



**EXTENDED NATURAL GAS ANALYSIS (\*DHA)**

**MAIN PAGE**

PRIMARY DB KEY:	<b>05-045-09022</b>	NAME/DESCRIP :	<b>GG3 CHEVRON 33-7D</b>
LEASE #:	<b>110170021</b>		<b>BRAIDEN HEAD</b>
FIELD/AREA:	<b>GRAND VALLEY - #31290</b>		
PROJECT NO. :	<b>202405125</b>	ANALYSIS NO. :	<b>02</b>
COMPANY NAME :	<b>CAERUS OIL &amp; GAS LLC</b>	ANALYSIS DATE:	<b>MAY 30, 2024 14:38</b>
OFFICE / BRANCH:	<b>PARACHUTE, CO</b>	SAMPLE DATE :	<b>MAY 23, 2024 7:30</b>
CUSTOMER REF:		TO:	
PRODUCER :	<b>CAERUS PICEANCE LLC</b>	EFFECTIVE DATE:	

**\*\*\*FIELD DATA\*\*\***

SAMPLE CYCLE:		SAMPLE TYPE:	<b>SPOT</b>
SAMPLE PRES. :	111 psig	PROBE :	<b>NO</b>
FLOW PRES. :	psig	CYLINDER NO. :	<b>ECA-743</b>
LAB PRES:	psig	SAMPLED BY :	<b>ALEX GALLEGOS</b>
SAMPLE TEMP. :	43 °f	SAMPLING COMPANY:	<b>CAERUS OIL &amp; GAS LLC</b>
AMBIENT TEMP.:	°f	H2S BY STAIN TUBE:	<b>- ppm mol</b>
H2O BY STAIN TUBE:	- #/mmcf	CO2 BY STAIN TUBE:	<b>- Mol %</b>
FIELD COMMENTS:			
LAB COMMENTS:			

<u>COMPONENT</u>	<u>MOLE %</u>	<u>MASS %</u>	<u>GPM @ 14.65</u>	<u>GPM @ 14.73</u>
HELIUM	0.00	0.00	---	---
HYDROGEN	0.01	0.00	---	---
OXYGEN/ARGON	0.00	0.00	---	---
NITROGEN	0.25	0.38	---	---
CARBON DIOXIDE	0.35	0.84	---	---
METHANE	91.4210	80.1488	---	---
ETHANE	4.0405	6.6395	1.0774	1.0833
PROPANE	2.1406	5.1583	0.5877	0.5909
I-BUTANE	0.3450	1.0958	0.1129	0.1136
N-BUTANE	0.6264	1.9896	0.1969	0.1980
I-PENTANE	0.2049	0.8068	0.0750	0.0753
N-PENTANE	0.1859	0.7329	0.0670	0.0673
HEXANES PLUS	0.4257	2.2083	0.1720	0.1723
<b>TOTALS</b>	<b>100.0000</b>	<b>100.0000</b>	<b>2.2889</b>	<b>2.3007</b>

<u>BTEX COMPONENTS</u>	<u>MOLE%</u>	<u>WT%</u>
BENZENE	0.0350	0.1494
TOLUENE	0.0276	0.1390
ETHYLBENZENE	0.0055	0.0319
XYLENES	0.0158	0.0917
<b>TOTAL BTEX</b>	<b>0.0839</b>	<b>0.4120</b>

