Propadiene		0.002%	
Propyne		0.002%	
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene		0.002%	
i-Butene		0.002%	
trans-2-Butene		0.002%	
cis-2-Butene		0.002%	
1,3-Butadiene	0.005%	0.002%	0.010%
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
1-Pentene		0.002%	
Hexane Plus	0.216%	0.002%	0.661%
Hydrogen Sulfide	0.00176%	0.00005%	0.00213%
Carbonyl Sulfide		0.00005%	
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

**Calculated Real Gas Properties per ASTM D3588-98**

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99920	0.99920
Compressibility Factor [z] (Sat.) =	0.99898	0.99898
Relative Density (Dry) =	0.9735	0.9735
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	138.7	139.0
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	136.3	136.6
Wobbe Index =	140.5	140.9
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	128.1	128.4
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	125.9	126.2

Notes: All blank values are below detection limit

N.A. - Not Analyzed



**TRACE SULFUR DETERMINATION BY ASTM D6228-98**

**Report Date:** 17-Jul-03  
**Client Name:** 61129-08  
**GTI Sample Number:** 031324-002  
**Sample Description:** Gas sample #2 6/26/2003 17:25  
**Date Analyzed:** 2-Jul-03  
**Analyst:** MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	17.6	Thiophene	0.90
Sulfur Dioxide		C1-Thiophenes	
Carbonyl Sulfide		C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan		Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide		Individual Unidentified	
Methyl Ethyl Sulfide		Sulfur Compounds	
Diethyl Sulfide		(all as monosulfides)	
Di-t-Butyl Sulfide		Compound #1	<u>0.03</u>
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.03
i-Propyl n-Propyl Disulfide		Total Identified:	18.5
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	18.5
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	1.16
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	1.10
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

**Notes:** Component Detection Limit:  
 0.5 ppmv for H2S, COS, and SO2  
 0.05 ppmv for all other compounds per sulfur  
 Underlined numbers are below standard detection limits, and are included for information only.  
 All blank values are below detection limit.  
 STP= 14.696psia, 0°C



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# CORE LABORATORIES

KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-002  
Sample Date: 6/26/03 5:25:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 2 (06-26-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
<b>Nitrogen Compounds, Comprehensive</b>					
Ammonia	1	ppm v / v	GC-Chemiluminesci	7/24/03	JB
Nitric Oxide	< 1	ppm v / v			
Nitrous Oxide	< 1	ppm v / v			
Hydrogen Cyanide	< 1	ppm v / v			
Methyl amine	< 1	ppm v / v			
Dimethyl amine	< 1	ppm v / v			
Acetonitrile	< 1	ppm v / v			
Acrylonitrile	< 1	ppm v / v			
Ethyl amine	< 1	ppm v / v			
Diethyl amine	< 1	ppm v / v			
Isopropyl Amine	< 1	ppm v / v			
Propionitrile	< 1	ppm v / v			
Tert-Butyl Amine	< 1	ppm v / v			
Isobutyl nitrile	< 1	ppm v / v			
Butyl nitrile	< 1	ppm v / v			
Pyridine	< 1	ppm v / v			
Pyrrole	< 1	ppm v / v			
Pentane nitrile	< 1	ppm v / v			
2-Methyl pyridine	< 1	ppm v / v			
2-Methyl pyrrole	< 1	ppm v / v			
3-Methyl pyrrole	< 1	ppm v / v			
3 & 4-Methyl pyridine	< 1	ppm v / v			
2,6-Dimethyl pyridine	< 1	ppm v / v			
2-Ethyl pyridine	< 1	ppm v / v			
2-Ethyl pyrrole	< 1	ppm v / v			
2,4-Dimethyl pyridine	< 1	ppm v / v			
2,4 Dimethyl pyrrole	< 1	ppm v / v			

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# CORE LABORATORIES

KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-002  
Sample Date: 6/26/03 5:25:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 2 (06-26-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
2,3-Dimethyl pyridine	< 1	ppm v / v			
4-Ethyl pyridine	< 1	ppm v / v			
3,5-Dimethyl pyridine	< 1	ppm v / v			
Aniline	< 1	ppm v / v			
Unidentified Nitrogen	< 1	ppm v / v			

Approved By:

Jean Waits  
Supervising Chemist



## Hydrocarbon Analysis in Xylene Wash

GTI Sample Number: 031324-003

Sample Description: Xylene Wash 6/26/2003

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	6.73	Indazole	<3.52
i-Propylbenzene	4,860	Methoxypropenylphenol/Dimethylnaphthalenes	66.2
Phenol/n-Propylbenzene	2,430	Ethenylnaphthalene	131
Benzeneamine	<3.52	Acenaphthylene	757
Benzonitrile/1-Ethyl-3-methylbenzene	3,930	Methyl-1,1'-biphenyl	36.5
1-Ethyl-4-methylbenzene	127	Naphthalenecarbonitriles	45.1
1-Ethyl-2-methylbenzene	260	Acenaphthene	53.9
Trimethylbenzenes	605	Dibenzofuran	99.6
Methyl phenols/Indene	319	Phenalene	59.8
Methylbenzonitriles	3.94	Fluorene	267
Tetramethylpiperidinone	<3.52	Methyldibenzofuran	25.0
Methylbenzofuran	13.6	Methylfluorenes	81.0
Dimethylphenols	6.16	Fluorenone	62.1
Methylindenes	44.3	Phenanthrene	625
Methylbenzaldehydes	22.8	Anthracene	174
Naphthalene	1,900	Carbazole	15.9
Quinoline	22.0	Benzoquinoline	21.3
Isoquinoline	15.0	Phenylnaphthalenes	96.2
Indole	20.1	Benzo[e]cinnoline	19.4
2-Methylnaphthalene	274	Methylphenanthrene or Methylanthracene	30.8
1-Methylnaphthalene	181	Methylphenanthrene or Methylanthracene	34.0
Methylquinolines	4.89	4H-Cyclopenta[d,e,f]phenanthrene	107
Methylindoles	<3.52	Fluoranthene	133
1,1'-Biphenyl	184	Pyrene	136
Ethylnaphthalenes	18.3	Methylpyrenes	65.6
Dihydroacenaphthylene	20.9	<b>Additional Unidentified Components</b>	<b>2,045</b>
		<b>Total :</b>	<b>20,456</b>

Total Filtered Solids (Dry Weight): 1.0 g (0.066% wt.)

Notes: Detection Limit = 3.52 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Toluene and Ethylbenzene were found in large quantities but not calculated, based on the assumption that the source of these components were from the Xylene solvent.

Due to the lack of a solvent blank to identify additional solvent impurities, all additional components found have been reported.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.

Results are based on a total volume of 1760 mLs.



## Analytical Report

Batch #: 031326.doc

Date: July 22, 2003

**Prepared for:**

**Andrew Kramer  
X5612**

**Gas Technology Institute  
1700 S. Mt Prospect Rd.  
Des Plaines, IL 60018**

**Project # 61129-08**

**Received Date: 7/2/2003**

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In April 2000, the Institute of Gas Technology (IGT) and the Gas Research Institute (GRI) combined to form the Gas Technology Institute (GTI).

Submitted by: Karen Crippen, (847) 768-0604  
Chemical Research Services

Technical Contact:

A handwritten signature in black ink, appearing to read "Alan G. Janos", written over a horizontal line.

Alan G. Janos, (847) 768-0603

## Analytical Report

Batch #: 031326.doc

Date: July 22, 2003

### Liquid Sample #1, 6/26/03 (031326-001)

#### Phase separation:

Organic phase (oil) recovered:	3.5 g	0.60 wt%
Aqueous phase recovered:	568.9 g	98.51 wt%
Solid phase recovered:	0.3 g	0.05 wt%
Loss (by difference):	<u>4.8 g</u>	0.83 wt%
Total sample received:	577.5 g	

Analyst: AGJ

#### Oil phase analyses (031326-003): (GC work only per A.K.)

#### Water phase analyses (031326-002):

Total Carbon: 0.38 wt%

Kjeldahl Nitrogen: 0.49 wt%

Ammonia: 0.28 wt%

Analyst: NJP, JS, AGJ



## Hydrocarbon Analysis in Oil Phase

GTI Sample Number: 031326-003

Sample Description: Liquid #1 6/26/03 Oil Phase

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	<0.317	Methoxypropenylphenol/Dimethylnaphthalenes	19.5
Ethylbenzene	<0.317	Ethenylnaphthalene	42.4
m,p-Xylene	<0.317	Acenaphthylene	235
o-Xylene	<0.317	Methyl-1,1'-biphenyl	11.5
Phenol	8.72	Naphthalenecarbonitriles	14.2
Benzeneamine	<0.317	Acenaphthene	16.5
Benzonitrile	<0.317	Dibenzofuran	32.9
Methylphenols/Indene	10.2	Phenalene	17.5
Methylbenzonitriles	0.564	Fluorene	97.8
Tetramethylpiperidinone	<0.317	Methyldibenzofuran	6.51
Methylbenzofuran	0.577	Methylfluorenes	18.8
Dimethylphenols	1.62	Fluorenone	8.88
Methylindenes	2.52	Phenanthrene	261
Methylbenzaldehydes	5.95	Anthracene	75.2
Naphthalene	234	Carbazole	3.12
Quinoline	6.01	Benzoquinoline	3.96
Isoquinoline	4.61	Phenylnaphthalenes	20.1
Indole	5.19	Benzo[c]innoline	3.40
2-Methylnaphthalene	72.2	Methylphenanthrene or Methylanthracene	6.52
1-Methylnaphthalene	49.4	Methylphenanthrene or Methylanthracene	6.87
Methylquinolines	3.18	4H-Cyclopenta[d,e,f]phenanthrene	21.9
Methylindoles	0.729	Fluoranthene	24.2
1,1'-Biphenyl	58.0	Pyrene	25.6
Ethyl-naphthalenes	5.94	Methylpyrenes	11.4
Dihydroacenaphthylene	6.67		
Indazole	<0.317	<b>Additional Unidentified Components</b>	<b>204</b>
		<b>Total :</b>	<b>1665</b>

Notes: Detection Limit = 0.317 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.

