



303-637-0150

EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PROJECT NO. :	201501070	ANALYSIS NO. :	01
COMPANY NAME :	EE3 LLC	ANALYSIS DATE:	JANUARY 21, 2015
ACCOUNT NO. :		SAMPLE DATE :	
PRODUCER :		CYLINDER NO. :	0592
LEASE NO. :		SAMPLED BY :	ZACHERY TERRILE
NAME/DESCRIP :	MUTUAL 4-30H		

FIELD DATA		SAMPLE TEMP. :	88
SAMPLE PRES. :	130	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	PROBE; POSSIBLE MOISTURE IN SAMPLE - EMPACT		

<u>COMPONENT</u>	<u>MOLE %</u>	<u>MASS %</u>	<u>GPM @ 14.650</u>	<u>GPM @ 14.730</u>
ALCOHOLS	3.4717	4.3590		
GLYCOLS	0.0013	0.0032		
HELIUM	0.03	0.00	---	---
HYDROGEN	0.14	0.01	---	---
OXYGEN/ARGON	0.01	0.01	---	---
NITROGEN	1.46	1.60	---	---
CARBON DIOXIDE	1.57	2.71	---	---
METHANE	64.07550	40.25160	---	---
ETHANE	11.8581	13.9624	3.1698	3.1871
PROPANE	9.7008	16.7506	2.6718	2.6864
I-BUTANE	1.0875	2.4751	0.3558	0.3577
N-BUTANE	3.4712	7.9004	1.0944	1.1003
I-PENTANE	0.8648	2.4358	0.3097	0.3113
N-PENTANE	0.9125	2.5781	0.3307	0.3325
HEXANES PLUS	1.3466	4.9538	0.5509	0.5538
<u>TOTALS</u>	<u>100.00000</u>	<u>100.00000</u>	<u>8.4831</u>	<u>8.5291</u>

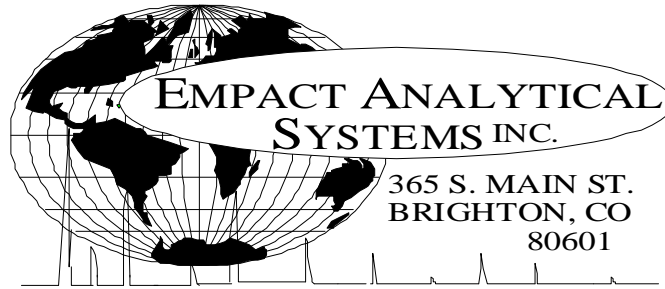
<u>BTEX COMPONENTS</u>	<u>MOLE%</u>	<u>WT%</u>		<u>BTU @ 14.650</u>	<u>14.730</u>
BENZENE	0.0289	0.0884	LOW NET DRY REAL :	1294.1 /scf	1301.2 /scf
TOLUENE	0.0214	0.0772	NET WET REAL :	1271.5 /scf	1278.5 /scf
ETHYLBENZENE	0.0023	0.0096	HIGH GROSS DRY REAL :	1421.7 /scf	1429.4 /scf
XYLENES	0.0094	0.0392	GROSS WET REAL :	1396.8 /scf	1404.6 /scf
<u>TOTAL BTEX</u>	<u>0.0620</u>	<u>0.2144</u>	NET DRY REAL :	19265.9 /lb	19371.2 /lb
			GROSS DRY REAL :	21159.3 /lb	21274.9 /lb

RELATIVE DENSITY (AIR=1):	0.8802
COMPRESSIBILITY FACTOR :	0.99472

(CALC: GPA STD 2145 & TP-17 @14.696 & 60 F)

*(DETAILED HYDROCARBON ANALYSIS/NJ 1993) : ASTM D6730

THIS DATA HAS BEEN ACQUIRED THROUGH APPLICATION OF CURRENT STATE-OF-THE-ART ANALYTICAL TECHNIQUES.
THE USE OF THIS INFORMATION IS THE RESPONSIBILITY OF THE USER. EMPACT ANALYTICAL SYSTEMS, ASSUMES NO
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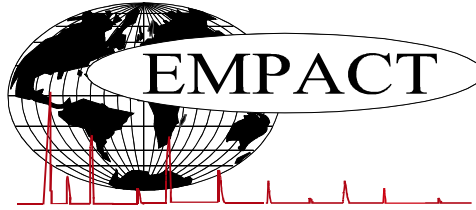
GLYCALC INFORMATION