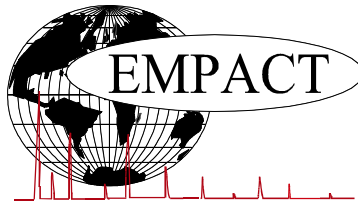
	<u>BTU @ 14.65</u>	<u>14.73</u>
<b>LHV NET DRY REAL :</b>	<b>1009.3 /scf</b>	<b>1014.8 /scf</b>
<b>NET WET REAL :</b>	<b>991.7 /scf</b>	<b>997.2 /scf</b>
<b>HHV GROSS DRY REAL :</b>	<b>1116.5 /scf</b>	<b>1122.5 /scf</b>
<b>GROSS WET REAL :</b>	<b>1097.0 /scf</b>	<b>1103.0 /scf</b>
<b>NET HEATING VALUE (60 °F ideal reaction):</b>	<b>20951.8 Btu/lbm</b>	<b>20951.8 Btu/lbm</b>
<b>GROSS HEATING VALUE (60°F ideal reaction):</b>	<b>23181.1 Btu/lbm</b>	<b>23181.1 Btu/lbm</b>
<b>RELATIVE DENSITY (AIR=1):</b>	<b>0.6310</b>	<b>0.6310</b>
<b>DENSITY</b>	<b>0.04822 lbm/scf</b>	<b>0.04822 lbm/scf</b>
<b>COMPRESSIBILITY FACTOR :</b>	<b>0.9975</b>	<b>0.9975</b>
<b>REGULAR WOBBE INDEX</b>	<b>1406.3</b>	<b>1406.3</b>

\*(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. :	202405125	ANALYSIS NO. :	02
COMPANY NAME :	CAERUS OIL & GAS LLC	ANALYSIS DATE :	MAY 30, 2024 14:38
ACCOUNT NO. :		SAMPLE DATE :	MAY 23, 2024 7:30
PRODUCER :	CAERUS PICEANCE LLC	CYLINDER NO. :	ECA-743
LEASE NO. :	110170021	SAMPLED BY :	ALEX GALLEGOS
NAME/DESCRIP :	GG3 CHEVRON 33-7D BRAIDEN HEAD		

***FIELD DATA***		SAMPLE TEMP. :	43
SAMPLE PRES. :	111	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	0.00	0.00
Hydrogen	0.01	0.00
Carbon Dioxide	0.35	0.84
Nitrogen	0.25	0.38
Methane	91.4210	80.1488
Ethane	4.0405	6.6395
Propane	2.1406	5.1583
Isobutane	0.3450	1.0958
n-Butane	0.6264	1.9896
Isopentane	0.1954	0.7704
n-Pentane	0.1859	0.7329
Cyclopentane	0.0095	0.0364
n-Hexane	0.0511	0.2407
Cyclohexane	0.0219	0.1007
Other Hexanes	0.0991	0.4641
Heptanes	0.0684	0.3730
Methylcyclohexane	0.0348	0.1867
2,2,4 Trimethylpentane	0.0000	0.0000
Benzene	0.0350	0.1494
Toluene	0.0276	0.1390
Ethylbenzene	0.0055	0.0319
Xylenes	0.0158	0.0917
C8+ Heavies	0.0665	0.4311
<u>Subtotal</u>	<u>100.00000</u>	<u>100.00000</u>
<u>Oxygen/Argon</u>	<u>0.00</u>	<u>0.00</u>
<b>Total</b>	<b>100.00000</b>	<b>100.00000</b>

	<b>Total</b>	<b>C6+</b>	<b>C8+</b>	<b>C10+</b>	
<b>Calculated Values BTU @</b>	<b>Sample</b>	<b>Fraction</b>	<b>Fraction</b>	<b>Fraction</b>	
<b>14.65</b>					
LHV Net Dry Real:	1009.3	4737.4	5701.3	7064.3	Btu/scf
Net Wet Real:	991.7	4654.6	5601.6	6940.8	Btu/scf
HHV Gross Dry Real:	1116.5	5077.7	6107.5	7599.3	Btu/scf
Gross Wet Real:	1097.0	4988.9	6000.7	7466.5	Btu/scf
<b>Other Calculated Values</b>					
Regualr Wobbe Index*	1406.3	2794.8	3060.4	3443.9	Btu/scf
Net Heating Value (60 °F ideal reaction):	20951.8	19095.5	19256.0	19092.0	Btu/lbm
Gross Heating Value (60°F ideal reaction):	23181.1	20470.0	20634.7	20539.5	Btu/lbm
Molar Mass (MW):	18.29975	94.911	115.597	141.746	g/mol
Relative Density (AIR=1):	0.6310	3.2775	3.9914	4.8941	SG
Density:	0.04822	0.25013	0.30463	0.37352	lbm/scf
Compressibility Factor:	0.9975	0.9933	0.9980	0.9994	Z
Liquid Volume real gas @:	<b>14.65</b>	17.7812	0.1715	0.0389	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**

