



EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PRIMARY DB KEY: **05-103-10782** NAME/DESCRIP : **FED 4S-95-1-21-DP**
 LEASE #: **INTERMEDIATE CSG**
 FIELD/AREA:

PROJECT NO. : **202506059** ANALYSIS NO. : **03**
 COMPANY NAME : **QB ENERGY OPERATING, LLC** ANALYSIS DATE: **JUNE 20, 2025 13:02**
 OFFICE / BRANCH: **PARACHUTE, CO** SAMPLE DATE : **MAY 19, 2025**
 CUSTOMER REF: TO:
 PRODUCER : **QB ENERGY OPERATING, LLC** EFFECTIVE DATE:

*****FIELD DATA*****

SAMPLE CYCLE: SAMPLE TYPE:
 SAMPLE PRES. : 1830 psig PROBE :
 FLOW PRES. : psig CYLINDER NO. : ECA-802
 LAB PRES: psig SAMPLED BY : NICK CROY
 SAMPLE TEMP. : 43 °f SAMPLING COMPANY: **QB ENERGY**
 AMBIENT TEMP.: °f H2S BY STAIN TUBE: — ppm mol
 H2O BY STAIN TUBE: - #/mmcf CO2 BY STAIN TUBE: - Mol %
 FIELD COMMENTS:
 LAB COMMENTS:

<u>COMPONENT</u>	<u>MOLE %</u>	<u>MASS %</u>	<u>GPM @</u>	
			<u>14.65</u>	<u>14.73</u>
ALCOHOLS	0.0008	0.0016	0.0000	0.0000
HELIUM	0.00	0.00	---	---
HYDROGEN	0.27	0.03	---	---
OXYGEN/ARGON	0.00	0.00	---	---
NITROGEN	0.08	0.13	---	---
CARBON DIOXIDE	0.02	0.05	---	---
METHANE	97.2455	92.7799	---	---
ETHANE	1.6583	2.9654	0.4415	0.4439
PROPANE	0.1878	0.4925	0.0519	0.0522
I-BUTANE	0.0576	0.1991	0.0190	0.0191
N-BUTANE	0.0288	0.0996	0.0090	0.0090
I-PENTANE	0.0244	0.1046	0.0090	0.0090
N-PENTANE	0.0101	0.0433	0.0040	0.0040
HEXANES PLUS	0.4167	3.1040	0.2050	0.2052
TOTALS	100.00000	100.00000	0.7394	0.7424

<u>BTEX COMPONENTS</u>	<u>MOLE%</u>	<u>WT%</u>
BENZENE	0.0169	0.0785
TOLUENE	0.0328	0.1797
ETHYLBENZENE	0.0026	0.0164
XYLENES	0.0269	0.1698
TOTAL BTEX	0.0792	0.4444

	<u>CALCULATED VALUES**</u>	
	<u>BTU @</u>	<u>BTU @</u>
	<u>14.65</u>	<u>14.73</u>
LHV NET DRY REAL :	944.2 /scf	949.4 /scf
NET WET REAL :	927.7 /scf	932.9 /scf
HHV GROSS DRY REAL :	1046.4 /scf	1052.1 /scf
GROSS WET REAL :	1028.1 /scf	1033.8 /scf
NET HEATING VALUE (60 °F ideal reaction):		21347.1 Btu/lbm
GROSS HEATING VALUE (60°F ideal reaction):		23674.6 Btu/lbm
RELATIVE DENSITY (AIR=1):		0.5793
DENSITY		0.04431 lbm/scf
COMPRESSIBILITY FACTOR :		0.9979
REGULAR WOBBE INDEX		1376.3

**(DETAILED HYDROCARBON ANALYSIS/NJ 1993)
 Mod ASTM D6730,GPA 2261 & GPA 2286.*

*** (CALC: GPA 2172, GPA 2145 & TP-17 @14.696 & 60 F)*

The data presented herein has been acquired by means of current analytical techniques and represents the judicious conclusion EMPACT Analytical Systems, Inc. Results of the analysis can be affected by the sampling conditions, therefore, are only warranted through proper lab protocol. EMPACT assumes no responsibility for interpretation or any consequences from application of the reported information and is the sole liability of the user. The reproduction in any media of this reported information may not be made, in portion or as a whole, without the written permission of EMPACT Analytical Systems, Inc.



