EMPACT

EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PRIMARY DB KEY:	**05-103-10782**	NAME/DESCRIP :	**FEDERAL 4S-95-1-21DP INTERMEDIATE CASING**
LEASE #:			
FIELD/AREA:			
PROJECT NO. :	**202501047**	ANALYSIS NO. :	**01**
COMPANY NAME :	**QB ENERGY OPERATING, LLC**	ANALYSIS DATE:	JANUARY 27, 2025 09:40
OFFICE / BRANCH:	PARACHUTE	SAMPLE DATE :	DECEMBER 16, 2024 15:00
CUSTOMER REF:		TO:	
PRODUCER :	QB ENERGY OPERATING, LLC	EFFECTIVE DATE:	
FIELD DATA			
SAMPLE CYCLE:		SAMPLE TYPE:	SPOT
SAMPLE PRES. :	1647 psig	PROBE :	
FLOW PRES. :	psig	CYLINDER NO. :	TBI-567
LAB PRES:	psig	SAMPLED BY :	ROWDY HAMILTON
SAMPLE TEMP. :	12 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:	***INT CSG Sample 2***		
LAB COMMENTS:			

COMPONENT	MOLE %	MASS %	GPM @ 14.65	GPM @ 14.73
ALCOHOLS	0.0002	0.0004	0.0000	0.0000
HELIUM	0.00	0.00	---	---
HYDROGEN	0.22	0.03	---	---
OXYGEN/ARGON	0.00	0.00	---	---
NITROGEN	0.10	0.16	---	---
CARBON DIOXIDE	0.02	0.05	---	---
METHANE	96.9947	91.3898	---	---
ETHANE	1.7703	3.1263	0.4715	0.4741
PROPANE	0.2223	0.5757	0.0609	0.0613
I-BUTANE	0.0711	0.2427	0.0230	0.0231
N-BUTANE	0.0402	0.1373	0.0130	0.0131
I-PENTANE	0.0264	0.1117	0.0090	0.0090
N-PENTANE	0.0081	0.0343	0.0030	0.0030
HEXANES PLUS	0.5267	4.1418	0.2870	0.2875
TOTALS	100.00000	100.00000	0.8674	0.8711

BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0182	0.0835
TOLUENE	0.0347	0.1878
ETHYLBENZENE	0.0029	0.0181
XYLENES	0.0283	0.1766
TOTAL BTEX	0.0841	0.4660

*(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 :	954.5 /scf	959.7 /scf
	NET WET REAL :	937.8 /scf	943.0 /scf
HHV	GROSS DRY REAL :	1058.1 /scf	1063.9 /scf
	GROSS WET REAL :	1039.6 /scf	1045.4 /scf
	NET HEATING VALUE (60 °F ideal reaction):		21297.6 Btu/lbm
	GROSS HEATING VALUE (60°F ideal reaction):		23613.1 Btu/lbm
	RELATIVE DENSITY (AIR=1):		0.5870
	DENSITY		0.04486 lbm/scf
	COMPRESSIBILITY FACTOR :		0.9979
	REGULAR WOBBE INDEX		1382.5

EMPACT

EXTENDED NATURAL GAS ANALYSIS (*DHA)

GLYCALC INFORMATION

PROJECT NO. :	202501047	ANALYSIS NO. :	01
COMPANY NAME :	QB ENERGY OPERATING, LLC	ANALYSIS DATE:	JANUARY 27, 2025 09:40
ACCOUNT NO. :		SAMPLE DATE :	DECEMBER 16, 2024 15:00
PRODUCER :	QB ENERGY OPERATING, LLC	CYLINDER NO. :	TBI-567
LEASE NO. :		SAMPLED BY :	ROWDY HAMILTON
NAME/DESCRIP :	FEDERAL 4S-95-1-21DP INTERMEDIATE CASING		
FIELD DATA		SAMPLE TEMP. :	12
SAMPLE PRES. :	1647	AMBIENT TEMP.:	
H2S BY STAIN TUBE:	_ *ppm mol*		
COMMENTS :	***SPOT*** ***INT CSG Sample 2***		

