EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PRIMARY DB KEY:	**05-103-10625**	NAME/DESCRIP :	**PICEANCE CREEK UNIT #T35X-11G7**
LEASE #:	**05-103-10625**		**RIO BLANCO #103**
FIELD/AREA:	**PICEANCE CREEK - #68800**		**BRAIDEN HEAD, INTER-CASING**
PROJECT NO. :	**202307022**	ANALYSIS NO. :	**02**
COMPANY NAME :	**CAERUS OIL & GAS LLC**	ANALYSIS DATE:	JULY 12, 2023 07:05
OFFICE / BRANCH:	PARACHUTE, CO	SAMPLE DATE :	JUNE 20, 2023
CUSTOMER REF:		TO:	
PRODUCER :	CAERUS PICEANCE LLC	EFFECTIVE DATE:	
FIELD DATA			
SAMPLE CYCLE:		SAMPLE TYPE:	SPOT
SAMPLE PRES. :	1589 psig	PROBE :	
FLOW PRES. :	psig	CYLINDER NO. :	ECA-701
LAB PRES:	psig	SAMPLED BY :	SHANE COLLETT
SAMPLE TEMP. :	72 ºf	SAMPLING COMPANY:	CAERUS OIL & GAS LLC
AMBIENT TEMP.:	ºf	H2S BY STAIN TUBE:	ppm mol
H2O BY STAIN TUBE:	#/mmcf	CO2 BY STAIN TUBE:	Mol %
FIELD COMMENTS:	***Federal Lease #8920003240***		
LAB COMMENTS:			

COMPONENT	MOLE %	MASS %	GPM @ 14.65	GPM @ 14.73
GLYCOLS	0.0011	0.0061	0.0010	0.0010
ALCOHOLS	0.0303	0.0543	0.0040	0.0040
HELIUM	0.00	0.00	---	---
HYDROGEN	0.05	0.01	---	---
OXYGEN/ARGON	0.00	0.00	---	---
NITROGEN	0.15	0.22	---	---
CARBON DIOXIDE	1.91	4.38	---	---
METHANE	90.3666	75.4821	---	---
ETHANE	4.2267	6.6175	1.1262	1.1323
PROPANE	1.0439	2.3967	0.2868	0.2884
I-BUTANE	0.2494	0.7548	0.0809	0.0814
N-BUTANE	0.2293	0.6939	0.0719	0.0723
I-PENTANE	0.1330	0.4990	0.0480	0.0482
N-PENTANE	0.0986	0.3704	0.0360	0.0362
HEXANES PLUS	1.5111	8.5152	0.6920	0.6944
TOTALS	100.00000	100.00000	2.3468	2.3582

BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0382	0.1554
TOLUENE	0.0734	0.3521
ETHYLBENZENE	0.0060	0.0332
XYLENES	0.0945	0.5223
TOTAL BTEX	0.2121	1.0630

*(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 :	1018.1 /scf	1023.6 /scf
	NET WET REAL :	1000.3 /scf	1005.8 /scf
HHV	GROSS DRY REAL :	1124.6 /scf	1130.7 /scf
	GROSS WET REAL :	1104.9 /scf	1111.0 /scf

NET HEATING VALUE (60 °F ideal reaction):	20161.8 Btu/lbm
GROSS HEATING VALUE (60°F ideal reaction):	22284.5 Btu/lbm
RELATIVE DENSITY (AIR=1):	0.6621
DENSITY	0.05060 lbm/scf
COMPRESSIBILITY FACTOR :	0.9976
REGULAR WOBBE INDEX	1383.1

EMPACT

EXTENDED NATURAL GAS ANALYSIS (*DHA)