## Analytical Report

Batch #: 031326.doc

Date: July 22, 2003

### Liquid Sample #2, 6/26/03 (031326-005)

#### Phase separation:

Organic phase (oil) recovered:	2.7 g	0.29 wt%
Aqueous phase recovered:	924.0 g	98.70 wt%
Solid phase recovered:	0.2 g	0.02 wt%
Loss (by difference):	<u>9.3 g</u>	0.99 wt%
Total sample received:	936.2 g	

Analyst: AGJ

#### Oil phase analyses (031326-007): (GC work only per A.K.)

#### Water phase analyses (031326-006):

Total Carbon:	0.28 wt%
Kjeldahl Nitrogen:	0.30 wt%
Ammonia:	0.21 wt%

Analyst: NJP, JS, AGJ

## Hydrocarbon Analysis in Oil Phase

GTI Sample Number: 031326-007

Sample Description: Liquid #2 6/26/03 Oil Phase

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	<0.238	Methoxypropenylphenol/Dimethylnaphthalenes	13.9
Ethylbenzene	<0.238	Ethenylnaphthalene	29.4
m,p-Xylene	<0.238	Acenaphthylene	170
o-Xylene	<0.238	Methyl-1,1'-biphenyl	8.06
Phenol	7.53	Naphthalenecarbonitriles	11.0
Benzeneamine	<0.238	Acenaphthene	11.5
Benzonitrile	0.501	Dibenzofuran	22.4
Methylphenols/Indene	16.3	Phenalene	11.7
Methylbenzonitriles	0.518	Fluorene	66.0
Tetramethylpiperidinone	<0.238	Methyldibenzofuran	4.14
Methylbenzofuran	0.997	Methylfluorenes	13.2
Dimethylphenols	1.02	Fluorenone	9.45
Methylindenes	2.94	Phenanthrene	187
Methylbenzaldehydes	4.05	Anthracene	50.4
Naphthalene	295	Carbazole	2.06
Quinoline	4.27	Benzoquinoline	2.55
Isoquinoline	3.13	Phenylnaphthalenes	12.6
Indole	4.00	Benzo[c]cinnoline	2.20
2-Methylnaphthalene	56.1	Methylphenanthrene or Methylanthracene	4.29
1-Methylnaphthalene	37.5	Methylphenanthrene or Methylanthracene	4.50
Methylquinolines	2.07	4H-Cyclopenta[d,e,f]phenanthrene	14.5
Methylindoles	0.565	Fluoranthene	16.6
1,1'-Biphenyl	41.6	Pyrene	17.3
Ethyl-naphthalenes	4.59	Methylpyrenes	9.57
Dihydroacenaphthylene	4.51		
Indazole	<0.238	<b>Additional Unidentified Components</b>	<b>139</b>
		<b>Total :</b>	<b>1321</b>

Notes: Detection Limit = 0.238 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.



Sample Login No: 031326-008  
Sample Description: Carbon ash fines, 6/26/03

Date: July 15, 2003

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received) w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	5.92	5.92	
Volatile Matter, %	15.02	15.02	15.96
Ash, %	25.64	25.57	
Fixed Carbon, % (by difference)	53.43	53.49	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash, %	27.18
Carbon, %	65.09
Hydrogen, %	0.52
Nitrogen, %	0.23
Sulfur, %	0.02
Oxygen, % (by difference)	6.96

Analyst: NJP





## Analytical Report

Batch # : 031327

August 06, 2003

### Prepared for:

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 7/2/2003**

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Submitted by: **Karen Crippen, (847) 768-0604**

**Chemical Research Services**

Technical contact for this report:

A handwritten signature in black ink, appearing to read "RB", is written over a horizontal line.

Russell Bora

(847) 768-0693




---

**Major Component Gas Analysis By Gas Chromatography**


---

**Report Date:** 17-Jul-03**Client Name:** 61129-08**GTI Sample Number:** 031327-001**Sample Description:** Gas sample #1 6/30/2003 16:00**Date Analyzed:** 15-Jul-03      **Analyst:** MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	11.4%	0.1%	0.83%
Carbon Dioxide	18.9%	0.03%	30.0%
Oxygen/Argon	1.23%	0.03%	1.43%
Nitrogen	51.2%	0.03%	51.8%
Carbon Monoxide	10.8%	0.03%	10.9%
Methane	4.64%	0.002%	2.69%
Ethane	0.24%	0.002%	0.26%
Ethene	1.14%	0.002%	1.16%
Ethyne	0.010%	0.002%	0.009%
Propane	0.010%	0.002%	0.016%
Propene	0.123%	0.002%	0.187%
Propadiene		0.002%	
Propyne	0.006%	0.002%	0.008%
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene	0.003%	0.002%	0.006%
i-Butene	0.006%	0.002%	0.012%
trans-2-Butene	0.005%	0.002%	0.010%
cis-2-Butene	0.003%	0.002%	0.007%
1,3-Butadiene	0.030%	0.002%	0.058%
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
Pentenes	0.003%	0.002%	0.009%
Hexane Plus	0.199%	0.002%	0.617%
Hydrogen Sulfide	0.00232%	0.00005%	0.00285%
Carbonyl Sulfide		0.00005%	
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

---

**Calculated Real Gas Properties per ASTM D3588-98**


---

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99913	0.99913
Compressibility Factor [z] (Sat.) =	0.99890	0.99890
Relative Density (Dry) =	0.9580	0.9580
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	155.4	155.8
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	152.8	153.1
Wobbe Index =	158.8	159.2
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	142.5	142.8
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	140.1	140.4

