EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PRIMARY DB KEY:	**05-103-10735**	NAME/DESCRIP :	**125190121 PCU 297-10A9**
LEASE #:	**COD52141**		**PRODUCTION CASING**
FIELD/AREA:	**PICEANCE CREEK**		

PROJECT NO. :	**202603079**	ANALYSIS NO. :	**02**
COMPANY NAME :	**QB ENERGY OPERATING, LLC**	ANALYSIS DATE:	APRIL 10, 2026 14:49
OFFICE / BRANCH:	PARACHUTE, CO	SAMPLE DATE :	MARCH 5, 2026
CUSTOMER REF:		TO:	
PRODUCER :	QB ENERGY OPERATING LLC	EFFECTIVE DATE:	

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

SAMPLE CYCLE:			SAMPLE TYPE:	SPOT
SAMPLE PRES. :	1484	psig	PROBE :	NO
FLOW PRES. :		psig	CYLINDER NO. :	ECA-795
LAB PRES:		psig	SAMPLED BY :	MIKE KELLEY
SAMPLE TEMP. :	40	ºf	SAMPLING COMPANY:	QB ENERGY
AMBIENT TEMP.:		ºf	H2S BY STAIN TUBE:	ppm mol
H2O BY STAIN TUBE:		#/mmcf	CO2 BY STAIN TUBE:	Mol %

FIELD COMMENTS:

LAB COMMENTS: ***Possible moisture observed.***

COMPONENT	MOLE %	MASS %	GPM @ 14.65	GPM @ 14.73
HELIUM	0.00	0.00	---	---
HYDROGEN	0.00	0.00	---	---
OXYGEN/ARGON	0.00	0.00	---	---
NITROGEN	0.08	0.12	---	---
CARBON DIOXIDE	5.14	12.43	---	---
METHANE	92.3769	81.3993	---	---
ETHANE	1.6788	2.7727	0.4476	0.4500
PROPANE	0.2404	0.5823	0.0659	0.0663
I-BUTANE	0.0751	0.2398	0.0250	0.0251
N-BUTANE	0.0401	0.1280	0.0130	0.0131
I-PENTANE	0.0289	0.1144	0.0100	0.0100
N-PENTANE	0.0160	0.0634	0.0060	0.0060
HEXANES PLUS	0.3238	2.1501	0.1510	0.1512
TOTALS	100.00000	100.00000	0.7185	0.7217

BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0086	0.0369
TOLUENE	0.0191	0.0967
ETHYLBENZENE	0.0016	0.0093
XYLENES	0.0224	0.1306
TOTAL BTEX	0.0517	0.2735

CALCULATED VALUES**

	BTU @	14.65	14.73	
LHV	NET DRY REAL :	895.6 /scf	900.5	/scf
	NET WET REAL :	879.9 /scf	884.8	/scf
HHV	GROSS DRY REAL :	992.9 /scf	998.4	/scf
	GROSS WET REAL :	975.5 /scf	981.0	/scf
	NET HEATING VALUE (60 °F ideal reaction):		18707.2	Btu/lbm
	GROSS HEATING VALUE (60°F ideal reaction):		20747.0	Btu/lbm
	RELATIVE DENSITY (AIR=1):		0.6271	
	DENSITY		0.04797	lbm/scf
	COMPRESSIBILITY FACTOR :		0.9978	
	REGULAR WOBBE INDEX		1255.1	

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

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

EMPACT

EXTENDED NATURAL GAS ANALYSIS (*DHA)