PROJECT NO. :	201501070	ANALYSIS NO. :	01
COMPANY NAME :	EE3 LLC	ANALYSIS DATE:	JANUARY 21, 2015
ACCOUNT NO. :		SAMPLE DATE :	
PRODUCER :		CYLINDER NO. :	0592
LEASE NO. :		SAMPLED BY :	ZACHERY TERRILE
NAME/DESCRIP :	MUTUAL 4-30H		

FIELD DATA		SAMPLE TEMP. :	88
SAMPLE PRES. :	130	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	PROBE; POSSIBLE MOISTURE IN SAMPLE - EMPACT		

<u>Componet</u>	<u>Mole %</u>	<u>Wt %</u>
Helium	0.03	0.00
Hydrogen	0.14	0.01
Carbon Dioxide	1.57	2.71
Nitrogen	1.46	1.60
Methane	64.07550	40.25160
Ethane	11.8581	13.9624
Propane	9.7008	16.7506
Isobutane	1.0875	2.4751
n-Butane	3.4712	7.9004
Isopentane	0.7702	2.1760
n-Pentane	0.9125	2.5781
Cyclopentane	0.0946	0.2598
n-Hexane	0.2415	0.8149
Cyclohexane	0.0747	0.2462
Other Hexanes	0.4610	1.5426
Heptanes	0.2642	1.0293
Methycyclohexane	0.0734	0.2822
2,2,4 Trimethylpentane	0.0002	0.0009
Benzene	0.0289	0.0884
Toluene	0.0214	0.0772
Ethylbenzene	0.0023	0.0096
Xylenes	0.0094	0.0392
C8+ Heavies	0.1696	0.8233
<u>Subtotal</u>	<u>96.51700</u>	<u>95.62780</u>
Oxygen/Argon	0.01	0.01
Alcohols	3.4717	4.3590
Glycols	0.0013	0.0032
Total	100.00000	100.00000

THE DATA PRESENTED HEREIN HAS BEEN ACQUIRED THROUGH JUDICIOUS APPLICATION OF CURRENT STATE-OF-THE ART ANALYTICAL TECHNIQUES. THE APPLICATIONS OF THIS INFORMATION IS THE RESPONSIBILITY OF THE USER. EMPACT ANALYTICAL SYSTEMS, INC. ASSUMES NO RESPONSIBILITY FOR ACCURACY OF THE REPORTED INFORMATION NOR ANY CONSEQUENCES OF IT'S APPLICATION.



**EXTENDED NATURAL GAS ANALYSIS (*DHA)
DHA COMPONENT LIST**

PROJECT NO. : 201501070
 COMPANY NAME : EE3 LLC
 ACCOUNT NO. :
 PRODUCER :
 LEASE NO. :
 NAME/DESCRIP : MUTUAL 4-30H

ANALYSIS NO. : 01
 ANALYSIS DATE: JANUARY 21, 2015
 SAMPLE DATE :
 CYLINDER NO. : 0592
 SAMPLED BY : ZACHERY TERRILE

FIELD DATA

SAMPLE PRES. : 130
 VAPOR PRES. :
 COMMENTS : PROBE; POSSIBLE MOISTURE IN SAMPLE - EMPACT

SAMPLE TEMP. : 88
 AMBIENT TEMP.:
 GRAVITY :