GLYCALC INFORMATION

PROJECT NO. :	202307022	ANALYSIS NO. :	02
COMPANY NAME :	CAERUS OIL & GAS LLC	ANALYSIS DATE:	JULY 12, 2023 07:05
ACCOUNT NO. :		SAMPLE DATE :	JUNE 20, 2023
PRODUCER :	CAERUS PICEANCE LLC	CYLINDER NO. :	ECA-701
LEASE NO. :	05-103-10625	SAMPLED BY :	SHANE COLLETT
NAME/DESCRIP :	PICEANCE CREEK UNIT #T35X-11G7 RIO BLANCO #103 BRAIDEN HEAD, INTER-CASING		
FIELD DATA		SAMPLE TEMP. :	72
SAMPLE PRES. :	1589	AMBIENT TEMP.:	
H2S BY STAIN TUBE:	_ *ppm mol*		
COMMENTS :	***SPOT*** ***Federal Lease #8920003240***		

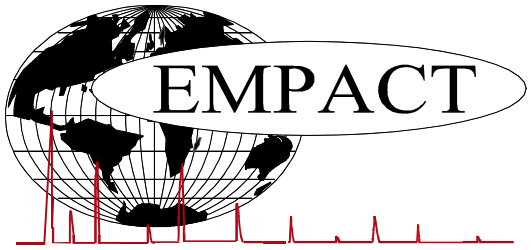
Componet	Mole %	Wt %
Helium	0.00	0.00
Hydrogen	0.05	0.01
Carbon Dioxide	1.91	4.38
Nitrogen	0.15	0.22
Methane	90.3666	75.4821
Ethane	4.2267	6.6175
Propane	1.0439	2.3967
Isobutane	0.2494	0.7548
n-Butane	0.2293	0.6939
Isopentane	0.1269	0.4767
n-Pentane	0.0986	0.3704
Cyclopentane	0.0061	0.0223
n-Hexane	0.0715	0.3208
Cyclohexane	0.0506	0.2218
Other Hexanes	0.1308	0.5835
Heptanes	0.1958	1.0181
Methylcyclohexane	0.1421	0.7264
2,2,4 Trimethylpentane	0.0001	0.0006
Benzene	0.0382	0.1554
Toluene	0.0734	0.3521
Ethylbenzene	0.0060	0.0332
Xylenes	0.0945	0.5223
C8+ Heavies	0.7081	4.5810
Subtotal	99.96860	99.93960
Oxygen/Argon	0.00	0.00
Glycols	0.0011	0.0061
Alcohols	0.0303	0.0543
Total	100.00000	100.00000

Calculated Values BTU @ **14.65**		**Total Sample**	**C6+ Fraction**	**C8+ Fraction**	**C10+ Fraction**	
LHV	Net Dry Real:	1018.1	5393.4	6065.2	7123.0	**Btu/scf**
	Net Wet Real:	1000.3	5299.1	5959.2	6998.5	**Btu/scf**
HHV	Gross Dry Real:	1124.6	5787.6	6512.0	7673.6	**Btu/scf**
	Gross Wet Real:	1104.9	5686.4	6398.2	7539.5	**Btu/scf**
Other Calculated Values						
Regualr Wobbe Index*		1383.1	2992.7	3178.2	3470.1	**Btu/scf**
Net Heating Value (60 °F ideal reaction):		20161.8	19134.0	19205.6	19073.0	**Btu/lbm**
Gross Heating Value (60°F ideal reaction):		22284.5	20537.1	20625.2	20545.8	**Btu/lbm**
Molar Mass (MW):		19.20409	108.228	121.988	142.401	**g/mol**
Relative Density (AIR=1):		0.6621	3.7364	4.2120	4.9166	**SG**
Density:		0.05060	0.28518	0.32145	0.37524	**lbm/scf**
Compressiblity Factor:		0.9976	0.9964	0.9985	0.9996	**Z**
Liquid Volume real gas @:	**14.65**	17.9177	0.6898	0.4107	0.0797	**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 uderscore) indicates there was no tube pulled for H2S.