GLYCALC INFORMATION

PROJECT NO. :	202603079	ANALYSIS NO. :	02
COMPANY NAME :	QB ENERGY OPERATING, LLC	ANALYSIS DATE:	APRIL 10, 2026 14:49
ACCOUNT NO. :		SAMPLE DATE :	MARCH 5, 2026
PRODUCER :	QB ENERGY OPERATING LLC	CYLINDER NO. :	ECA-795
LEASE NO. :	COD52141	SAMPLED BY :	MIKE KELLEY
NAME/DESCRIP :	125190121 PCU 297-10A9 PRODUCTION CASING		
FIELD DATA		SAMPLE TEMP. :	40
SAMPLE PRES. :	1484	AMBIENT TEMP.:	
H2S BY STAIN TUBE:	_ *ppm mol*		
COMMENTS :	***SPOT*** ***NO PROBE*** ***Possible moisture observed.***		

Componet	Mole %	Wt %
Helium	0.00	0.00
Hydrogen	0.00	0.00
Carbon Dioxide	5.14	12.43
Nitrogen	0.08	0.12
Methane	92.3769	81.3993
Ethane	1.6788	2.7727
Propane	0.2404	0.5823
Isobutane	0.0751	0.2398
n-Butane	0.0401	0.1280
Isopentane	0.0280	0.1109
n-Pentane	0.0160	0.0634
Cyclopentane	0.0009	0.0035
n-Hexane	0.0102	0.0483
Cyclohexane	0.0068	0.0314
Other Hexanes	0.0235	0.1107
Heptanes	0.0253	0.1384
Methylcyclohexane	0.0171	0.0922
2,2,4 Trimethylpentane	0.0001	0.0006
Benzene	0.0086	0.0369
Toluene	0.0191	0.0967
Ethylbenzene	0.0016	0.0093
Xylenes	0.0224	0.1306
C8+ Heavies	0.1891	1.4550
Subtotal	100.00000	100.00000
Oxygen/Argon	0.00	0.00
Total	100.00000	100.00000

Calculated Values BTU @ **14.65**		**Total Sample**	**C6+ Fraction**	**C8+ Fraction**	**C10+ Fraction**	
LHV	Net Dry Real:	895.6	6029.2	6804.1	7672.0	**Btu/scf**
	Net Wet Real:	879.9	5923.8	6685.2	7537.9	**Btu/scf**
HHV	Gross Dry Real:	992.9	6484.0	7330.8	8297.6	**Btu/scf**
	Gross Wet Real:	975.5	6370.7	7202.6	8152.5	**Btu/scf**
Other Calculated Values						
Regualr Wobbe Index*		1255.1	3176.3	3387.9	3629.0	**Btu/scf**
Net Heating Value (60 °F ideal reaction):		18707.2	18981.0	19019.4	19031.0	**Btu/lbm**
Gross Heating Value (60°F ideal reaction):		20747.0	20409.1	20483.9	20573.5	**Btu/lbm**
Molar Mass (MW):		18.20588	120.944	136.293	152.3	**g/mol**
Relative Density (AIR=1):		0.6271	4.1756	4.7050	5.2585	**SG**
Density:		0.04797	0.31870	0.35915	0.40133	**lbm/scf**
Compressiblity Factor:		0.9978	0.9979	0.9993	0.9998	**Z**
Liquid Volume real gas @:	**14.65**	17.1701	0.1505	0.1107	0.0668	**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.

EMPACT Analytical Systems Inc. 365 S Main St Brighton, CO 80601 303-637-0150

EXTENDED NATURAL GAS ANALYSIS (*DHA)

DHA COMPONENT LIST

PRIMARY DB KEY:	**05-103-10735**	NAME/DESCRIP :	**125190121 PCU 297-10A9**
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FIELD/AREA:	**PICEANCE CREEK**		
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PRODUCER :	QB ENERGY OPERATING LLC	EFFECTIVE DATE:	
FIELD DATA			
SAMPLE CYCLE:		SAMPLE TYPE:	SPOT
SAMPLE PRES. :	1484 psig	PROBE :	NO
FLOW PRES. :	psig	CYLINDER NO. :	ECA-795
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SAMPLE TEMP. :	40 ºf	SAMPLING COMPANY:	QB ENERGY
AMBIENT TEMP.:	ºf	H2S BY STAIN TUBE:	_ ppm mol
H2O BY STAIN TUBE:	- #/mmcf	CO2 BY STAIN TUBE:	- Mol %
FIELD COMMENTS:			
LAB COMMENTS:	***Possible moisture observed.***		