Componet	Mole %	Wt %
Helium	0.00	0.00
Hydrogen	0.22	0.03
Carbon Dioxide	0.02	0.05
Nitrogen	0.10	0.16
Methane	96.9947	91.3898
Ethane	1.7703	3.1263
Propane	0.2223	0.5757
Isobutane	0.0711	0.2427
n-Butane	0.0402	0.1373
Isopentane	0.0255	0.1080
n-Pentane	0.0081	0.0343
Cyclopentane	0.0009	0.0037
n-Hexane	0.0060	0.0304
Cyclohexane	0.0078	0.0385
Other Hexanes	0.0218	0.1097
Heptanes	0.0234	0.1372
Methylcyclohexane	0.0224	0.1292
2,2,4 Trimethylpentane	0.0001	0.0007
Benzene	0.0182	0.0835
Toluene	0.0347	0.1878
Ethylbenzene	0.0029	0.0181
Xylenes	0.0283	0.1766
C8+ Heavies	0.3611	3.2301
Subtotal	99.99980	99.99960
Oxygen/Argon	0.00	0.00
Alcohols	0.0002	0.0004
Total	100.00000	100.00000

Calculated Values BTU @ **14.65**		**Total Sample**	**C6+ Fraction**	**C8+ Fraction**	**C10+ Fraction**	
LHV	Net Dry Real:	954.5	6673.0	7437.9	8130.3	**Btu/scf**
	Net Wet Real:	937.8	6556.3	7307.9	7988.2	**Btu/scf**
HHV	Gross Dry Real:	1058.1	7187.6	8029.4	8800.2	**Btu/scf**
	Gross Wet Real:	1039.6	7062.0	7889.0	8646.4	**Btu/scf**
Other Calculated Values						
Regualr Wobbe Index*		1382.5	3350.4	3554.7	3742.5	**Btu/scf**
Net Heating Value (60 °F ideal reaction):		21297.6	18506.7	18482.7	18431.4	**Btu/lbm**
Gross Heating Value (60°F ideal reaction):		23613.1	19929.6	19948.8	19946.1	**Btu/lbm**
Molar Mass (MW):		17.02648	133.833	148.581	161.112	**g/mol**
Relative Density (AIR=1):		0.5870	4.6204	5.1306	5.5626	**SG**
Density:		0.04486	0.35269	0.39154	0.42455	**lbm/scf**
Compressiblity Factor:		0.9979	0.9988	0.9996	0.9999	**Z**
Liquid Volume real gas @:	**14.65**	17.2488	0.2861	0.2363	0.1894	**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

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FIELD COMMENTS:	***INT CSG Sample 2***		
LAB COMMENTS:			

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ **14.65**	GPM @ **14.73**
Hydrogen	---	0.22	0.03	---	---
Nitrogen	---	0.10	0.16	---	---
Carbon Dioxide	---	0.02	0.05	---	---
Methane	P1	96.9947	91.3898	---	---
Ethane	P2	1.7703	3.1263	0.472	0.474
Propane	P3	0.2223	0.5757	0.061	0.061
i-Butane	I4	0.0711	0.2427	0.023	0.023
Methanol	X1	0.0002	0.0004	0.000	0.000
n-Butane	P4	0.0401	0.1369	0.013	0.013
2,2-Dimethylpropane	I5	0.0025	0.0106	0.001	0.001
i-Pentane	I5	0.0230	0.0974	0.008	0.008
UnknownC4s	U4	0.0001	0.0004	0.000	0.000
n-Pentane	P5	0.0080	0.0339	0.003	0.003
2,2-Dimethylbutane	I6	0.0031	0.0157	0.001	0.001
Cyclopentane	N5	0.0009	0.0037	0.000	0.000
2,3-Dimethylbutane	I6	0.0020	0.0101	0.001	0.001
2-Methylpentane	I6	0.0068	0.0344	0.003	0.003
3-Methylpentane	I6	0.0042	0.0213	0.002	0.002
UnknownC5s	U5	0.0001	0.0004	0.000	0.000
n-Hexane	P6	0.0060	0.0304	0.002	0.002
2,2-Dimethylpentane	I7	0.0006	0.0035	0.000	0.000
Methylcyclopentane	N6	0.0056	0.0277	0.002	0.002
2,4-Dimethylpentane	I7	0.0008	0.0047	0.000	0.000
2,2,3-Trimethylbutane	I7	0.0003	0.0018	0.000	0.000
Benzene	A6	0.0182	0.0835	0.005	0.005
3,3-Dimethylpentane	I7	0.0006	0.0035	0.000	0.000
Cyclohexane	N6	0.0078	0.0385	0.003	0.003
2-Methylhexane	I7	0.0039	0.0230	0.002	0.002

