



EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PRIMARY DB KEY: **05-103-11266** NAME/DESCRIP : **120196002 YCF XOM 2-42-1**
 LEASE #: **COC-68957A** **BRADEN HEAD**
 FIELD/AREA: **YELLOW CREEK**

PROJECT NO. : **202603075** ANALYSIS NO. : **01**
 COMPANY NAME : **QB ENERGY OPERATING, LLC** ANALYSIS DATE: **APRIL 09, 2026 06:49**
 OFFICE / BRANCH: **PARACHUTE, CO** SAMPLE DATE : **MARCH 4, 2026**
 CUSTOMER REF: TO:
 PRODUCER : **QB ENERGY OPERATING LLC** EFFECTIVE DATE:

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

SAMPLE CYCLE: SAMPLE TYPE: **SPOT**
 SAMPLE PRES. : 545 psig PROBE : **NO**
 FLOW PRES. : psig CYLINDER NO. : **ECA-815**
 LAB PRES: psig SAMPLED BY : **MIKE KELLEY**
 SAMPLE TEMP. : 57 °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 @ 14.65</u>	<u>GPM @ 14.73</u>
HELIUM	0.01	0.00	---	---
HYDROGEN	0.01	0.00	---	---
OXYGEN/ARGON	0.01	0.02	---	---
NITROGEN	0.33	0.53	---	---
CARBON DIOXIDE	0.05	0.12	---	---
METHANE	91.9984	83.8398	---	---
ETHANE	5.4196	9.2573	1.4450	1.4529
PROPANE	1.3668	3.4237	0.3757	0.3778
I-BUTANE	0.2800	0.9245	0.0909	0.0914
N-BUTANE	0.3559	1.1751	0.1119	0.1125
I-PENTANE	0.0990	0.4058	0.0360	0.0362
N-PENTANE	0.0546	0.2238	0.0200	0.0201
HEXANES PLUS	0.0157	0.0800	0.0060	0.0060
TOTALS	100.0000	100.0000	2.0855	2.0969

<u>BTEX COMPONENTS</u>	<u>MOLE%</u>	<u>WT%</u>
BENZENE	0.0000	0.0000
TOLUENE	0.0000	0.0000
ETHYLBENZENE	0.0000	0.0000
XYLENES	0.0000	0.0000
TOTAL BTEX	0.0000	0.0000

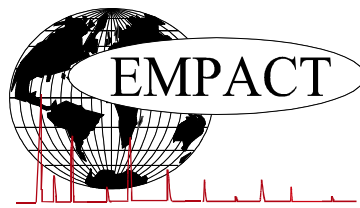
	<u>CALCULATED VALUES**</u>	
	<u>BTU @ 14.65</u>	<u>14.73</u>
LHV NET DRY REAL :	980.6 /scf	986.0 /scf
NET WET REAL :	963.5 /scf	968.9 /scf
HHV GROSS DRY REAL :	1086.4 /scf	1092.4 /scf
GROSS WET REAL :	1067.4 /scf	1073.4 /scf
NET HEATING VALUE (60 °F ideal reaction):	21158.4 Btu/lbm	21158.4 Btu/lbm
GROSS HEATING VALUE (60°F ideal reaction):	23435.7 Btu/lbm	23435.7 Btu/lbm
RELATIVE DENSITY (AIR=1):	0.6078	0.6078
DENSITY	0.04639 lbm/scf	0.04639 lbm/scf
COMPRESSIBILITY FACTOR :	0.9976	0.9976
REGULAR WOBBE INDEX	1394.5	1394.5

*(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. :	202603075	ANALYSIS NO. :	01
COMPANY NAME :	QB ENERGY OPERATING, LLC	ANALYSIS DATE:	APRIL 09, 2026 06:49
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2026
PRODUCER :	QB ENERGY OPERATING LLC	CYLINDER NO. :	ECA-815
LEASE NO. :	COC-68957A	SAMPLED BY :	MIKE KELLEY
NAME/DESCRIP :	120196002 YCF XOM 2-42-1 BRADEN HEAD		