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ **14.65**	GPM @ **14.73**
Nitrogen	---	0.08	0.12	---	---
Carbon Dioxide	---	5.14	12.43	---	---
Methane	P1	92.3769	81.3993	---	---
Ethane	P2	1.6788	2.7727	0.448	0.450
Propane	P3	0.2404	0.5823	0.066	0.066
i-Butane	I4	0.0751	0.2398	0.025	0.025
n-Butane	P4	0.0401	0.1280	0.013	0.013
2,2-Dimethylpropane	I5	0.0020	0.0079	0.001	0.001
i-Pentane	I5	0.0260	0.1030	0.009	0.009
n-Pentane	P5	0.0160	0.0634	0.006	0.006
2,2-Dimethylbutane	I6	0.0025	0.0118	0.001	0.001
Cyclopentane	N5	0.0009	0.0035	0.000	0.000
2,3-Dimethylbutane	I6	0.0022	0.0104	0.001	0.001
2-Methylpentane	I6	0.0087	0.0412	0.004	0.004
3-Methylpentane	I6	0.0051	0.0242	0.002	0.002
n-Hexane	P6	0.0102	0.0483	0.004	0.004
2,2-Dimethylpentane	I7	0.0007	0.0038	0.000	0.000
Methylcyclopentane	N6	0.0050	0.0231	0.002	0.002
2,4-Dimethylpentane	I7	0.0008	0.0044	0.000	0.000
2,2,3-Trimethylbutane	I7	0.0003	0.0016	0.000	0.000
Benzene	A6	0.0086	0.0369	0.002	0.002
3,3-Dimethylpentane	I7	0.0005	0.0027	0.000	0.000
Cyclohexane	N6	0.0068	0.0314	0.002	0.002
2-Methylhexane	I7	0.0041	0.0226	0.002	0.002
2,3-Dimethylpentane	I7	0.0011	0.0060	0.000	0.000
1,1-Dimethylcyclopentane	N7	0.0008	0.0043	0.000	0.000
3-Methylhexane	I7	0.0039	0.0215	0.002	0.002