PRIMARY DB KEY:	<b>05-045-09022</b>	NAME/DESCRIP :	<b>GG3 CHEVRON 33-7D</b>
LEASE #:	<b>110170021</b>		<b>BRAIDEN HEAD</b>
FIELD/AREA:	<b>GRAND VALLEY - #31290</b>		
PROJECT NO. :	<b>202405125</b>	ANALYSIS NO. :	<b>02</b>
COMPANY NAME :	<b>CAERUS OIL &amp; GAS LLC</b>	ANALYSIS DATE:	<b>MAY 30, 2024 14:38</b>
OFFICE / BRANCH:	<b>PARACHUTE, CO</b>	SAMPLE DATE :	<b>MAY 23, 2024 7:30</b>
CUSTOMER REF:		TO:	
PRODUCER :	<b>CAERUS PICEANCE LLC</b>	EFFECTIVE DATE:	
<b>***FIELD DATA***</b>			
SAMPLE CYCLE:		SAMPLE TYPE:	<b>SPOT</b>
SAMPLE PRES. :	111 psig	PROBE :	<b>NO</b>
FLOW PRES. :	psig	CYLINDER NO. :	<b>ECA-743</b>
LAB PRES:	psig	SAMPLED BY :	<b>ALEX GALLEGOS</b>
SAMPLE TEMP. :	43 °f	SAMPLING COMPANY:	<b>CAERUS OIL &amp; GAS LLC</b>
AMBIENT TEMP.:	°f	H2S BY STAIN TUBE:	<b>- ppm mol</b>
H2O BY STAIN TUBE:	- #/mmcf	CO2 BY STAIN TUBE:	<b>- Mol %</b>
FIELD COMMENTS:			
LAB COMMENTS:			

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.65	GPM @ 14.73
Helium	---	0.00	0.00	---	---
Hydrogen	---	0.01	0.00	---	---
Oxygen/Argon	---	0.00	0.00	---	---
Nitrogen	---	0.25	0.38	---	---
Carbon Dioxide	---	0.35	0.84	---	---
Methane	P1	91.4210	80.1488	---	---
Ethane	P2	4.0405	6.6395	1.077	1.083
Propane	P3	2.1406	5.1583	0.588	0.591
i-Butane	I4	0.3450	1.0958	0.113	0.114
n-Butane	P4	0.6264	1.9896	0.197	0.198
2,2-Dimethylpropane	I5	0.0015	0.0059	0.001	0.001
i-Pentane	I5	0.1939	0.7645	0.071	0.071
n-Pentane	P5	0.1859	0.7329	0.067	0.067
2,2-Dimethylbutane	I6	0.0018	0.0085	0.001	0.001
Cyclopentane	N5	0.0095	0.0364	0.003	0.003
2,3-Dimethylbutane	I6	0.0066	0.0311	0.003	0.003
2-Methylpentane	I6	0.0437	0.2058	0.018	0.018
3-Methylpentane	I6	0.0228	0.1074	0.009	0.009
n-Hexane	P6	0.0511	0.2407	0.021	0.021
2,2-Dimethylpentane	I7	0.0006	0.0033	0.000	0.000
Methylcyclopentane	N6	0.0242	0.1113	0.009	0.009
2,4-Dimethylpentane	I7	0.0017	0.0093	0.001	0.001
2,2,3-Trimethylbutane	I7	0.0003	0.0016	0.000	0.000
Benzene	A6	0.0350	0.1494	0.010	0.010
3,3-Dimethylpentane	I7	0.0004	0.0022	0.000	0.000
Cyclohexane	N6	0.0219	0.1007	0.007	0.007
2-Methylhexane	I7	0.0107	0.0586	0.005	0.005
2,3-Dimethylpentane	I7	0.0039	0.0214	0.002	0.002
1,1-Dimethylcyclopentane	N7	0.0017	0.0091	0.001	0.001
3-Methylhexane	I7	0.0106	0.0580	0.005	0.005