Notes: All blank values are below detection limit

N.A. - Not Analyzed



**TRACE SULFUR DETERMINATION BY ASTM D6228-98**

**Report Date:** 17-Jul-03

**Client Name:** 61129-08

**GTI Sample Number:** 031327-001

**Sample Description:** Gas sample #1 6/30/2003 16:00

**Date Analyzed:** 2-Jul-03

**Analyst:** MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	23.2	Thiophene	1.17
Sulfur Dioxide		C1-Thiophenes	0.13
Carbonyl Sulfide		C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan		Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide		Individual Unidentified	
Methyl Ethyl Sulfide		Sulfur Compounds	
Diethyl Sulfide		(all as monosulfides)	
Di-t-Butyl Sulfide			
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.00
i-Propyl n-Propyl Disulfide		Total Identified:	24.5
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	24.5
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	1.53
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	1.45
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

**Notes:** Component Detection Limit:  
 0.5 ppmv for H2S, COS, and SO2  
 0.05 ppmv for all other compounds per sulfur  
 All blank values are below detection limit.  
 STP= 14.696psia, 0°C



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# CORE LABORATORIES

KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-003  
Sample Date: 6/30/03 4:00:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 1 (06-30-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
<b>Nitrogen Compounds, Comprehensive</b>					
Ammonia	1	ppm v / v	GC-Chemiluminesci	7/24/03	JB
Nitric Oxide	< 1	ppm v / v			
Nitrous Oxide	< 1	ppm v / v			
Hydrogen Cyanide	< 1	ppm v / v			
Methyl amine	< 1	ppm v / v			
Dimethyl amine	< 1	ppm v / v			
Acetonitrile	1	ppm v / v			
Acrylonitrile	< 1	ppm v / v			
Ethyl amine	< 1	ppm v / v			
Diethyl amine	< 1	ppm v / v			
Isopropyl Amine	< 1	ppm v / v			
Propionitrile	2	ppm v / v			
Tert-Butyl Amine	< 1	ppm v / v			
Isobutyl nitrile	< 1	ppm v / v			
Butyl nitrile	< 1	ppm v / v			
Pyridine	< 1	ppm v / v			
Pyrrole	< 1	ppm v / v			
Pentane nitrile	< 1	ppm v / v			
2-Methyl pyridine	< 1	ppm v / v			
2-Methyl pyrrole	< 1	ppm v / v			
3-Methyl pyrrole	< 1	ppm v / v			
3 & 4-Methyl pyridine	< 1	ppm v / v			
2,6-Dimethyl pyridine	< 1	ppm v / v			
2-Ethyl pyridine	< 1	ppm v / v			
2-Ethyl pyrrole	< 1	ppm v / v			
2,4-Dimethyl pyridine	< 1	ppm v / v			
2,4 Dimethyl pyrrole	< 1	ppm v / v			

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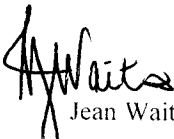
KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-003  
Sample Date: 6/30/03 4:00:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 1 (06-30-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
2,3-Dimethyl pyridine	< 1	ppm v / v			
4-Ethyl pyridine	< 1	ppm v / v			
3,5-Dimethyl pyridine	< 1	ppm v / v			
Aniline	< 1	ppm v / v			
Unidentified Nitrogen	< 1	ppm v / v			

Approved By:

  
Jean Waits  
Supervising Chemist

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**Major Component Gas Analysis By Gas Chromatography**


---

**Report Date:** 17-Jul-03**Client Name:** 61129-08**GTI Sample Number:** 031327-002**Sample Description:** Gas sample #2 6/30/2003 16:45**Date Analyzed:** 15-Jul-03**Analyst:** MAD

Component	Mol %	Det. Limit	Weight %
Helium		0.1%	
Hydrogen	10.9%	0.1%	0.79%
Carbon Dioxide	18.8%	0.03%	29.7%
Oxygen/Argon	1.13%	0.03%	1.31%
Nitrogen	52.6%	0.03%	52.8%
Carbon Monoxide	10.4%	0.03%	10.4%
Methane	4.44%	0.002%	2.55%
Ethane	0.22%	0.002%	0.24%
Ethene	1.10%	0.002%	1.11%
Ethyne	0.018%	0.002%	0.017%
Propane	0.009%	0.002%	0.014%
Propene	0.116%	0.002%	0.176%
Propadiene	0.002%	0.002%	0.003%
Propyne	0.007%	0.002%	0.011%
i-Butane		0.002%	
n-Butane		0.002%	
1-Butene	0.002%	0.002%	0.004%
i-Butene	0.006%	0.002%	0.011%
trans-2-Butene	0.004%	0.002%	0.007%
cis-2-Butene	0.003%	0.002%	0.005%
1,3-Butadiene	0.033%	0.002%	0.064%
i-Pentane		0.002%	
n-Pentane		0.002%	
neo-Pentane		0.002%	
Pentenes	0.004%	0.002%	0.010%
Hexane Plus	0.204%	0.002%	0.630%
Hydrogen Sulfide	0.00242%	0.00005%	0.00296%
Carbonyl Sulfide		0.00005%	
<b>Total</b>	<b>100.0%</b>		<b>100.0%</b>