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1c,3-Dimethylcyclopentane	N7	0.0010	0.0054	0.000	0.000
1t,3-Dimethylcyclopentane	N7	0.0009	0.0048	0.000	0.000
3-Ethylpentane	I7	0.0002	0.0011	0.000	0.000
1t,2-Dimethylcyclopentane	N7	0.0015	0.0081	0.001	0.001
2,2,4-Trimethylpentane	I8	0.0001	0.0006	0.000	0.000
n-Heptane	P7	0.0084	0.0462	0.004	0.004
1c,2-Dimethylcyclopentane	N7	0.0004	0.0021	0.000	0.000
Methylcyclohexane	N7	0.0171	0.0922	0.007	0.007
2,2-Dimethylhexane	I8	0.0005	0.0031	0.000	0.000
1,1,3-Trimethylcyclopentane	N7	0.0001	0.0006	0.000	0.000
Ethylcyclopentane	N7	0.0006	0.0032	0.000	0.000
2,5-Dimethylhexane	I8	0.0007	0.0044	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0006	0.0038	0.000	0.000
2,4-Dimethylhexane	I8	0.0001	0.0006	0.000	0.000
1c,2t,4-Trimethylcyclopentane	N8	0.0003	0.0019	0.000	0.000
3,3-Dimethylhexane	I8	0.0002	0.0013	0.000	0.000
Toluene	A7	0.0191	0.0967	0.006	0.006
2,3-Dimethylhexane	I8	0.0005	0.0031	0.000	0.000
2-Methyl-3-ethylpentane	I8	0.0001	0.0006	0.000	0.000
2-Methylheptane	I8	0.0033	0.0207	0.002	0.002
4-Methylheptane	I8	0.0011	0.0069	0.001	0.001
3-Methyl-3-ethylpentane	I8	0.0001	0.0006	0.000	0.000
3,4-Dimethylhexane	I8	0.0001	0.0006	0.000	0.000
3-Methylheptane	I8	0.0028	0.0176	0.001	0.001
1c,2t,3-Trimethylcyclopentane	N8	0.0035	0.0216	0.002	0.002
3-Ethylhexane	I8	0.0003	0.0019	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0015	0.0092	0.001	0.001
1,1-Dimethylcyclohexane	N8	0.0006	0.0037	0.000	0.000
2,2,5-Trimethylhexane	I9	0.0001	0.0007	0.000	0.000
3c-Ethylmethylcyclopentane	N8	0.0001	0.0006	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0001	0.0006	0.000	0.000
2t-Ethylmethylcyclopentane	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	0.0012	0.0074	0.001	0.001
1t,3-Dimethylcyclohexane	N8	0.0005	0.0031	0.000	0.000
n-Octane	P8	0.0087	0.0546	0.004	0.004
1c,4-Dimethylcyclohexane	N8	0.0010	0.0061	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-Dimethylheptane	I9	0.0005	0.0035	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0015	0.0104	0.001	0.001
2,2,3-Trimethylhexane	I9	0.0005	0.0035	0.000	0.000
2,4-Dimethylheptane	I9	0.0001	0.0007	0.000	0.000
Ethylcyclohexane	N8	0.0013	0.0080	0.001	0.001
n-Propylcyclopentane	N8	0.0006	0.0037	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0004	0.0027	0.000	0.000
2,5-Dimethylheptane	I9	0.0012	0.0085	0.001	0.001
3,3-Dimethylheptane	I9	0.0003	0.0021	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0001	0.0007	0.000	0.000
Ethylbenzene	I8	0.0016	0.0093	0.001	0.001
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	0.0151	0.0880	0.006	0.006
1,4-Dimethylbenzene (p-Xylene)	A8	0.0050	0.0292	0.002	0.002
3,4-Dimethylheptane	I9	0.0002	0.0014	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0001	0.0007	0.000	0.000
4-Ethylheptane	I9	0.0003	0.0021	0.000	0.000
4-Methyloctane	I9	0.0016	0.0113	0.001	0.001
2-Methyloctane	I9	0.0024	0.0169	0.001	0.001
1c,2t,3-Trimethylcyclohexane	N9	0.0001	0.0007	0.000	0.000