1c,3-Dimethylcyclopentane	N7	0.0039	0.0209	0.002	0.002
1t,3-Dimethylcyclopentane	N7	0.0035	0.0188	0.002	0.002
3-Ethylpentane	I7	0.0005	0.0027	0.000	0.000
1t,2-Dimethylcyclopentane	N7	0.0062	0.0333	0.003	0.003
n-Heptane	P7	0.0218	0.1194	0.010	0.010
1c,2-Dimethylcyclopentane	N7	0.0005	0.0027	0.000	0.000
Methylcyclohexane	N7	0.0348	0.1867	0.014	0.014
2,2-Dimethylhexane	I8	0.0011	0.0069	0.001	0.001
1,1,3-Trimethylcyclopentane	N7	0.0005	0.0031	0.000	0.000
Ethylcyclopentane	N7	0.0016	0.0086	0.001	0.001
2,5-Dimethylhexane	I8	0.0009	0.0056	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0009	0.0056	0.000	0.000
2,4-Dimethylhexane	I8	0.0001	0.0006	0.000	0.000
1c,2t,4-Trimethylcyclopentane	N8	0.0014	0.0086	0.001	0.001
3,3-Dimethylhexane	I8	0.0002	0.0013	0.000	0.000
2,3,3-Trimethylpentane	I8	0.0001	0.0006	0.000	0.000
Toluene	A7	0.0276	0.1390	0.009	0.009
2,3-Dimethylhexane	I8	0.0011	0.0069	0.001	0.001
2-Methyl-3-ethylpentane	I8	0.0004	0.0025	0.000	0.000
2-Methylheptane	I8	0.0056	0.0350	0.003	0.003
4-Methylheptane	I8	0.0013	0.0081	0.001	0.001
3-Methyl-3-ethylpentane	I8	0.0002	0.0013	0.000	0.000
3,4-Dimethylhexane	I8	0.0003	0.0019	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0001	0.0006	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0001	0.0006	0.000	0.000
3-Methylheptane	I8	0.0031	0.0193	0.002	0.002
1c,2t,3-Trimethylcyclopentane	N8	0.0046	0.0282	0.002	0.002
3-Ethylhexane	I8	0.0005	0.0031	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0021	0.0129	0.001	0.001
1,1-Dimethylcyclohexane	N8	0.0007	0.0043	0.000	0.000
2,2,5-Trimethylhexane	I9	0.0001	0.0007	0.000	0.000
3c-Ethylmethylcyclopentane	N8	0.0004	0.0025	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0004	0.0025	0.000	0.000
2t-Ethylmethylcyclopentane	N8	0.0006	0.0037	0.000	0.000
1,1-Methylethylcyclopentane	N8	0.0001	0.0006	0.000	0.000
2,2,4-Trimethylhexane	I9	0.0001	0.0007	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0019	0.0116	0.001	0.001
1t,3-Dimethylcyclohexane	N8	0.0001	0.0006	0.000	0.000
n-Octane	P8	0.0093	0.0580	0.005	0.005
1c,4-Dimethylcyclohexane	N8	0.0011	0.0067	0.001	0.001
i-Propylcyclopentane	I8	0.0001	0.0006	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0002	0.0014	0.000	0.000
2,2,3,4-Tetramethylpentane	I9	0.0001	0.0007	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0001	0.0007	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0001	0.0006	0.000	0.000
2,2-Dimethylheptane	I9	0.0007	0.0049	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0032	0.0221	0.002	0.002
2,2,3-Trimethylhexane	I9	0.0005	0.0035	0.000	0.000
2,4-Dimethylheptane	I9	0.0002	0.0014	0.000	0.000
4,4-Dimethylheptane	I9	0.0001	0.0007	0.000	0.000
Ethylcyclohexane	N8	0.0029	0.0178	0.001	0.001
n-Propylcyclopentane	N8	0.0010	0.0061	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0003	0.0021	0.000	0.000
2,5-Dimethylheptane	I9	0.0014	0.0098	0.001	0.001
3,3-Dimethylheptane	I9	0.0002	0.0014	0.000	0.000
3,5-Dimethylheptane	I9	0.0001	0.0007	0.000	0.000
2,6-Dimethylheptane	I9	0.0002	0.0014	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0002	0.0014	0.000	0.000
Ethylbenzene	I8	0.0055	0.0319	0.002	0.002
1c,2t,4t-Trimethylcyclohexane	N9	0.0001	0.0007	0.000	0.000
2,3-Dimethylheptane	I9	0.0001	0.0007	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0092	0.0534	0.004	0.004
1,4-Dimethylbenzene (p-Xylene)	A8	0.0021	0.0122	0.001	0.001