**EXTENDED NATURAL GAS ANALYSIS (*DHA)
GLYCALC INFORMATION**

PROJECT NO. :	202506059	ANALYSIS NO. :	03
COMPANY NAME :	QB ENERGY OPERATING, LLC	ANALYSIS DATE:	JUNE 20, 2025 13:02
ACCOUNT NO. :		SAMPLE DATE :	MAY 19, 2025
PRODUCER :	QB ENERGY OPERATING, LLC	CYLINDER NO. :	ECA-802
LEASE NO. :		SAMPLED BY :	NICK CROY
NAME/DESCRIP :	FED 4S-95-1-21-DP INTERMEDIATE CSG		

FIELD DATA

SAMPLE PRES. :	1830	SAMPLE TEMP. :	43
H2S BY STAIN TUBE:	— ppm mol	AMBIENT TEMP.:	
COMMENTS :			

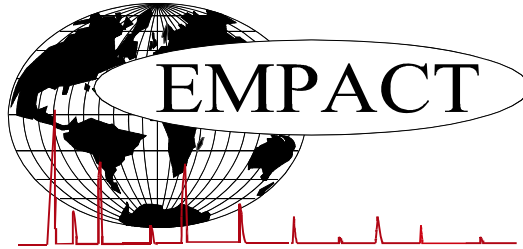
<u>Componet</u>	<u>Mole %</u>	<u>Wt %</u>
Helium	0.00	0.00
Hydrogen	0.27	0.03
Carbon Dioxide	0.02	0.05
Nitrogen	0.08	0.13
Methane	97.2455	92.7799
Ethane	1.6583	2.9654
Propane	0.1878	0.4925
Isobutane	0.0576	0.1991
n-Butane	0.0288	0.0996
Isopentane	0.0236	0.1013
n-Pentane	0.0101	0.0433
Cyclopentane	0.0008	0.0033
n-Hexane	0.0059	0.0302
Cyclohexane	0.0077	0.0385
Other Hexanes	0.0219	0.1116
Heptanes	0.0244	0.1453
Methylcyclohexane	0.0211	0.1232
2,2,4 Trimethylpentane	0.0001	0.0007
Benzene	0.0169	0.0785
Toluene	0.0328	0.1797
Ethylbenzene	0.0026	0.0164
Xylenes	0.0269	0.1698
C8+ Heavies	0.2564	2.2101
<u>Subtotal</u>	<u>99.99920</u>	<u>99.99840</u>
Oxygen/Argon	0.00	0.00
Alcohols	0.0008	0.0016
<u>Total</u>	<u>100.00000</u>	<u>100.00000</u>

Calculated Values BTU @		Total	C6+	C8+	C10+
	LHV Net Dry Real:	944.2	6220.3	7028.9	7874.7 Btu/scf
	Net Wet Real:	927.7	6111.6	6906.0	7737.0 Btu/scf
	HHV Gross Dry Real:	1046.4	6688.3	7576.1	8518.0 Btu/scf
	Gross Wet Real:	1028.1	6571.4	7443.7	8369.1 Btu/scf

Other Calculated Values					
Regualr Wobbe Index*	1376.3	3222.3	3444.9	3681.0	Btu/scf
Net Heating Value (60 °F ideal reaction):	21347.1	18738.1	18765.1	18754.1	Btu/lbm
Gross Heating Value (60 °F ideal reaction):	23674.6	20141.3	20220.3	20276.7	Btu/lbm
Molar Mass (MW):	16.81564	125.159	140.835	156.025	g/mol
Relative Density (AIR=1):	0.5793	4.3211	4.8618	5.3868	SG
Density:	0.04431	0.32982	0.37111	0.41115	lbm/scf
Compressibility Factor:	0.9979	0.9983	0.9995	0.9998	Z
Liquid Volume real gas @:	17.1661	0.2044	0.1565	0.1087	gal/1000 scf

* The Wobbe pressure base in the number considered is based upon the given Pb of the HHV above.
 #DIV/0 or 0 (zero) will appear in the Calculated Value Section when there is no C6+, C8+ or C10+ in the sample to calculate these factors.
 BDL - Below Detection Limit. The H2S LOS has a detection limit of 0.25 ppm. A _ (an underscore) indicates there was no tube pulled for H2S.