3-Ethylheptane	I9	0.0001	0.0007	0.000	0.000
3-Methyloctane	I9	0.0003	0.0021	0.000	0.000
1c,2t,4c-Trimethylcyclohexane	I9	0.0023	0.0159	0.001	0.001
1,1,2-Trimethylcyclohexane	N9	0.0001	0.0007	0.000	0.000
3,3-Diethylpentane	I9	0.0001	0.0007	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0023	0.0134	0.001	0.001
i-Butylcyclopentane	N9	0.0009	0.0063	0.000	0.000
UnknownC8s	U8	0.0002	0.0013	0.000	0.000
n-Nonane	P9	0.0117	0.0824	0.007	0.007
1,1-Methylethylcyclohexane	N9	0.0006	0.0042	0.000	0.000
i-Propylbenzene	A9	0.0002	0.0013	0.000	0.000
i-Propylcyclohexane	N9	0.0002	0.0014	0.000	0.000
2,2-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
2,4-Dimethyloctane	I10	0.0005	0.0039	0.000	0.000
2,6-Dimethyloctane	I10	0.0002	0.0015	0.000	0.000
2,5-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
n-Butylcyclopentane	N9	0.0008	0.0055	0.000	0.000
3,3-Dimethyloctane	I10	0.0007	0.0055	0.000	0.000
n-Propylbenzene	A9	0.0015	0.0099	0.001	0.001
3,6-Dimethyloctane	I10	0.0004	0.0031	0.000	0.000
3-Methyl-5-ethylheptane	I10	0.0001	0.0008	0.000	0.000
1,3-Methylethylbenzene	A9	0.0019	0.0125	0.001	0.001
1,4-Methylethylbenzene	A9	0.0007	0.0046	0.000	0.000
1,3,5-Trimethylbenzene	A9	0.0046	0.0304	0.002	0.002
2,3-Dimethyloctane	I10	0.0007	0.0055	0.000	0.000
5-Methylnonane	I10	0.0016	0.0125	0.001	0.001
1,2-Methylethylbenzene	A9	0.0021	0.0138	0.001	0.001
2-Methylnonane	I10	0.0005	0.0039	0.000	0.000
3-Ethyloctane	I10	0.0003	0.0024	0.000	0.000
3-Methylnonane	I10	0.0018	0.0141	0.001	0.001
t-Butylbenzene	A10	0.0038	0.0280	0.002	0.002
i-Butylcyclohexane	N10	0.0005	0.0038	0.000	0.000
1t-Methyl-2-n-propylcyclohexane	I10	0.0002	0.0015	0.000	0.000
i-Butylbenzene	A10	0.0001	0.0007	0.000	0.000
sec-Butylbenzene	A10	0.0001	0.0007	0.000	0.000
UnknownC9s	U9	0.0046	0.0324	0.003	0.003
n-Decane	P10	0.0141	0.1102	0.009	0.009
1,2,3-Trimethylbenzene	A9	0.0001	0.0007	0.000	0.000
1,3-Methyl-i-propylbenzene	A10	0.0005	0.0037	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0004	0.0030	0.000	0.000
Sec-Butylcyclohexane	A10	0.0001	0.0008	0.000	0.000
1,2-Methyl-i-propylbenzene	A10	0.0002	0.0015	0.000	0.000
3-Ethylnonane	I10	0.0006	0.0052	0.000	0.000
1,3-Diethylbenzene	A10	0.0018	0.0133	0.001	0.001
1,3-Methyl-n-propylbenzene	A10	0.0008	0.0059	0.000	0.000
1,4-Diethylbenzene	A10	0.0008	0.0059	0.000	0.000
1,4-Methyl-n-propylbenzene	A10	0.0002	0.0015	0.000	0.000
n-Butylbenzene	A10	0.0010	0.0074	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0003	0.0022	0.000	0.000
1,2-Diethylbenzene	A10	0.0009	0.0066	0.000	0.000
t-Decahydronaphthalene	A9	0.0003	0.0025	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0011	0.0081	0.001	0.001
1,3-Dimethyl-4-ethylbenzene	A10	0.0024	0.0177	0.002	0.002
1,2-Dimethyl-4-ethylbenzene	A10	0.0006	0.0044	0.000	0.000
1,3-Dimethyl-2-ethylbenzene	A10	0.0016	0.0118	0.001	0.001
1,2-Dimethyl-3-ethylbenzene	A10	0.0003	0.0022	0.000	0.000
1,2-Ethyl-i-propylbenzene	A10	0.0004	0.0032	0.000	0.000
1,4-Methyl-t-butylbenzene	A11	0.0004	0.0032	0.000	0.000
UnknownC10s	U10	0.0038	0.0297	0.002	0.002
n-Undecane	P11	0.0179	0.1537	0.012	0.012
1,4-Ethyl-i-propylbenzene	A11	0.0004	0.0032	0.000	0.000