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
Helium	---	0.03	0.00	---	---
Hydrogen	---	0.14	0.01	---	---
Oxygen/Argon	---	0.01	0.01	---	---
Nitrogen	---	1.46	1.60	---	---
Carbon Dioxide	---	1.57	2.71	---	---
Methane	P1	64.07550	40.25160	---	---
Ethane	P2	11.8581	13.9624	3.170	3.187
Propane	P3	9.7008	16.7506	2.672	2.686
i-Butane	I4	1.0875	2.4751	0.356	0.358
Methanol	X1	3.4665	4.3495	0.442	0.444
n-Butane	P4	3.4712	7.9004	1.094	1.100
2,2-Dimethylpropane	I5	0.0028	0.0079	0.001	0.001
Ethanol	X2	0.0050	0.0090	0.001	0.001
i-Pentane	I5	0.7674	2.1681	0.281	0.282
i-Propanol	X3	0.0002	0.0005	0.000	0.000
n-Pentane	P5	0.9124	2.5778	0.331	0.333
2,2-Dimethylbutane	I6	0.0024	0.0081	0.001	0.001
Cyclopentane	N5	0.0946	0.2598	0.028	0.028
2,3-Dimethylbutane	I6	0.0168	0.0567	0.007	0.007
2-Methylpentane	I6	0.1752	0.5912	0.073	0.074
3-Methylpentane	I6	0.0995	0.3358	0.041	0.041
UnknownC5s	U5	0.0001	0.0003	0.000	0.000
n-Hexane	P6	0.2415	0.8149	0.099	0.100
2,2-Dimethylpentane	I7	0.0003	0.0012	0.000	0.000
Methylcyclopentane	N6	0.1670	0.5504	0.059	0.060
2,4-Dimethylpentane	I7	0.0074	0.0291	0.003	0.003
2,2,3-Trimethylbutane	I7	0.0002	0.0008	0.000	0.000
Benzene	A6	0.0289	0.0884	0.008	0.008
3,3-Dimethylpentane	I7	0.0006	0.0024	0.000	0.000
Cyclohexane	N6	0.0747	0.2462	0.025	0.025
2-Methylhexane	I7	0.0319	0.1252	0.015	0.015
2,3-Dimethylpentane	I7	0.0143	0.0561	0.006	0.006
1,1-Dimethylcyclopentane	N7	0.0106	0.0408	0.004	0.004
3-Methylhexane	I7	0.0371	0.1456	0.017	0.017
1c,3-Dimethylcyclopentane	N7	0.0222	0.0854	0.010	0.010
Ethylene glycol	GL2	0.0013	0.0032	0.001	0.001
1t,3-Dimethylcyclopentane	N7	0.0198	0.0761	0.009	0.009
3-Ethylpentane	I7	0.0033	0.0130	0.001	0.001
1t,2-Dimethylcyclopentane	N7	0.0362	0.1392	0.017	0.017
2,2,4-Trimethylpentane	I8	0.0002	0.0009	0.000	0.000
UnknownC6s	U6	0.0001	0.0004	0.000	0.000