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DHA COMPONENT LIST

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 AMBIENT TEMP.: °f H2S BY STAIN TUBE: - ppm mol
 H2O BY STAIN TUBE: - #/mmcf CO2 BY STAIN TUBE: - Mol %
 FIELD COMMENTS:
 LAB COMMENTS:

COMPONENT	PIANO #	MOLE %	MASS %	GPM @	
				14.65	14.73
Hydrogen	---	0.27	0.03	---	---
Nitrogen	---	0.08	0.13	---	---
Carbon Dioxide	---	0.02	0.05	---	---
Methane	P1	97.2455	92.7799	---	---
Ethane	P2	1.6583	2.9654	0.442	0.444
Propane	P3	0.1878	0.4925	0.052	0.052
i-Butane	I4	0.0576	0.1991	0.019	0.019
Methanol	X1	0.0008	0.0016	0.000	0.000
n-Butane	P4	0.0288	0.0996	0.009	0.009
2,2-Dimethylpropane	I5	0.0027	0.0116	0.001	0.001
i-Pentane	I5	0.0209	0.0897	0.008	0.008
n-Pentane	P5	0.0099	0.0425	0.004	0.004
2,2-Dimethylbutane	I6	0.0032	0.0164	0.001	0.001
Cyclopentane	N5	0.0008	0.0033	0.000	0.000
2,3-Dimethylbutane	I6	0.0021	0.0108	0.001	0.001
2-Methylpentane	I6	0.0070	0.0359	0.003	0.003
3-Methylpentane	I6	0.0042	0.0215	0.002	0.002
UnknownC5s	U5	0.0002	0.0008	0.000	0.000
n-Hexane	P6	0.0059	0.0302	0.002	0.002
2,2-Dimethylpentane	I7	0.0007	0.0042	0.000	0.000
Methylcyclopentane	N6	0.0054	0.0270	0.002	0.002
2,4-Dimethylpentane	I7	0.0010	0.0060	0.000	0.000
2,2,3-Trimethylbutane	I7	0.0004	0.0024	0.000	0.000
Benzene	A6	0.0169	0.0785	0.005	0.005
3,3-Dimethylpentane	I7	0.0007	0.0042	0.000	0.000
Cyclohexane	N6	0.0077	0.0385	0.003	0.003
2-Methylhexane	I7	0.0026	0.0155	0.001	0.001
2,3-Dimethylpentane	I7	0.0024	0.0143	0.001	0.001

1,1-Dimethylcyclopentane	N7	0.0009	0.0052	0.000	0.000
3-Methylhexane	I7	0.0038	0.0227	0.002	0.002
1c,3-Dimethylcyclopentane	N7	0.0012	0.0070	0.001	0.001
1t,3-Dimethylcyclopentane	N7	0.0011	0.0064	0.001	0.001
3-Ethylpentane	I7	0.0002	0.0012	0.000	0.000
1t,2-Dimethylcyclopentane	N7	0.0018	0.0105	0.001	0.001
2,2,4-Trimethylpentane	I8	0.0001	0.0007	0.000	0.000
n-Heptane	P7	0.0059	0.0352	0.003	0.003
1c,2-Dimethylcyclopentane	N7	0.0002	0.0012	0.000	0.000
Methylcyclohexane	N7	0.0211	0.1232	0.008	0.008
2,2-Dimethylhexane	I8	0.0002	0.0014	0.000	0.000
1,1,3-Trimethylcyclopentane	N7	0.0006	0.0040	0.000	0.000
Ethylcyclopentane	N7	0.0008	0.0047	0.000	0.000
2,5-Dimethylhexane	I8	0.0008	0.0054	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0008	0.0054	0.000	0.000
2,4-Dimethylhexane	I8	0.0001	0.0007	0.000	0.000
1c,2t,4-Trimethylcyclopentane	N8	0.0003	0.0020	0.000	0.000
3,3-Dimethylhexane	I8	0.0003	0.0020	0.000	0.000
Toluene	A7	0.0328	0.1797	0.011	0.011
2,3-Dimethylhexane	I8	0.0005	0.0034	0.000	0.000
2-Methyl-3-ethylpentane	I8	0.0002	0.0014	0.000	0.000
2-Methylheptane	I8	0.0032	0.0218	0.002	0.002
4-Methylheptane	I8	0.0011	0.0075	0.001	0.001
3-Methyl-3-ethylpentane	I8	0.0002	0.0014	0.000	0.000
3,4-Dimethylhexane	I8	0.0001	0.0007	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0001	0.0007	0.000	0.000
3-Methylheptane	I8	0.0018	0.0123	0.001	0.001
1c,2t,3-Trimethylcyclopentane	N8	0.0048	0.0321	0.002	0.002
3-Ethylhexane	I8	0.0009	0.0061	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0021	0.0140	0.001	0.001
1,1-Dimethylcyclohexane	N8	0.0008	0.0054	0.000	0.000
2,2,5-Trimethylhexane	I9	0.0001	0.0008	0.000	0.000
3c-Ethylmethylcyclopentane	N8	0.0002	0.0013	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0002	0.0013	0.000	0.000
2t-Ethylmethylcyclopentane	N8	0.0002	0.0013	0.000	0.000
2,2,4-Trimethylhexane	I9	0.0001	0.0008	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0015	0.0100	0.001	0.001
1t,3-Dimethylcyclohexane	N8	0.0001	0.0007	0.000	0.000
UnknownC7s	U7	0.0001	0.0006	0.000	0.000
n-Octane	P8	0.0065	0.0441	0.003	0.003
1c,4-Dimethylcyclohexane	N8	0.0018	0.0120	0.001	0.001
i-Propylcyclopentane	I8	0.0001	0.0007	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0003	0.0023	0.000	0.000
2,2,3,4-Tetramethylpentane	I9	0.0001	0.0008	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0001	0.0008	0.000	0.000
2,2-Dimethylheptane	I9	0.0003	0.0023	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0020	0.0150	0.001	0.001
2,2,3-Trimethylhexane	I9	0.0005	0.0038	0.000	0.000
2,4-Dimethylheptane	I9	0.0006	0.0046	0.000	0.000
4,4-Dimethylheptane	I9	0.0001	0.0008	0.000	0.000
Ethylcyclohexane	N8	0.0013	0.0087	0.001	0.001
n-Propylcyclopentane	N8	0.0008	0.0054	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0003	0.0023	0.000	0.000
2,5-Dimethylheptane	I9	0.0016	0.0122	0.001	0.001
3,3-Dimethylheptane	I9	0.0005	0.0038	0.000	0.000
3,5-Dimethylheptane	I9	0.0002	0.0016	0.000	0.000
2,6-Dimethylheptane	I9	0.0002	0.0016	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0004	0.0030	0.000	0.000
Ethylbenzene	I8	0.0026	0.0164	0.001	0.001
1c,2t,4t-Trimethylcyclohexane	N9	0.0001	0.0008	0.000	0.000
2,3-Dimethylheptane	I9	0.0001	0.0008	0.000	0.000