1,2,4,5-Tetramethylbenzene	A11	0.0004	0.0030	0.000	0.000
1,2-Methyl-n-butylbenzene	A11	0.0003	0.0024	0.000	0.000
1,2,3,5-Tetramethylbenzene	A11	0.0006	0.0044	0.000	0.000
1,2-Methyl-t-butylbenzene	A11	0.0005	0.0041	0.000	0.000
5-Methylindan	A11	0.0001	0.0007	0.000	0.000
4-Methylindan	A11	0.0008	0.0058	0.001	0.001
1,2-Ethyl-n-propylbenzene	A11	0.0008	0.0065	0.001	0.001
2-Methylindan	A11	0.0001	0.0007	0.000	0.000
1,3-Methyl-n-butylbenzene	A11	0.0001	0.0008	0.000	0.000
1,3-Di-i-propylbenzene	A11	0.0003	0.0027	0.000	0.000
sec-Pentylbenzene	A11	0.0009	0.0073	0.000	0.000
n-Pentylbenzene	A11	0.0012	0.0098	0.001	0.001
1,2-Di-n-propylbenzene	A11	0.0011	0.0098	0.001	0.001
1,4-Di-i-propylbenzene	A11	0.0004	0.0036	0.000	0.000
Tetrahydronaphthalene	A10	0.0005	0.0036	0.000	0.000
Naphthalene	A10	0.0021	0.0148	0.002	0.002
1-t-Butyl-3,5-dimethylbenzene	A12	0.0003	0.0027	0.000	0.000
1,4-Ethyl-t-butylbenzene	A11	0.0003	0.0027	0.000	0.000
1,3-Di-n-propylbenzene	A12	0.0006	0.0053	0.000	0.000
UnknownC11s	U11	0.0056	0.0481	0.004	0.004
n-Dodecane	P12	0.0259	0.2423	0.019	0.019
1,3,5-Triethylbenzene	A12	0.0036	0.0321	0.002	0.002
1,2,4-Triethylbenzene	A12	0.0003	0.0027	0.000	0.000
n-Hexylbenzene	A12	0.0001	0.0009	0.000	0.000
1,2,3,4,5-Pentamethylbenzene	A13	0.0001	0.0008	0.000	0.000
UnknownC12s	U12	0.0060	0.0515	0.004	0.004
n-Tridecane	P13	0.0002	0.0020	0.000	0.000
n-Tetradecane	P14	0.0001	0.0011	0.000	0.000
UnknownC14s	U14	0.0002	0.0022	0.000	0.000
n-Pentadecane	P15	0.0001	0.0011	0.000	0.000
TOTAL		100.00000	100.00000	0.7185	0.7217

BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0086	0.0369
TOLUENE	0.0191	0.0967
ETHYLBENZENE	0.0016	0.0093
XYLENES	0.0224	0.1306
TOTAL BTEX	0.0517	0.2735

*(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 :	895.6 /scf	900.5 /scf	
	NET WET REAL :	879.9 /scf	884.8 /scf	
HHV	GROSS DRY REAL :	992.9 /scf	998.4 /scf	
	GROSS WET REAL :	975.5 /scf	981.0 /scf	
	NET HEATING VALUE (60 °F ideal reaction):		18707.2	Btu/lbm
	GROSS HEATING VALUE (60°F ideal reaction):		20747.0	Btu/lbm
	RELATIVE DENSITY (AIR=1):		0.6271	
	DENSITY		0.04797	lb/scf
	COMPRESSIBILITY FACTOR :		0.9978	
	REGULAR WOBBE INDEX		1255.1	

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

Net Dry Ideal BTU	6035.3 /scf	Relative Density - SG (Air=1)	4.1756	***C6+factors***
Gross Dry Ideal BTU	6490.6 /scf	**Z** Compressiblity Factor	0.99788	0.99682
Net Dry Ideal BTU	18981 /lb	Density Factor	318.695 lbm/1000 ft3	
Gross Dry Ideal BTU	20409.1 /lb	Molar Mass or MW	120.944 g/mol	
		Volume Liquid Ideal gas	0.151 scf/gal	19.7

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.