---

**Calculated Real Gas Properties per ASTM D3588-98**


---

Temp. (°F) =	60.0	60.0
Press. (psia) =	14.696	14.73
Compressibility Factor [z] (Dry) =	0.99913	0.99913
Compressibility Factor [z] (Sat.) =	0.99891	0.99890
Relative Density (Dry) =	0.9626	0.9626
Gross HV (Dry) (Btu/ft <sup>3</sup> ) =	149.8	150.2
Gross HV (Sat.) (Btu/ft <sup>3</sup> ) =	147.3	147.6
Wobbe Index =	152.7	153.1
Net HV (Dry) (Btu/ft <sup>3</sup> ) =	137.4	137.7
Net HV (Sat.) (Btu/ft <sup>3</sup> ) =	135.0	135.4

Notes: All blank values are below detection limit

N.A. - Not Analyzed



## TRACE SULFUR DETERMINATION BY ASTM D6228-98

Report Date: 17-Jul-03

Client Name: 61129-08

GTI Sample Number: 031327-002

Sample Description: Gas sample #2 6/30/2003 16:45

Date Analyzed: 2-Jul-03

Analyst: MAD

Component Name	PPMV	Component Name	PPMV
Hydrogen Sulfide	24.2	Thiophene	1.03
Sulfur Dioxide		C1-Thiophenes	0.10
Carbonyl Sulfide		C2-Thiophenes	
Carbon Disulfide		C3-Thiophenes	
Methyl Mercaptan	0.05	Benzothiophene	
Ethyl Mercaptan		C1-Benzothiophenes	
i-Propyl Mercaptan		C2-Benzothiophenes	
n-Propyl Mercaptan			
t-Butyl Mercaptan		Thiophane	
		Thiophenol	
Dimethyl Sulfide		Individual Unidentified	
Methyl Ethyl Sulfide		Sulfur Compounds	
Diethyl Sulfide		(all as monosulfides)	
Di-t-Butyl Sulfide			
Dimethyl Disulfide			
Methyl Ethyl Disulfide			
Methyl i-Propyl Disulfide			
Diethyl Disulfide			
Methyl n-Propyl Disulfide			
Methyl t-Butyl Disulfide			
Ethyl i-Propyl Disulfide			
Ethyl n-Propyl Disulfide			
Ethyl t-Butyl Disulfide			
Di-i-Propyl Disulfide		Total Unidentified:	0.00
i-Propyl n-Propyl Disulfide		Total Identified:	25.4
Di-n-Propyl Disulfide			
i-Propyl t-Butyl Disulfide		<b>Total Sulfur Content</b>	
n-Propyl t-Butyl Disulfide		As molar PPM	25.4
Di-t-Butyl Disulfide		As Grains/100 SCF @ STP	1.59
Dimethyl Trisulfide		As Grains/100 SCF @ 14.73	1.50
Diethyl Trisulfide		psia, 60°F	
Di-t-Butyl Trisulfide			

## Notes:

Component Detection Limit:

0.5 ppmv for H<sub>2</sub>S, COS, and SO<sub>2</sub>

0.05 ppmv for all other compounds per sulfur

All blank values are below detection limit.

STP= 14.696psia, 0°C



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# CORE LABORATORIES

KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-004  
Sample Date: 6/30/03 4:45:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 2 (06-30-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
<b>Nitrogen Compounds, Comprehensive</b>					
Ammonia	1	ppm v / v	GC-Chemiluminesci	7/24/03	JB
Nitric Oxide	< 1	ppm v / v			
Nitrous Oxide	< 1	ppm v / v			
Hydrogen Cyanide	< 1	ppm v / v			
Methyl amine	< 1	ppm v / v			
Dimethyl amine	< 1	ppm v / v			
Acetonitrile	< 1	ppm v / v			
Acrylonitrile	< 1	ppm v / v			
Ethyl amine	< 1	ppm v / v			
Diethyl amine	< 1	ppm v / v			
Isopropyl Amine	< 1	ppm v / v			
Propionitrile	2	ppm v / v			
Tert-Butyl Amine	< 1	ppm v / v			
Isobutyl nitrile	< 1	ppm v / v			
Butyl nitrile	< 1	ppm v / v			
Pyridine	< 1	ppm v / v			
Pyrrole	< 1	ppm v / v			
Pentane nitrile	< 1	ppm v / v			
2-Methyl pyridine	< 1	ppm v / v			
2-Methyl pyrrole	< 1	ppm v / v			
3-Methyl pyrrole	< 1	ppm v / v			
3 & 4-Methyl pyridine	< 1	ppm v / v			
2,6-Dimethyl pyridine	< 1	ppm v / v			
2-Ethyl pyridine	< 1	ppm v / v			
2-Ethyl pyrrole	< 1	ppm v / v			
2,4-Dimethyl pyridine	< 1	ppm v / v			
2,4 Dimethyl pyrrole	< 1	ppm v / v			

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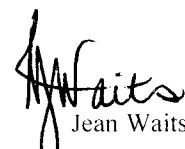
KAREN CRIPPEN  
GAS TECHNOLOGY INSTITUTE  
1700 S MOUNT PROSPECT RD  
DES PLAINES, IL 60018

Sample Number: 132956-004  
Sample Date: 6/30/03 4:45:00 PM  
Date Reported: 7/24/03  
Date Received: 7/18/03  
Sample ID: Sample # 2 (06-30-03)  
Description:

## Analytical Report

Test	Result	Units	Method	Date	Analyst
2,3-Dimethyl pyridine	< 1	ppm v / v			
4-Ethyl pyridine	< 1	ppm v / v			
3,5-Dimethyl pyridine	< 1	ppm v / v			
Aniline	< 1	ppm v / v			
Unidentified Nitrogen	< 1	ppm v / v			

Approved By:

  
Jean Waits  
Supervising Chemist

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## Hydrocarbon Analysis in Xylene Wash

GTI Sample Number: 031327-003

Sample Description: Xylene Wash 6/30/2003

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	7.07	Indazole	<1.84
i-Propylbenzene	5,380	Methoxypropenylphenol/Dimethylnaphthalenes	41.5
Phenol/n-Propylbenzene	2,740	Ethenylnaphthalene	57.7
Benzeneamine	<1.84	Acenaphthylene	356
Benzonitrile/1-Ethyl-3-methylbenzene	4,290	Methyl-1,1'-biphenyl	19.4
1-Ethyl-4-methylbenzene	206	Naphthalenecarbonitriles	16.6
1-Ethyl-2-methylbenzene	284	Acenaphthene	44.4
Trimethylbenzenes	632	Dibenzofuran	55.5
Methyl phenols/Indene	202	Phenylene	30.5
Methylbenzotriles	<1.84	Fluorene	140
Tetramethylpiperidinone	<1.84	Methylidibenzofuran	11.7
Methylbenzofuran	12.6	Methylfluorenes	34.1
Dimethylphenols	15.7	Fluorenone	20.7
Methylindenes	35.1	Phenanthrene	368
Methylbenzaldehydes	23.0	Anthracene	85.9
Naphthalene	739	Carbazole	6.33
Quinoline	8.54	Benzoquinoline	7.99
Isoquinoline	7.25	Phenylnaphthalenes	48.0
Indole	11.2	Benzo[e]cinoline	8.85
2-Methylnaphthalene	175	Methylphenanthrene or Methylanthracene	15.7
1-Methylnaphthalene	120	Methylphenanthrene or Methylanthracene	16.4
Methylquinolines	<1.84	4H-Cyclopenta[d,e,f]phenanthrene	52.8
Methylindoles	<1.84	Fluoranthene	97.6
1,1'-Biphenyl	101	Pyrene	111
Ethylnaphthalenes	20.3	Methylpyrenes	48.0
Dihydroacenaphthylene	9.02	<b>Additional Unidentified Components</b>	<b>1,566</b>
		<b>Total :</b>	<b>18,279</b>

Total Filtered Solids (Dry Weight): 0.2 g (0.013% wt.)

Notes: Detection Limit = 1.84 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Toluene and Ethylbenzene were found in large quantities but not calculated, based on the assumption that the source of these components were from the Xylene solvent.

Due to the lack of a solvent blank to identify additional solvent impurities, all additional components found have been reported.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.

Results are based on a total volume of 1840 mLs.



## Analytical Report

Batch #: 031328.doc

Date: July 18, 2003

**Prepared for:**

**Andrew Kramer  
X5612**

**Gas Technology Institute  
1700 S. Mt Prospect Rd.  
Des Plaines, IL 60018**

**Project # 61129-08**

**Received Date: 7/2/2003**

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In April 2000, the Institute of Gas Technology (IGT) and the Gas Research Institute (GRI) combined to form the Gas Technology Institute (GTI).

Submitted by: Karen Crippen, (847) 768-0604  
Chemical Research Services

Technical Contact:

A handwritten signature in black ink that reads "Alan G. Janos". The signature is written in a cursive style and is positioned above a horizontal line.

Alan G. Janos, (847) 768-0603

## Analytical Report

Batch #: 031328.doc

Date: July 18, 2003

### Liquid Sample #1, 6/30/03 (031328-001)

#### Phase separation:

Organic phase (oil) recovered:	7.6 g	1.26 wt%
Aqueous phase recovered:	581.6 g	96.78 wt%
Solid phase recovered:	0.5 g	0.08 wt%
Loss (by difference):	<u>11.3 g</u>	1.88 wt%
Total sample received:	601.0 g	