1,3-Dimethylbenzene (m-Xylene)	A8	0.0177	0.1117	0.007	0.007
1,4-Dimethylbenzene (p-Xylene)	A8	0.0054	0.0341	0.002	0.002
3,4-Dimethylheptane	I9	0.0010	0.0076	0.001	0.001
3,4-Dimethylheptane (2)	I9	0.0003	0.0023	0.000	0.000
4-Ethylheptane	I9	0.0003	0.0023	0.000	0.000
4-Methyloctane	I9	0.0018	0.0137	0.001	0.001
2-Methyloctane	I9	0.0026	0.0198	0.001	0.001
1c,2t,3-Trimethylcyclohexane	N9	0.0001	0.0008	0.000	0.000
3-Ethylheptane	I9	0.0001	0.0008	0.000	0.000
3-Methyloctane	I9	0.0004	0.0030	0.000	0.000
1c,2t,4c-Trimethylcyclohexane	I9	0.0025	0.0188	0.001	0.001
1,1,2-Trimethylcyclohexane	N9	0.0001	0.0008	0.000	0.000
3,3-Diethylpentane	I9	0.0001	0.0008	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0038	0.0240	0.001	0.001
i-Butylcyclopentane	N9	0.0012	0.0090	0.001	0.001
UnknownC8s	U8	0.0003	0.0020	0.000	0.000
n-Nonane	P9	0.0098	0.0748	0.006	0.006
1,1-Methylethylcyclohexane	N9	0.0003	0.0023	0.000	0.000
i-Propylbenzene	A9	0.0003	0.0021	0.000	0.000
i-Propylcyclohexane	N9	0.0003	0.0023	0.000	0.000
2,2-Dimethyloctane	I10	0.0003	0.0026	0.000	0.000
2,4-Dimethyloctane	I10	0.0004	0.0034	0.000	0.000
2,6-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
2,5-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
n-Butylcyclopentane	N9	0.0009	0.0068	0.001	0.001
3,3-Dimethyloctane	I10	0.0007	0.0060	0.000	0.000
n-Propylbenzene	A9	0.0018	0.0129	0.001	0.001
3,6-Dimethyloctane	I10	0.0014	0.0118	0.001	0.001
3-Methyl-5-ethylheptane	I10	0.0002	0.0017	0.000	0.000
1,3-Methylethylbenzene	A9	0.0044	0.0315	0.002	0.002
1,4-Methylethylbenzene	A9	0.0019	0.0136	0.001	0.001
1,3,5-Trimethylbenzene	A9	0.0051	0.0365	0.002	0.002
2,3-Dimethyloctane	I10	0.0009	0.0076	0.001	0.001
5-Methylnonane	I10	0.0016	0.0136	0.001	0.001
1,2-Methylethylbenzene	A9	0.0012	0.0086	0.001	0.001
2-Methylnonane	I10	0.0024	0.0203	0.001	0.001
3-Ethylheptane	I10	0.0006	0.0051	0.000	0.000
3-Methylnonane	I10	0.0019	0.0161	0.001	0.001
1,2,4-Trimethylbenzene	A9	0.0002	0.0014	0.000	0.000
t-Butylbenzene	A10	0.0071	0.0567	0.003	0.003
i-Butylcyclohexane	N10	0.0005	0.0042	0.000	0.000
1t-Methyl-2-n-propylcyclohexane	I10	0.0002	0.0017	0.000	0.000
i-Butylbenzene	A10	0.0005	0.0040	0.000	0.000
sec-Butylbenzene	A10	0.0003	0.0024	0.000	0.000
UnknownC9s	U9	0.0042	0.0321	0.002	0.002
n-Decane	P10	0.0119	0.1007	0.007	0.007
1,2,3-Trimethylbenzene	A9	0.0014	0.0100	0.001	0.001
1,3-Methyl-i-propylbenzene	A10	0.0005	0.0040	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0003	0.0024	0.000	0.000
Sec-Butylcyclohexane	A10	0.0009	0.0075	0.001	0.001
1,2-Methyl-i-propylbenzene	A10	0.0015	0.0120	0.001	0.001
3-Ethylheptane	I10	0.0005	0.0046	0.000	0.000
1,3-Diethylbenzene	A10	0.0002	0.0016	0.000	0.000
1,3-Methyl-n-propylbenzene	A10	0.0001	0.0008	0.000	0.000
1,4-Diethylbenzene	A10	0.0010	0.0080	0.000	0.000
1,4-Methyl-n-propylbenzene	A10	0.0012	0.0096	0.001	0.001
n-Butylbenzene	A10	0.0004	0.0032	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0016	0.0128	0.001	0.001
1,2-Diethylbenzene	A10	0.0007	0.0056	0.000	0.000
t-Decahydronaphthalene	A9	0.0002	0.0018	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0014	0.0112	0.001	0.001
1,3-Dimethyl-4-ethylbenzene	A10	0.0005	0.0040	0.000	0.000