3,4-Dimethylheptane	I9	0.0001	0.0007	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0005	0.0035	0.000	0.000
4-Ethylheptane	I9	0.0002	0.0014	0.000	0.000
4-Methyloctane	I9	0.0010	0.0070	0.001	0.001
2-Methyloctane	I9	0.0016	0.0112	0.001	0.001
1c,2t,3-Trimethylcyclohexane	N9	0.0002	0.0014	0.000	0.000
3-Ethylheptane	I9	0.0002	0.0014	0.000	0.000
3-Methyloctane	I9	0.0003	0.0021	0.000	0.000
1c,2t,4c-Trimethylcyclohexane	I9	0.0016	0.0110	0.001	0.001
1,1,2-Trimethylcyclohexane	N9	0.0001	0.0007	0.000	0.000
3,3-Diethylpentane	I9	0.0002	0.0014	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0045	0.0261	0.002	0.002
i-Butylcyclopentane	N9	0.0011	0.0076	0.001	0.001
n-Nonane	P9	0.0027	0.0189	0.002	0.002
1,1-Methylethylcyclohexane	N9	0.0005	0.0034	0.000	0.000
i-Propylbenzene	A9	0.0008	0.0053	0.000	0.000
i-Propylcyclohexane	N9	0.0003	0.0021	0.000	0.000
2,2-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
2,6-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
n-Butylcyclopentane	N9	0.0004	0.0027	0.000	0.000
3,6-Dimethyloctane	I10	0.0005	0.0039	0.000	0.000
1,3-Methylethylbenzene	A9	0.0002	0.0013	0.000	0.000
1,4-Methylethylbenzene	A9	0.0001	0.0007	0.000	0.000
t-Butylbenzene	A10	0.0001	0.0007	0.000	0.000
UnknownC9s	U9	0.0022	0.0154	0.001	0.001
UnknownC10s	U10	0.0007	0.0055	0.000	0.000
<b>TOTAL</b>		<b>100.0000</b>	<b>100.0000</b>	<b>2.2889</b>	<b>2.3007</b>

**CALCULATED VALUES\*\***

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.65	14.73
BENZENE	0.0350	0.1494	LHV NET DRY REAL :	1009.3 /scf	1014.8 /scf
TOLUENE	0.0276	0.1390	NET WET REAL :	991.7 /scf	997.2 /scf
ETHYLBENZENE	0.0055	0.0319	HHV GROSS DRY REAL :	1116.5 /scf	1122.5 /scf
XYLENES	0.0158	0.0917	GROSS WET REAL :	1097.0 /scf	1103.0 /scf
TOTAL BTEX	0.0839	0.4120	NET HEATING VALUE (60 °F ideal reaction):		20951.8 Btu/lbm
			GROSS HEATING VALUE (60°F ideal reaction):		23181.1 Btu/lbm
			RELATIVE DENSITY (AIR=1):		0.6310
			DENSITY		0.04822 lb/scf
			COMPRESSIBILITY FACTOR :		0.9975
			REGULAR WOBBE INDEX		1406.3

\*(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	4720.5 /scf	Relative Density - SG (Air=1)	3.2775	<b>C6+ factors</b>
Gross Dry Ideal BTU	5059.6 /scf	Z Compressibility Factor	0.99331	0.99246
Net Dry Ideal BTU	19095.5 /lb	Density Factor	250.129 lbm/1000 ft3	
Gross Dry Ideal BTU	20470 /lb	Molar Mass or MW	94.911 g/mol	
		Volume Liquid Ideal gas	0.172 scf/gal	24.4

**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.**

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.