n-Heptane	P7	0.0711	0.2790	0.033	0.033
1c,2-Dimethylcyclopentane	N7	0.0032	0.0123	0.001	0.001
Methylcyclohexane	N7	0.0734	0.2822	0.029	0.029
2,2-Dimethylhexane	I8	0.0080	0.0358	0.004	0.004
Ethylcyclopentane	N7	0.0059	0.0227	0.002	0.002
2,5-Dimethylhexane	I8	0.0018	0.0081	0.001	0.001
2,2,3-Trimethylpentane	I8	0.0003	0.0013	0.000	0.000
2,4-Dimethylhexane	I8	0.0027	0.0121	0.001	0.001
1c,2t,4-Trimethylcyclopentane	N8	0.0057	0.0251	0.003	0.003
3,3-Dimethylhexane	I8	0.0004	0.0018	0.000	0.000
1t,2c,4-Trimethylcyclopentane	N8	0.0064	0.0281	0.003	0.003
2,3,4-Trimethylpentane	I8	0.0010	0.0045	0.000	0.000
Toluene	A7	0.0214	0.0772	0.007	0.007
2,3-Dimethylhexane	I8	0.0022	0.0098	0.001	0.001
2-Methyl-3-ethylpentane	I8	0.0011	0.0049	0.001	0.001
1,1,2-Trimethylcyclopentane	N8	0.0002	0.0009	0.000	0.000
2-Methylheptane	I8	0.0125	0.0559	0.006	0.006
4-Methylheptane	I8	0.0033	0.0148	0.002	0.002
3-Methyl-3-ethylpentane	I8	0.0004	0.0018	0.000	0.000
3,4-Dimethylhexane	I8	0.0004	0.0018	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0003	0.0013	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0003	0.0013	0.000	0.000
3-Methylheptane	I8	0.0057	0.0255	0.003	0.003
1c,2t,3-Trimethylcyclopentane	N8	0.0105	0.0461	0.005	0.005
3-Ethylhexane	I8	0.0013	0.0058	0.001	0.001
1t,4-Dimethylcyclohexane	N8	0.0038	0.0167	0.002	0.002
1,1-Dimethylcyclohexane	N8	0.0011	0.0048	0.000	0.000
3c-Ethylmethylcyclopentane	N8	0.0001	0.0004	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0013	0.0057	0.001	0.001
2t-Ethylmethylcyclopentane	N8	0.0011	0.0048	0.001	0.001
1,1-Methylethylcyclopentane	N8	0.0036	0.0158	0.002	0.002
2,2,4-Trimethylhexane	I9	0.0004	0.0020	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0055	0.0242	0.003	0.003
UnknownC7s	U7	0.0001	0.0004	0.000	0.000
n-Octane	P8	0.0164	0.0733	0.008	0.008
1c,4-Dimethylcyclohexane	N8	0.0020	0.0088	0.001	0.001
i-Propylcyclopentane	I8	0.0002	0.0009	0.000	0.000
2,4,4-Trimethylhexane	I9	0.0002	0.0010	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0004	0.0020	0.000	0.000
2,2,3,4-Tetramethylpentane	I9	0.0001	0.0005	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0002	0.0010	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0016	0.0071	0.001	0.001
2,2-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0048	0.0237	0.002	0.002
2,2,3-Trimethylhexane	I9	0.0016	0.0080	0.001	0.001
2,4-Dimethylheptane	I9	0.0006	0.0030	0.000	0.000
4,4-Dimethylheptane	I9	0.0002	0.0010	0.000	0.000
Ethylcyclohexane	N8	0.0024	0.0105	0.001	0.001
n-Propylcyclopentane	N8	0.0011	0.0048	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0002	0.0010	0.000	0.000
2,5-Dimethylheptane	I9	0.0003	0.0015	0.000	0.000
3,3-Dimethylheptane	I9	0.0004	0.0020	0.000	0.000
3,5-Dimethylheptane	I9	0.0002	0.0010	0.000	0.000
2,6-Dimethylheptane	I9	0.0003	0.0015	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0005	0.0025	0.000	0.000
Ethylbenzene	I8	0.0023	0.0096	0.001	0.001
1c,2t,4t-Trimethylcyclohexane	N9	0.0011	0.0054	0.001	0.001
2,3-Dimethylheptane	I9	0.0010	0.0050	0.001	0.001
1,3-Dimethylbenzene (m-Xylene)	A8	0.0043	0.0179	0.002	0.002
1,4-Dimethylbenzene (p-Xylene)	A8	0.0029	0.0121	0.001	0.001
3,4-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0006	0.0030	0.000	0.000
4-Ethylheptane	I9	0.0002	0.0010	0.000	0.000
4-Methyloctane	I9	0.0010	0.0050	0.001	0.001