1,2-Dimethyl-4-ethylbenzene	A10	0.0030	0.0240	0.001	0.001
1,3-Dimethyl-2-ethylbenzene	A10	0.0029	0.0231	0.001	0.001
1,2-Dimethyl-3-ethylbenzene	A10	0.0006	0.0048	0.000	0.000
1,2-Ethyl-i-propylbenzene	A10	0.0013	0.0115	0.001	0.001
1,4-Methyl-t-butylbenzene	A11	0.0001	0.0009	0.000	0.000
UnknownC10s	U10	0.0079	0.0668	0.005	0.005
n-Undecane	P11	0.0149	0.1385	0.010	0.010
1,4-Ethyl-i-propylbenzene	A11	0.0012	0.0106	0.001	0.001
1,2,4,5-Tetramethylbenzene	A11	0.0018	0.0144	0.001	0.001
1,2-Methyl-n-butylbenzene	A11	0.0002	0.0018	0.000	0.000
1,2,3,5-Tetramethylbenzene	A11	0.0017	0.0136	0.001	0.001
1,2-Methyl-t-butylbenzene	A11	0.0011	0.0097	0.001	0.001
5-Methylindan	A11	0.0011	0.0086	0.001	0.001
4-Methylindan	A11	0.0007	0.0055	0.000	0.000
1,2-Ethyl-n-propylbenzene	A11	0.0005	0.0044	0.000	0.000
2-Methylindan	A11	0.0009	0.0071	0.001	0.001
1,3-Methyl-n-butylbenzene	A11	0.0003	0.0026	0.000	0.000
1,3-Di-i-propylbenzene	A11	0.0005	0.0048	0.000	0.000
sec-Pentylbenzene	A11	0.0003	0.0026	0.000	0.000
n-Pentylbenzene	A11	0.0030	0.0265	0.002	0.002
1,2-Di-n-propylbenzene	A11	0.0016	0.0155	0.001	0.001
1,4-Di-i-propylbenzene	A11	0.0007	0.0068	0.001	0.001
Tetrahydronaphthalene	A10	0.0009	0.0071	0.001	0.001
Naphthalene	A10	0.0026	0.0198	0.002	0.002
1-t-Butyl-3,5-dimethylbenzene	A12	0.0005	0.0048	0.000	0.000
1,4-Ethyl-t-butylbenzene	A11	0.0009	0.0087	0.001	0.001
1,3-Di-n-propylbenzene	A12	0.0009	0.0087	0.001	0.001
UnknownC11s	U11	0.0075	0.0697	0.005	0.005
n-Dodecane	P12	0.0144	0.1459	0.010	0.010
1,3,5-Triethylbenzene	A12	0.0008	0.0077	0.000	0.000
1,2,4-Triethylbenzene	A12	0.0008	0.0077	0.000	0.000
1,4-Methyl-n-pentylbenzene	A12	0.0018	0.0174	0.001	0.001
n-Hexylbenzene	A12	0.0022	0.0212	0.001	0.001
1,2,3,4,5-Pentamethylbenzene	A13	0.0006	0.0053	0.000	0.000
2-Methylnaphthalene	A11	0.0010	0.0084	0.001	0.001
1-Methylnaphthalene	A11	0.0008	0.0068	0.001	0.001
UnknownC12s	U12	0.0195	0.1813	0.013	0.013
n-Tridecane	P13	0.0088	0.0965	0.007	0.007
UnknownC13s	U13	0.0160	0.1754	0.012	0.012
n-Tetradecane	P14	0.0015	0.0177	0.001	0.001
UnknownC14s	U14	0.0045	0.0531	0.004	0.004
n-Pentadecane	P15	0.0003	0.0038	0.000	0.000
UnknownC15s	U15	0.0004	0.0051	0.000	0.000
n-Hexadecane	P16	0.0001	0.0014	0.000	0.000
TOTAL		100.00000	100.00000	0.7394	0.7424

BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0169	0.0785
TOLUENE	0.0328	0.1797
ETHYLBENZENE	0.0026	0.0164
XYLENES	0.0269	0.1698
TOTAL BTEX	0.0792	0.4444

*(DETAILED HYDROCARBON ANALYSIS/NJ 1993)
 Mod ASTM D6730, GPA 2261 & GPA 2286.

** (CALC: GPA 2172, GPA 2145 & TP-17 @14.696 & 60 F)

CALCULATED VALUES**

BTU @	14.65	14.73
LHV NET DRY REAL :	944.2 /scf	949.4 /scf
NET WET REAL :	927.7 /scf	932.9 /scf
HHV GROSS DRY REAL :	1046.4 /scf	1052.1 /scf
GROSS WET REAL :	1028.1 /scf	1033.8 /scf
NET HEATING VALUE (60 °F ideal reaction):		21347.1 Btu/lbm
GROSS HEATING VALUE (60°F ideal reaction):		23674.6 Btu/lbm
RELATIVE DENSITY (AIR=1):		0.5793
DENSITY		0.04431 lb/scf
COMPRESSIBILITY FACTOR :		0.9979
REGULAR WOBBE INDEX		1376.3

C6+ Fraction of DHA Gas Analysis @60°F, 14.696 psia

Net Dry Ideal BTU	<u>6229.5</u> /scf	Relative Density - SG (Air=1)	<u>4.3211</u>	C6+ factors
Gross Dry Ideal BTU	<u>6698.2</u> /scf	Z Compressibility Factor	<u>0.99834</u>	<u>0.99743</u>
Net Dry Ideal BTU	<u>18738.1</u> /lb	Density Factor	<u>329.821</u> lbm/1000 ft3	
Gross Dry Ideal BTU	<u>20141.3</u> /lb	Molar Mass or MW	<u>125.159</u> g/mol	
		Volume Liquid Ideal gas	<u>0.205</u> scf/gal	<u>19.3</u>

**This hexanes plus fraction may be applied in place of published C6+ factors. The Z & GPM need additional calc for C6+ factors.
 #DIV/0 or 0 (zero) will appear in this section when there is no hexanes plus in the sample to calculate C6+ factors.**

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