FIELD DATA		SAMPLE TEMP. :	57
SAMPLE PRES. :	545	AMBIENT TEMP.:	
H2S BY STAIN TUBE:	—		
COMMENTS :	<i>SPOT</i> <i>ppm mol</i> <i>NO PROBE</i>		

<u>Componet</u>	<u>Mole %</u>	<u>Wt %</u>
Helium	0.01	0.00
Hydrogen	0.01	0.00
Carbon Dioxide	0.05	0.12
Nitrogen	0.33	0.53
Methane	91.9984	83.8398
Ethane	5.4196	9.2573
Propane	1.3668	3.4237
Isobutane	0.2800	0.9245
n-Butane	0.3559	1.1751
Isopentane	0.0989	0.4054
n-Pentane	0.0546	0.2238
Cyclopentane	0.0001	0.0004
n-Hexane	0.0033	0.0161
Cyclohexane	0.0001	0.0005
Other Hexanes	0.0111	0.0544
Heptanes	0.0004	0.0024
Methylcyclohexane	0.0001	0.0006
2,2,4 Trimethylpentane	0.0000	0.0000
Benzene	0.0000	0.0000
Toluene	0.0000	0.0000
Ethylbenzene	0.0000	0.0000
Xylenes	0.0000	0.0000
C8+ Heavies	0.0007	0.0060
<u>Subtotal</u>	<u>99.99000</u>	<u>99.98000</u>
<u>Oxygen/Argon</u>	<u>0.01</u>	<u>0.02</u>
Total	100.00000	100.00000

	Total	C6+	C8+	C10+
Calculated Values BTU @	Sample	Fraction	Fraction	Fraction
14.65				
LHV Net Dry Real:	980.6	4602.8	7566.1	7747.4 Btu/scf
Net Wet Real:	963.5	4522.3	7433.8	7612.0 Btu/scf
HHV Gross Dry Real:	1086.4	4968.8	8146.6	8340.9 Btu/scf
Gross Wet Real:	1067.4	4881.9	8004.2	8195.1 Btu/scf
Other Calculated Values				
Regualr Wobbe Index*	1394.5	2798.6	3586.3	3628.0 Btu/scf
Net Heating Value (60 °F ideal reaction):	21158.4	19329.1	19156.4	19149.3 Btu/lbm
Gross Heating Value (60°F ideal reaction):	23435.7	20867.7	20626.1	20616.0 Btu/lbm
Molar Mass (MW):	17.60383	89.429	150.299	153.974 g/mol
Relative Density (AIR=1):	0.6078	3.0878	5.1894	5.3163 SG
Density:	0.04639	0.23567	0.39606	0.40574 lbm/scf
Compressibility Factor:	0.9976	0.9866	0.9997	0.9998 Z
Liquid Volume real gas @:	14.65	17.6366	0.006	0 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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EXTENDED NATURAL GAS ANALYSIS (*DHA)

DHA COMPONENT LIST

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 FIELD/AREA: **YELLOW CREEK**

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 PRODUCER : **QB ENERGY OPERATING LLC** EFFECTIVE DATE:
*****FIELD DATA*****
 SAMPLE CYCLE: SAMPLE TYPE: **SPOT**
 SAMPLE PRES. : **545** psig PROBE : **NO**
 FLOW PRES. : psig CYLINDER NO. : **ECA-815**
 LAB PRES: psig SAMPLED BY : **MIKE KELLEY**
 SAMPLE TEMP. : **57** °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:

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.65	GPM @ 14.73
Helium	---	0.01	0.00	---	---
Hydrogen	---	0.01	0.00	---	---
Oxygen/Argon	---	0.01	0.02	---	---
Nitrogen	---	0.33	0.53	---	---
Carbon Dioxide	---	0.05	0.12	---	---
Methane	P1	91.9984	83.8398	---	---
Ethane	P2	5.4196	9.2573	1.445	1.453
Propane	P3	1.3668	3.4237	0.376	0.378
i-Butane	I4	0.2800	0.9245	0.091	0.091
n-Butane	P4	0.3558	1.1748	0.112	0.113
2,2-Dimethylpropane	I5	0.0039	0.0160	0.001	0.001
i-Pentane	I5	0.0950	0.3894	0.035	0.035
UnknownC4s	U4	0.0001	0.0003	0.000	0.000
n-Pentane	P5	0.0546	0.2238	0.020	0.020
2,2-Dimethylbutane	I6	0.0015	0.0073	0.001	0.001
Cyclopentane	N5	0.0001	0.0004	0.000	0.000
2,3-Dimethylbutane	I6	0.0013	0.0064	0.001	0.001
2-Methylpentane	I6	0.0058	0.0284	0.002	0.002
3-Methylpentane	I6	0.0023	0.0113	0.001	0.001
n-Hexane	P6	0.0033	0.0161	0.001	0.001
Methylcyclopentane	N6	0.0002	0.0010	0.000	0.000
2,4-Dimethylpentane	I7	0.0001	0.0006	0.000	0.000
Cyclohexane	N6	0.0001	0.0005	0.000	0.000
2-Methylhexane	I7	0.0001	0.0006	0.000	0.000
3-Methylhexane	I7	0.0001	0.0006	0.000	0.000
n-Heptane	P7	0.0001	0.0006	0.000	0.000
Methylcyclohexane	N7	0.0001	0.0006	0.000	0.000

n-Nonane	P9	0.0001	0.0007	0.000	0.000
n-Decane	P10	0.0001	0.0008	0.000	0.000
UnknownC10s	U10	0.0001	0.0008	0.000	0.000
n-Undecane	P11	0.0002	0.0018	0.000	0.000
UnknownC11s	U11	0.0001	0.0009	0.000	0.000
n-Dodecane	P12	0.0001	0.0010	0.000	0.000
TOTAL		100.00000	100.00000	2.0855	2.0969

CALCULATED VALUES**

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.65	14.73
BENZENE	0.0000	0.0000	LHV NET DRY REAL :	980.6 /scf	986.0 /scf
TOLUENE	0.0000	0.0000	NET WET REAL :	963.5 /scf	968.9 /scf
ETHYLBENZENE	0.0000	0.0000	HHV GROSS DRY REAL :	1086.4 /scf	1092.4 /scf
XYLENES	0.0000	0.0000	GROSS WET REAL :	1067.4 /scf	1073.4 /scf
TOTAL BTEX	0.0000	0.0000	NET HEATING VALUE (60 °F ideal reaction):		21158.4 Btu/lbm
			GROSS HEATING VALUE (60°F ideal reaction):		23435.7 Btu/lbm
			RELATIVE DENSITY (AIR=1):		0.6078
			DENSITY		0.04639 lb/scf
			COMPRESSIBILITY FACTOR :		0.9976
			REGULAR WOBBE INDEX		1394.5

*(DETAILED HYDROCARBON ANALYSIS/NJ 1993)

Mod ASTM D6730, GPA 2261 & GPA 2286.

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

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

Net Dry Ideal BTU	<u>4555.5</u> /scf	Relative Density - SG (Air=1)	<u>3.0878</u>	C6+factors
Gross Dry Ideal BTU	<u>4917.8</u> /scf	Z Compressibility Factor	<u>0.98663</u>	<u>0.98605</u>
Net Dry Ideal BTU	<u>19329.1</u> /lb	Density Factor	<u>235.666</u> lbm/1000 ft3	
Gross Dry Ideal BTU	<u>20867.7</u> /lb	Molar Mass or MW	<u>89.429</u> g/mol	
		Volume Liquid Ideal gas	<u>0.006</u> scf/gal	<u>23.8</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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