EXTENDED NATURAL GAS ANALYSIS (*DHA)

DHA COMPONENT LIST

PRIMARY DB KEY:	**05-103-10625**	NAME/DESCRIP :	**PICEANCE CREEK UNIT #T35X-11G7**
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PRODUCER :	CAERUS PICEANCE LLC	EFFECTIVE DATE:	
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SAMPLE PRES. :	1589 psig	PROBE :	
FLOW PRES. :	psig	CYLINDER NO. :	ECA-701
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H2O BY STAIN TUBE:	#/mmcf	CO2 BY STAIN TUBE:	Mol %
FIELD COMMENTS:	***Federal Lease #8920003240***		
LAB COMMENTS:			

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ **14.65**	GPM @ **14.73**
Helium	---	0.00	0.00	---	---
Hydrogen	---	0.05	0.01	---	---
Oxygen/Argon	---	0.00	0.00	---	---
Nitrogen	---	0.15	0.22	---	---
Carbon Dioxide	---	1.91	4.38	---	---
Methane	P1	90.3666	75.4821	---	---
Ethane	P2	4.2267	6.6175	1.126	1.132
Propane	P3	1.0439	2.3967	0.287	0.288
i-Butane	I4	0.2494	0.7548	0.081	0.081
Methanol	X1	0.0284	0.0474	0.004	0.004
n-Butane	P4	0.2293	0.6939	0.072	0.072
2,2-Dimethylpropane	I5	0.0052	0.0195	0.002	0.002
Ethanol	X2	0.0001	0.0003	0.000	0.000
i-Pentane	I5	0.1217	0.4572	0.044	0.044
i-Propanol	X3	0.0004	0.0012	0.000	0.000
n-Pentane	P5	0.0986	0.3704	0.036	0.036
t-Butanol	X4	0.0013	0.0050	0.000	0.000
2,2-Dimethylbutane	I6	0.0080	0.0359	0.003	0.003
Cyclopentane	N5	0.0061	0.0223	0.002	0.002
2,3-Dimethylbutane	I6	0.0117	0.0525	0.005	0.005
2-Methylpentane	I6	0.0497	0.2230	0.021	0.021
3-Methylpentane	I6	0.0291	0.1306	0.012	0.012
n-Hexane	P6	0.0715	0.3208	0.029	0.029
2-Butanol	X4	0.0001	0.0004	0.000	0.000
2,2-Dimethylpentane	I7	0.0024	0.0125	0.001	0.001
Methylcyclopentane	N6	0.0323	0.1415	0.011	0.011
2,4-Dimethylpentane	I7	0.0051	0.0266	0.002	0.002
2,2,3-Trimethylbutane	I7	0.0013	0.0068	0.001	0.001
Benzene	A6	0.0382	0.1554	0.011	0.011
3,3-Dimethylpentane	I7	0.0022	0.0115	0.001	0.001