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2,3-Dimethylpentane	I7	0.0011	0.0065	0.000	0.000
1,1-Dimethylcyclopentane	N7	0.0009	0.0052	0.000	0.000
3-Methylhexane	I7	0.0037	0.0218	0.002	0.002
1c,3-Dimethylcyclopentane	N7	0.0012	0.0069	0.001	0.001
1t,3-Dimethylcyclopentane	N7	0.0011	0.0063	0.001	0.001
3-Ethylpentane	I7	0.0002	0.0012	0.000	0.000
1t,2-Dimethylcyclopentane	N7	0.0018	0.0104	0.001	0.001
2,2,4-Trimethylpentane	I8	0.0001	0.0007	0.000	0.000
UnknownC6s	U6	0.0001	0.0005	0.000	0.000
n-Heptane	P7	0.0059	0.0347	0.003	0.003
1c,2-Dimethylcyclopentane	N7	0.0003	0.0017	0.000	0.000
Methylcyclohexane	N7	0.0224	0.1292	0.009	0.009
2,2-Dimethylhexane	I8	0.0006	0.0041	0.000	0.000
1,1,3-Trimethylcyclopentane	N7	0.0002	0.0013	0.000	0.000
Ethylcyclopentane	N7	0.0007	0.0041	0.000	0.000
2,5-Dimethylhexane	I8	0.0008	0.0053	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0007	0.0047	0.000	0.000
2,4-Dimethylhexane	I8	0.0001	0.0007	0.000	0.000
1c,2t,4-Trimethylcyclopentane	N8	0.0004	0.0026	0.000	0.000
3,3-Dimethylhexane	I8	0.0003	0.0020	0.000	0.000
2,3,4-Trimethylpentane	I8	0.0001	0.0007	0.000	0.000
2,3,3-Trimethylpentane	I8	0.0001	0.0007	0.000	0.000
Toluene	A7	0.0347	0.1878	0.012	0.012
2,3-Dimethylhexane	I8	0.0006	0.0041	0.000	0.000
2-Methyl-3-ethylpentane	I8	0.0001	0.0007	0.000	0.000
2-Methylheptane	I8	0.0031	0.0208	0.002	0.002
4-Methylheptane	I8	0.0010	0.0067	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
3-Methylheptane	I8	0.0029	0.0194	0.001	0.001
1c,2t,3-Trimethylcyclopentane	N8	0.0043	0.0284	0.002	0.002
3-Ethylhexane	I8	0.0002	0.0014	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0019	0.0125	0.001	0.001
1,1-Dimethylcyclohexane	N8	0.0008	0.0053	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	0.0001	0.0007	0.000	0.000
2t-Ethylmethylcyclopentane	N8	0.0001	0.0007	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0014	0.0092	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	0.0069	0.0463	0.004	0.004
1c,4-Dimethylcyclohexane	N8	0.0011	0.0072	0.001	0.001
i-Propylcyclopentane	I8	0.0001	0.0007	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0002	0.0015	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0001	0.0008	0.000	0.000
2,2-Dimethylheptane	I9	0.0004	0.0030	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0016	0.0119	0.001	0.001
2,2,3-Trimethylhexane	I9	0.0003	0.0022	0.000	0.000
2,4-Dimethylheptane	I9	0.0002	0.0015	0.000	0.000
Ethylcyclohexane	N8	0.0010	0.0066	0.000	0.000
n-Propylcyclopentane	N8	0.0006	0.0039	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0005	0.0037	0.000	0.000
2,5-Dimethylheptane	I9	0.0014	0.0106	0.001	0.001
3,3-Dimethylheptane	I9	0.0004	0.0030	0.000	0.000
3,5-Dimethylheptane	I9	0.0002	0.0015	0.000	0.000
2,6-Dimethylheptane	I9	0.0002	0.0015	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0001	0.0008	0.000	0.000
Ethylbenzene	I8	0.0029	0.0181	0.001	0.001
2,3-Dimethylheptane	I9	0.0001	0.0008	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0182	0.1135	0.007	0.007