2-Methyloctane	I9	0.0011	0.0055	0.001	0.001
1c,2t,3-Trimethylcyclohexane	N9	0.0002	0.0010	0.000	0.000
3-Ethylheptane	I9	0.0005	0.0025	0.000	0.000
3-Methyloctane	I9	0.0013	0.0065	0.001	0.001
1c,2t,4c-Trimethylcyclohexane	I9	0.0001	0.0005	0.000	0.000
1,1,2-Trimethylcyclohexane	N9	0.0002	0.0010	0.000	0.000
3,3-Diethylpentane	I9	0.0002	0.0010	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0022	0.0092	0.001	0.001
i-Butylcyclopentane	N9	0.0011	0.0054	0.001	0.001
UnknownC8s	U8	0.0007	0.0031	0.000	0.000
n-Nonane	P9	0.0039	0.0196	0.002	0.002
1,1-Methylethylcyclohexane	N9	0.0007	0.0035	0.000	0.000
i-Propylbenzene	A9	0.0008	0.0038	0.000	0.000
i-Propylcyclohexane	N9	0.0002	0.0010	0.000	0.000
2,2-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
2,4-Dimethyloctane	I10	0.0002	0.0011	0.000	0.000
2,6-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
2,5-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
n-Butylcyclopentane	N9	0.0008	0.0040	0.000	0.000
3,3-Dimethyloctane	I10	0.0002	0.0011	0.000	0.000
n-Propylbenzene	A9	0.0007	0.0033	0.000	0.000
3,6-Dimethyloctane	I10	0.0005	0.0028	0.000	0.000
3-Methyl-5-ethylheptane	I10	0.0006	0.0033	0.000	0.000
1,3-Methylethylbenzene	A9	0.0009	0.0042	0.001	0.001
1,4-Methylethylbenzene	A9	0.0004	0.0019	0.000	0.000
1,3,5-Trimethylbenzene	A9	0.0006	0.0028	0.000	0.000
2,3-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
5-Methylnonane	I10	0.0005	0.0028	0.000	0.000
1,2-Methylethylbenzene	A9	0.0008	0.0038	0.000	0.000
2-Methylnonane	I10	0.0001	0.0006	0.000	0.000
3-Ethylheptane	I10	0.0001	0.0006	0.000	0.000
3-Methylnonane	I10	0.0003	0.0017	0.000	0.000
t-Butylbenzene	A10	0.0014	0.0074	0.001	0.001
i-Butylcyclohexane	N10	0.0003	0.0016	0.000	0.000
1t-Methyl-2-n-propylcyclohexane	I10	0.0001	0.0006	0.000	0.000
i-Butylbenzene	A10	0.0001	0.0005	0.000	0.000
sec-Butylbenzene	A10	0.0001	0.0005	0.000	0.000
UnknownC9s	U9	0.0033	0.0166	0.002	0.002
n-Decane	P10	0.0017	0.0095	0.001	0.001
1,2,3-Trimethylbenzene	A9	0.0006	0.0028	0.000	0.000
1,3-Methyl-i-propylbenzene	A10	0.0002	0.0011	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0002	0.0011	0.000	0.000
Sec-Butylcyclohexane	A10	0.0004	0.0022	0.000	0.000
1,2-Methyl-i-propylbenzene	A10	0.0003	0.0016	0.000	0.000
3-Ethylnonane	I10	0.0001	0.0006	0.000	0.000
1,3-Diethylbenzene	A10	0.0001	0.0005	0.000	0.000
1,3-Methyl-n-propylbenzene	A10	0.0001	0.0005	0.000	0.000
1,4-Methyl-n-propylbenzene	A10	0.0004	0.0021	0.000	0.000
n-Butylbenzene	A10	0.0002	0.0011	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0003	0.0016	0.000	0.000
1,2-Diethylbenzene	A10	0.0002	0.0011	0.000	0.000
t-Decahydronaphthalene	A9	0.0002	0.0012	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0002	0.0011	0.000	0.000
1,4-Dimethyl-2-ethylbenzene	A10	0.0002	0.0011	0.000	0.000
1,3-Dimethyl-4-ethylbenzene	A10	0.0003	0.0016	0.000	0.000
1,3-Dimethyl-2-ethylbenzene	A10	0.0002	0.0011	0.000	0.000
1,2-Dimethyl-3-ethylbenzene	A10	0.0003	0.0016	0.000	0.000
1,2-Ethyl-i-propylbenzene	A10	0.0002	0.0012	0.000	0.000
1,4-Methyl-t-butylbenzene	A11	0.0002	0.0012	0.000	0.000
UnknownC10s	U10	0.0034	0.0190	0.002	0.002
n-Undecane	P11	0.0011	0.0067	0.001	0.001
1,4-Ethyl-i-propylbenzene	A11	0.0001	0.0006	0.000	0.000
1,2,4,5-Tetramethylbenzene	A11	0.0004	0.0021	0.000	0.000
1,2-Methyl-n-butylbenzene	A11	0.0001	0.0006	0.000	0.000