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Cyclohexane	N6	0.0506	0.2218	0.017	0.017
2-Methylhexane	I7	0.0269	0.1403	0.012	0.012
2,3-Dimethylpentane	I7	0.0098	0.0511	0.004	0.004
1,1-Dimethylcyclopentane	N7	0.0057	0.0292	0.002	0.002
3-Methylhexane	I7	0.0276	0.1440	0.013	0.013
1c,3-Dimethylcyclopentane	N7	0.0079	0.0404	0.004	0.004
1t,3-Dimethylcyclopentane	N7	0.0073	0.0373	0.003	0.003
3-Ethylpentane	I7	0.0008	0.0042	0.000	0.000
1t,2-Dimethylcyclopentane	N7	0.0128	0.0654	0.006	0.006
2,2,4-Trimethylpentane	I8	0.0001	0.0006	0.000	0.000
n-Heptane	P7	0.0784	0.4090	0.036	0.036
1c,2-Dimethylcyclopentane	N7	0.0015	0.0076	0.001	0.001
Methylcyclohexane	N7	0.1421	0.7264	0.057	0.057
2,2-Dimethylhexane	I8	0.0040	0.0238	0.002	0.002
1,1,3-Trimethylcyclopentane	N7	0.0014	0.0082	0.001	0.001
Ethylcyclopentane	N7	0.0046	0.0235	0.002	0.002
2,5-Dimethylhexane	I8	0.0054	0.0321	0.003	0.003
2,2,3-Trimethylpentane	I8	0.0053	0.0315	0.003	0.003
1c,2t,4-Trimethylcyclopentane	N8	0.0033	0.0193	0.002	0.002
3,3-Dimethylhexane	I8	0.0020	0.0119	0.001	0.001
2,3,4-Trimethylpentane	I8	0.0002	0.0012	0.000	0.000
2,3,3-Trimethylpentane	I8	0.0002	0.0012	0.000	0.000
Toluene	A7	0.0734	0.3521	0.025	0.025
2,3-Dimethylhexane	I8	0.0047	0.0280	0.002	0.002
2-Methyl-3-ethylpentane	I8	0.0009	0.0054	0.000	0.000
1,1,2-Trimethylcyclopentane	N8	0.0002	0.0011	0.000	0.000
2-Methylheptane	I8	0.0316	0.1880	0.016	0.016
4-Methylheptane	I8	0.0085	0.0506	0.004	0.004
3-Methyl-3-ethylpentane	I8	0.0018	0.0107	0.001	0.001
3,4-Dimethylhexane	I8	0.0008	0.0047	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0003	0.0018	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0003	0.0018	0.000	0.000
3-Methylheptane	I8	0.0207	0.1231	0.011	0.011
1c,2t,3-Trimethylcyclopentane	N8	0.0349	0.2039	0.018	0.018
3-Ethylhexane	I8	0.0021	0.0125	0.001	0.001
1t,4-Dimethylcyclohexane	N8	0.0140	0.0818	0.007	0.007
1,1-Dimethylcyclohexane	N8	0.0056	0.0327	0.003	0.003
2,2,5-Trimethylhexane	I9	0.0008	0.0054	0.000	0.000
3c-Ethylmethylcyclopentane	N8	0.0013	0.0076	0.001	0.001
3t-Ethylmethylcyclopentane	N8	0.0011	0.0064	0.001	0.001
2t-Ethylmethylcyclopentane	N8	0.0013	0.0076	0.001	0.001
1,1-Methylethylcyclopentane	N8	0.0003	0.0018	0.000	0.000
2,2,4-Trimethylhexane	I9	0.0005	0.0033	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0116	0.0678	0.006	0.006
1t,3-Dimethylcyclohexane	N8	0.0006	0.0035	0.000	0.000
UnknownC7s	U7	0.0001	0.0005	0.000	0.000
n-Octane	P8	0.0841	0.5002	0.043	0.043
1c,4-Dimethylcyclohexane	N8	0.0171	0.0999	0.009	0.009
i-Propylcyclopentane	I8	0.0006	0.0035	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0018	0.0120	0.001	0.001
2,2,3,4-Tetramethylpentane	I9	0.0003	0.0020	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0006	0.0040	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0002	0.0011	0.000	0.000
2,2-Dimethylheptane	I9	0.0053	0.0354	0.003	0.003
1,1,4-Trimethylcyclohexane	N9	0.0141	0.0927	0.007	0.007
2,2,3-Trimethylhexane	I9	0.0039	0.0260	0.002	0.002
2,4-Dimethylheptane	I9	0.0009	0.0060	0.001	0.001
4,4-Dimethylheptane	I9	0.0002	0.0013	0.000	0.000
Ethylcyclohexane	N8	0.0179	0.1046	0.008	0.008
n-Propylcyclopentane	N8	0.0053	0.0310	0.002	0.002
1c,3c,5-Trimethylcyclohexane	N9	0.0013	0.0085	0.001	0.001
2,5-Dimethylheptane	I9	0.0137	0.0915	0.008	0.008
3,3-Dimethylheptane	I9	0.0022	0.0147	0.001	0.001