Analyst: AGJ

#### Oil phase analyses (031328-003): (GC work only per A.K.)

#### Water phase analyses (031328-002):

Total Carbon:	0.61 wt%
Kjeldahl Nitrogen:	0.23 wt%
Ammonia:	0.25 wt%

Analyst: NJP, JS, AGJ



### Hydrocarbon Analysis in Oil Phase

GTI Sample Number: 031328-003

Sample Description: Liquid 6/30/03 Oil Phase

Tentative Component ID	Estimated mg recovered	Tentative Component ID	Estimated mg recovered
Hydroxymethylpentanone	<0.670	Methoxypropenylphenol/Dimethylnaphthalenes	74.4
Ethylbenzene	<0.670	Ethenylnaphthalene	101
m,p-Xylene	<0.670	Acenaphthylene	571
o-Xylene	<0.670	Methyl-1,1'-biphenyl	36.2
Phenol	18.9	Naphthalenecarbonitriles	26.2
Benzeneamine	<0.670	Acenaphthene	77.6
Benzonitrile	<0.670	Dibenzofuran	96.8
Methylphenols/Indene	46.1	Phenalene	45.5
Methylbenzonitriles	1.29	Fluorene	242
Tetramethylpiperidinone	<0.670	Methyldibenzofuran	18.8
Methylbenzofuran	2.63	Methylfluorenes	43.3
Dimethylphenols	14.6	Fluorenone	28.3
Methylindenes	14.5	Phenanthrene	713
Methylbenzaldehydes	19.9	Anthracene	154
Naphthalene	403	Carbazole	5.58
Quinoline	12.9	Benzoquinoline	6.65
Isoquinoline	11.4	Phenylnaphthalenes	43.4
Indole	13.4	Benzo[c]cinnoline	5.81
2-Methylnaphthalene	221	Methylphenanthrene or Methylanthracene	14.6
1-Methylnaphthalene	161	Methylphenanthrene or Methylanthracene	14.7
Methylquinolines	5.61	4H-Cyclopenta[d,e,f]phenanthrene	46.9
Methylindoles	1.64	Fluoranthene	66.5
1,1'-Biphenyl	183	Pyrene	70.4
Ethyl-naphthalenes	35.7	Methylpyrenes	24.3
Dihydroacenaphthylene	14.4		
Indazole	<0.670	<b>Additional Unidentified Components</b>	<b>477</b>
		<b>Total :</b>	<b>4185</b>

Notes: Detection Limit = 0.670 mg

All components were tentatively identified by comparing the mass spectra of each component to a mass spectral reference library.

Estimated concentrations of all components were calculated as naphthalene by GC-FID.



Sample Login No: 031328-004

Date: July 15, 2003

Sample Description: Carbon ash fines, 6/30/03

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received) w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	4.64	4.64	
Volatile Matter, %	26.47	26.47	27.75
Ash, %	37.46	37.32	
Fixed Carbon, % (by difference)	31.44	31.58	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash, %	39.13
Carbon, %	49.67
Hydrogen, %	0.61
Nitrogen, %	0.24
Sulfur, %	0.12
Oxygen, % (by difference)	10.24

Analyst: NJP



## Analytical Report

Batch # : 031303

July 15, 2003

**Prepared for:**

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 6/26/2003**

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Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report:

A handwritten signature in black ink, appearing to read "Alan G. Janos", is written over a horizontal line.

Alan G. Janos

(847) 768-0603



Sample Login No: 031303-001

Date: July 15, 2003

Sample Description: 6/12/03 Carbon Ash Fines

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u> <u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	6.98	6.98	
Volatile Matter, %	16.08	16.08	17.29
Ash, %	41.58	41.46	
Fixed Carbon, % (by difference)	35.36	35.48	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash, %	44.57
Carbon, %	47.46
Hydrogen, %	0.51
Nitrogen, %	1.05
Sulfur, %	0.05
Oxygen, % (by difference)	6.36

Analyst: NJP





## Analytical Report

Batch # : 031305

July 15, 2003

**Prepared for:**

**Andrew Kramer, Process Engineering**

**Phone: (847) 768-0612**

**Project #: 61129-08**

**Received Date: 6/26/2003**

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Submitted by: Karen Crippen, (847) 768-0604

Chemical Research Services

Technical contact for this report: \_\_\_\_\_

A handwritten signature in cursive script that reads "Alan G. Janos". The signature is written in black ink and is positioned above a horizontal line.

Alan G. Janos

(847) 768-0603



Sample Login No: 031305-001

Date: July 15, 2003

Sample Description: Test #14, Solids, 6/18/03

<b>Proximate Analysis</b>	<u>(As received)</u>	<u>(As received)</u> <u>w/SO3 correction</u>	<u>(Dry basis)</u>
Moisture, %	5.27	5.27	
Volatile Matter, %	14.25	14.25	15.04
Ash, %	25.64	25.50	
Fixed Carbon, % (by difference)	54.85	54.98	

<b>Ultimate Analysis</b>	<u>(Dry basis)</u>
Ash, %	26.92
Carbon, %	65.44
Hydrogen, %	1.16
Nitrogen, %	0.60
Sulfur, %	0.10
Oxygen, % (by difference)	5.78

Analyst: NJP

**Appendix D**  
**Typical Material & Energy Balance**

**Material & Energy Balance - Test Run 15, April 15, 2003**

Input	Lbs/hr	C	H	O	N	S	Ash	Total	Btu	Adj Value	Original values	
Biomass (Menard)		1.37	0.18	1.30	0.00	0.000	0.01	2.87	23,132	3		
Moisture			0.00	0.00				0.00				
Steam			0.10	0.84				0.94	1,371		7.11 g/min	
Air				1.41	4.65			6.06	1,907	35	35.5 l/min	
Nitrogen (purge)					0.66			0.66		2.9	4 l/min	
Total Input		1.37	0.28	3.56	5.31	0.00	0.01	10.54	26,411			
Output		C	H	O	N	S	Ash	Total	Btu			
Solids		0.02					0.01	0.03	379			
Oils/Tars		0.01						0.015	225			
Condensate			0.14	1.10				1.23	2,216	1019	840	
Product Gas Composition								0				
Lbs/hr	0.34261										2.057 cfm	
H2	11.95%	10.200%	13.700%		0.08			0.08				
CO	10.11%	8.620%	11.600%	0.42		0.55		0.97				
CO2	17.00%	16.200%	17.800%	0.70		1.86		2.56				
CH4	3.59%	2.610%	4.570%	0.15	0.05			0.20				
C2H4	0.46%	0.420%	0.509%	0.04	0.01			0.04				
C2H6	0.11%	0.076%	0.148%	0.01	0.00			0.01				
C3H8	0.02%	0.015%	0.020%	0.00	0.00			0.00				
C6H6	0.13%	0.135%	0.120%	0.03	0.00			0.03				
N2	55.45%	60.500%	50.400%			5.32		5.32	6,840			
O2	0.45%	0.34%	0.56%		0.05			0.05				
H2S	0.0154%	0.0159%	0.0149%		0.0001		0.0017	0.0018				
COS	0.0009%	0.0008%	0.0009%	0.0000		0.0000	0.0001	0.0002				
Total Gas	Average	Sample 1	Sample 2	1.34	0.14	2.47	5.32	0.00	0.00	9.28	20,358	
Total Output		1.38	0.28	3.57	5.32	0.00	0.01	10.56	30,019			
Balance (Output/Input), %		-0.16%	0.27%	-0.28%	-0.09%	0.00%	0.00%	-0.18%				