1,2,3,5-Tetramethylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2-Methyl-t-butylbenzene	A11	0.0001	0.0006	0.000	0.000
1,2-Ethyl-n-propylbenzene	A11	0.0002	0.0012	0.000	0.000
2-Methylindan	A11	0.0002	0.0010	0.000	0.000
1,3-Di-i-propylbenzene	A11	0.0001	0.0006	0.000	0.000
sec-Pentylbenzene	A11	0.0001	0.0006	0.000	0.000
n-Pentylbenzene	A11	0.0001	0.0006	0.000	0.000
1t-M-2-(4MP)cyclopentane	P12	0.0001	0.0007	0.000	0.000
1,2-Di-n-propylbenzene	A11	0.0002	0.0013	0.000	0.000
1,4-Di-i-propylbenzene	A11	0.0002	0.0013	0.000	0.000
Tetrahydronaphthalene	A10	0.0001	0.0005	0.000	0.000
Naphthalene	A10	0.0002	0.0010	0.000	0.000
1-t-Butyl-3,5-dimethylbenzene	A12	0.0001	0.0006	0.000	0.000
1,4-Ethyl-t-butylbenzene	A11	0.0001	0.0006	0.000	0.000
1,3-Di-n-propylbenzene	A12	0.0007	0.0045	0.001	0.001
UnknownC11s	U11	0.0026	0.0159	0.002	0.002
n-Dodecane	P12	0.0008	0.0053	0.001	0.001
1,3,5-Triethylbenzene	A12	0.0002	0.0013	0.000	0.000
1,2,4-Triethylbenzene	A12	0.0002	0.0013	0.000	0.000
1,4-Methyl-n-pentylbenzene	A12	0.0001	0.0006	0.000	0.000
n-Hexylbenzene	A12	0.0004	0.0026	0.000	0.000
1,2,3,4,5-Pentamethylbenzene	A13	0.0001	0.0006	0.000	0.000
2-Methylnaphthalene	A11	0.0003	0.0017	0.000	0.000
1-Methylnaphthalene	A11	0.0002	0.0011	0.000	0.000
UnknownC12s	U12	0.0020	0.0123	0.001	0.001
n-Tridecane	P13	0.0005	0.0036	0.000	0.000
UnknownC13s	U13	0.0016	0.0116	0.001	0.001
n-Tetradecane	P14	0.0004	0.0031	0.000	0.000
UnknownC14s	U14	0.0016	0.0124	0.001	0.001
n-Pentadecane	P15	0.0003	0.0025	0.000	0.000
UnknownC15s	U15	0.0010	0.0083	0.001	0.001
n-Hexadecane	P16	0.0001	0.0009	0.000	0.000
UnknownC16s	U16	0.0001	0.0009	0.000	0.000
UnknownC19s	U19	0.0001	0.0011	0.000	0.000
UnknownC20s	U20	0.0001	0.0011	0.000	0.000
TOTAL		100.00000	100.00000	8.9271	8.9755

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.650	14.730
BENZENE	0.0289	0.0884	LOW NET DRY REAL :	1294.1 /scf	1301.2 /scf
TOLUENE	0.0214	0.0772	NET WET REAL :	1271.5 /scf	1278.5 /scf
ETHYLBENZENE	0.0023	0.0096	HIGH GROSS DRY REAL :	1421.7 /scf	1429.4 /scf
XYLENES	0.0094	0.0392	GROSS WET REAL :	1396.8 /scf	1404.6 /scf
TOTAL BTEX	0.0620	0.2144	NET DRY REAL :	19265.9 /lb	19371.2 /lb
			GROSS DRY REAL :	21159.3 /lb	21274.9 /lb

RELATIVE DENSITY (AIR=1): 0.8802
 COMPRESSIBILITY FACTOR : 0.99472

(CALC: GPA STD 2145 & TP-17 @14.696 & 60 F)

*(DETAILED HYDROCARBON ANALYSIS/NJ 1993) ; ASTM D6730

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