3,5-Dimethylheptane	I9	0.0005	0.0033	0.000	0.000
2,6-Dimethylheptane	I9	0.0007	0.0047	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0048	0.0316	0.002	0.002
Ethylbenzene	I8	0.0060	0.0332	0.002	0.002
1c,2t,4t-Trimethylcyclohexane	N9	0.0002	0.0013	0.000	0.000
2,3-Dimethylheptane	I9	0.0004	0.0027	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0599	0.3311	0.023	0.023
1,4-Dimethylbenzene (p-Xylene)	A8	0.0230	0.1271	0.009	0.009
3,4-Dimethylheptane	I9	0.0008	0.0054	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0016	0.0107	0.001	0.001
4-Ethylheptane	I9	0.0015	0.0100	0.001	0.001
4-Methyloctane	I9	0.0110	0.0735	0.006	0.006
2-Methyloctane	I9	0.0174	0.1162	0.010	0.010
1c,2t,3-Trimethylcyclohexane	N9	0.0006	0.0040	0.000	0.000
3-Ethylheptane	I9	0.0005	0.0033	0.000	0.000
3-Methyloctane	I9	0.0025	0.0167	0.001	0.001
1c,2t,4c-Trimethylcyclohexane	I9	0.0168	0.1104	0.009	0.009
1,1,2-Trimethylcyclohexane	N9	0.0007	0.0046	0.000	0.000
3,3-Diethylpentane	I9	0.0014	0.0094	0.001	0.001
1,2-Dimethylbenzene (o-Xylene)	A8	0.0116	0.0641	0.004	0.004
i-Butylcyclopentane	N9	0.0081	0.0533	0.004	0.004
n-Nonane	P9	0.0876	0.5850	0.049	0.049
1,1-Methylethylcyclohexane	N9	0.0046	0.0302	0.003	0.003
i-Propylbenzene	A9	0.0019	0.0119	0.001	0.001
i-Propylcyclohexane	N9	0.0019	0.0125	0.001	0.001
2,2-Dimethyloctane	I10	0.0011	0.0082	0.001	0.001
2,4-Dimethyloctane	I10	0.0008	0.0059	0.000	0.000
2,6-Dimethyloctane	I10	0.0032	0.0237	0.002	0.002
2,5-Dimethyloctane	I10	0.0007	0.0052	0.000	0.000
n-Butylcyclopentane	N9	0.0065	0.0427	0.004	0.004
3,3-Dimethyloctane	I10	0.0050	0.0370	0.003	0.003
n-Propylbenzene	A9	0.0123	0.0770	0.005	0.005
Diethylene glycol	GL4	0.0011	0.0061	0.001	0.001
3,6-Dimethyloctane	I10	0.0040	0.0296	0.002	0.002
3-Methyl-5-ethylheptane	I10	0.0017	0.0126	0.001	0.001
1,3-Methylethylbenzene	A9	0.0101	0.0632	0.006	0.006
1,4-Methylethylbenzene	A9	0.0036	0.0225	0.002	0.002
1,3,5-Trimethylbenzene	A9	0.0158	0.0989	0.007	0.007
2,3-Dimethyloctane	I10	0.0024	0.0178	0.001	0.001
1,2-Methylethylbenzene	A9	0.0077	0.0482	0.004	0.004
2-Methylnonane	I10	0.0095	0.0704	0.006	0.006
3-Ethyloctane	I10	0.0019	0.0141	0.001	0.001
3-Methylnonane	I10	0.0061	0.0452	0.004	0.004
1,2,4-Trimethylbenzene	A9	0.0001	0.0006	0.000	0.000
t-Butylbenzene	A10	0.0125	0.0874	0.006	0.006
i-Butylcyclohexane	N10	0.0018	0.0131	0.001	0.001
1t-Methyl-2-n-propylcyclohexane	I10	0.0008	0.0058	0.000	0.000
i-Butylbenzene	A10	0.0007	0.0049	0.000	0.000
sec-Butylbenzene	A10	0.0004	0.0028	0.000	0.000
UnknownC9s	U9	0.0116	0.0775	0.007	0.007
n-Decane	P10	0.0348	0.2578	0.021	0.021
1,2,3-Trimethylbenzene	A9	0.0016	0.0100	0.001	0.001
1,3-Methyl-i-propylbenzene	A10	0.0005	0.0035	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0051	0.0357	0.003	0.003
1,2-Methyl-i-propylbenzene	A10	0.0012	0.0084	0.001	0.001
3-Ethylnonane	I10	0.0004	0.0033	0.000	0.000
1,3-Diethylbenzene	A10	0.0002	0.0014	0.000	0.000
1,3-Methyl-n-propylbenzene	A10	0.0016	0.0112	0.001	0.001
1,4-Methyl-n-propylbenzene	A10	0.0002	0.0014	0.000	0.000
n-Butylbenzene	A10	0.0010	0.0070	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0003	0.0021	0.000	0.000
1,2-Diethylbenzene	A10	0.0004	0.0028	0.000	0.000
t-Decahydronaphthalene	A9	0.0002	0.0016	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0008	0.0056	0.001	0.001