**Material & Energy Balance - Test Run 27, June 26, 2003**

Input	Lbs/hr	C	H	O	N	S	Ash		Total	Btu	Adj Value	Original values
Biomass (Menard)		1.65	0.21	1.56	0.00	0.0003	0.0151		3.44	27,726	3.3	
Moisture			0.00	0.00					0.00			
Steam			0.09	0.73					0.82	1,202		6.23 g/min
Air				1.55	5.11				6.65	2,096	35	39 l/min
Nitrogen (purge)					0.35				0.35		2.3	2.1 l/min
Total Input		1.65	0.30	3.85	5.46	0.0003	0.02		11.27	31,023		
Output		C	H	O	N	S	Ash	NH3	Total	Btu		
Solids		0.04		0.00			0.0151		0.06	663		
Oils/Tars		0.05	0.00	0.00	0.00				0.055	824	0	
Condensate		0.01	0.16	1.29				0.0041	1.46	2,612	577	770
Product Gas Composition										0		
	Lbs/hr	0.345507										2.075 cfm
H2	7.000%		0.05						0.05			
CO	12.500%	0.52		0.69					1.21			
CO2	16.200%	0.67		1.79					2.46			
CH4	4.520%	0.19	0.06						0.25			
C2H4	1.300%	0.11	0.02						0.13			
C2H6	0.160%	0.01	0.00						0.02			
C3H8	0.033%	0.00	0.00						0.01			
C6H6 plus	0.220%	0.05	0.00						0.06			
N2	56.500%				5.47				5.47	7164		
O2	0.65%			0.07					0.07			
H2S	0.0023%		0.0000			0.0002			0.0003			
COS									0.0000			
Total Gas	Sample 1	1.56	0.14	2.55	5.47	0.0002	0.00		9.71	20,531		
Total Output		1.65	0.30	3.85	5.47	0.0002	0.02		11.28	31,793		
Balance (Output/Input), %		-0.07%	-0.07%	-0.04%	-0.14%		0.00%		-0.05%			

**Material & Energy Balance - Test Run 29, June 30, 2003**

Input	Lbs/hr	C	H	O	N	S	Ash		Total	Btu	Adj Value	Original values
Biomass		1.46	0.19	1.39	0.0031	0.000	0.01		3.05	24,583	3.2	
Moisture			0.00	0.00					0.00			
Steam			0.09	0.72					0.81	1,179		6.1143 g/min
Air				1.29	4.26				5.54	1,746	28	32.5 l/min
Nitrogen (purge)					0.12				0.12		2.3	0.75 l/min
Total Input		1.46	0.28	3.40	4.38	0.00	0.01		9.54	27,508		
Output		C	H	O	N	S	Ash	NH3	Total	Btu		
Solids		0.017		0.004		0.0000	0.01		0.034	408		
Oils/Tars		0.055	0.005	0.00					0.062	934	0.01	
Condensate		0.007	0.13	1.01	0.076			0.0029	1.14	2,048	582	690 gms
Product Gas Composition										0		
Lbs/hr	0.301099											1.8083 cfm
H2	11.15%	11.40%	10.90%		0.07				0.07			
CO	10.60%	10.80%	10.40%	0.383		0.51			0.89			
CO2	18.85%	18.90%	18.80%	0.681		1.82			2.50			
CH4	4.54%	4.64%	4.44%	0.164	0.05				0.22			
C2H4	1.120%	1.140%	1.10%	0.081	0.01				0.09			
C2H6	0.230%	0.240%	0.22%	0.017	0.00				0.02			
C3H8	0.120%	0.123%	0.12%	0.013	0.00				0.02			
C6H6 plus	0.20%	0.20%	0.20%	0.044	0.00				0.05			
N2	51.90%	51.20%	52.60%			4.38			4.38	6,106		
O2	0.51%	0.57%	0.46%			0.05			0.05			
H2S	0.0024%	0.0023%	0.0024%		0.0000		0.0002		0.0002			
COS												
Total Gas	Average	Sample 1	Sample 2	1.382	0.15	2.38	4.38	0.0002	0.00	8.28	17,892	
Total Output				1.462	0.28	3.40	4.45	0.0003	0.01	9.52	27,389	
Balance (Output/Input), %				-0.002	-0.20%	0.15%	-1.49%	13.17%	0.00%	0.19%		