1,3-Dimethyl-4-ethylbenzene	A10	0.0015	0.0105	0.001	0.001
1,2-Dimethyl-4-ethylbenzene	A10	0.0004	0.0028	0.000	0.000
1,3-Dimethyl-2-ethylbenzene	A10	0.0001	0.0007	0.000	0.000
1,2-Dimethyl-3-ethylbenzene	A10	0.0001	0.0007	0.000	0.000
1,2-Ethyl-i-propylbenzene	A10	0.0002	0.0016	0.000	0.000
1,4-Methyl-t-butylbenzene	A11	0.0002	0.0016	0.000	0.000
UnknownC10s	U10	0.0208	0.1541	0.013	0.013
n-Undecane	P11	0.0039	0.0318	0.003	0.003
1,4-Ethyl-i-propylbenzene	A11	0.0001	0.0008	0.000	0.000
1,2,4,5-Tetramethylbenzene	A11	0.0002	0.0014	0.000	0.000
1,2-Methyl-n-butylbenzene	A11	0.0001	0.0008	0.000	0.000
1,2,3,5-Tetramethylbenzene	A11	0.0001	0.0007	0.000	0.000
1,2-Di-n-propylbenzene	A11	0.0001	0.0008	0.000	0.000
1,4-Di-i-propylbenzene	A11	0.0001	0.0008	0.000	0.000
Tetrahydronaphthalene	A10	0.0001	0.0007	0.000	0.000
UnknownC11s	U11	0.0107	0.0871	0.007	0.007
n-Dodecane	P12	0.0002	0.0018	0.000	0.000
1,3,5-Triethylbenzene	A12	0.0001	0.0008	0.000	0.000
UnknownC12s	U12	0.0012	0.0098	0.001	0.001
TOTAL		100.00000	100.00000	2.3468	2.3582

BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0382	0.1554
TOLUENE	0.0734	0.3521
ETHYLBENZENE	0.0060	0.0332
XYLENES	0.0945	0.5223
TOTAL BTEX	0.2121	1.0630

**(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 :	1018.1 /scf	1023.6	/scf
	NET WET REAL :	1000.3 /scf	1005.8	/scf
HHV	GROSS DRY REAL :	1124.6 /scf	1130.7	/scf
	GROSS WET REAL :	1104.9 /scf	1111.0	/scf
	NET HEATING VALUE (60 °F ideal reaction):		20161.8	Btu/lbm
	GROSS HEATING VALUE (60°F ideal reaction):		22284.5	Btu/lbm
	RELATIVE DENSITY (AIR=1):		0.6621	
	DENSITY		0.05060	lb/scf
	COMPRESSIBILITY FACTOR :		0.9976	
	REGULAR WOBBE INDEX		1383.1	

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

Net Dry Ideal BTU	5390.9 /scf	Relative Density - SG (Air=1)	3.7364		***C6+factors***
Gross Dry Ideal BTU	5784.9 /scf	**Z** Compressiblity Factor	0.99641		0.99562
Net Dry Ideal BTU	19134 /lb	Density Factor	285.18	lbm/1000 ft3	
Gross Dry Ideal BTU	20537.1 /lb	Molar Mass or MW	108.228	g/mol	
		Volume Liquid Ideal gas	0.692	scf/gal	21.9

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.