1,4-Dimethylbenzene (p-Xylene)	A8	0.0061	0.0381	0.002	0.002
3,4-Dimethylheptane	I9	0.0001	0.0008	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0001	0.0008	0.000	0.000
4-Ethylheptane	I9	0.0003	0.0022	0.000	0.000
4-Methyloctane	I9	0.0018	0.0136	0.001	0.001
2-Methyloctane	I9	0.0025	0.0189	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	0.0186	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	0.0040	0.0250	0.002	0.002
i-Butylcyclopentane	N9	0.0013	0.0096	0.001	0.001
UnknownC8s	U8	0.0001	0.0007	0.000	0.000
n-Nonane	P9	0.0103	0.0776	0.006	0.006
1,1-Methylethylcyclohexane	N9	0.0007	0.0052	0.000	0.000
i-Propylbenzene	A9	0.0005	0.0035	0.000	0.000
i-Propylcyclohexane	N9	0.0002	0.0015	0.000	0.000
2,2-Dimethyloctane	I10	0.0001	0.0008	0.000	0.000
2,4-Dimethyloctane	I10	0.0005	0.0042	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	0.0008	0.0059	0.000	0.000
3,3-Dimethyloctane	I10	0.0009	0.0075	0.001	0.001
n-Propylbenzene	A9	0.0017	0.0120	0.001	0.001
3,6-Dimethyloctane	I10	0.0013	0.0109	0.001	0.001
3-Methyl-5-ethylheptane	I10	0.0001	0.0008	0.000	0.000
1,3-Methylethylbenzene	A9	0.0053	0.0374	0.003	0.003
1,4-Methylethylbenzene	A9	0.0023	0.0162	0.001	0.001
1,3,5-Trimethylbenzene	A9	0.0052	0.0367	0.002	0.002
2,3-Dimethyloctane	I10	0.0009	0.0075	0.001	0.001
5-Methylnonane	I10	0.0020	0.0167	0.001	0.001
1,2-Methylethylbenzene	A9	0.0022	0.0155	0.001	0.001
2-Methylnonane	I10	0.0021	0.0176	0.001	0.001
3-Ethyloctane	I10	0.0003	0.0025	0.000	0.000
3-Methylnonane	I10	0.0026	0.0217	0.002	0.002
1,2,4-Trimethylbenzene	A9	0.0001	0.0007	0.000	0.000
t-Butylbenzene	A10	0.0093	0.0733	0.005	0.005
i-Butylcyclohexane	N10	0.0006	0.0049	0.000	0.000
1t-Methyl-2-n-propylcyclohexane	I10	0.0002	0.0016	0.000	0.000
i-Butylbenzene	A10	0.0003	0.0024	0.000	0.000
sec-Butylbenzene	A10	0.0002	0.0016	0.000	0.000
UnknownC9s	U9	0.0020	0.0151	0.001	0.001
n-Decane	P10	0.0167	0.1396	0.010	0.010
1,2,3-Trimethylbenzene	A9	0.0020	0.0141	0.001	0.001
1,3-Methyl-i-propylbenzene	A10	0.0002	0.0016	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0003	0.0024	0.000	0.000
Sec-Butylcyclohexane	A10	0.0006	0.0049	0.000	0.000
1,2-Methyl-i-propylbenzene	A10	0.0023	0.0182	0.001	0.001
3-Ethylnonane	I10	0.0017	0.0156	0.001	0.001
1,3-Diethylbenzene	A10	0.0004	0.0032	0.000	0.000
1,3-Methyl-n-propylbenzene	A10	0.0005	0.0039	0.000	0.000
1,4-Diethylbenzene	A10	0.0026	0.0205	0.001	0.001
1,4-Methyl-n-propylbenzene	A10	0.0011	0.0087	0.001	0.001
n-Butylbenzene	A10	0.0024	0.0189	0.001	0.001
1,3-Dimethyl-5-ethylbenzene	A10	0.0008	0.0063	0.001	0.001
1,2-Diethylbenzene	A10	0.0008	0.0063	0.000	0.000
t-Decahydronaphthalene	A9	0.0002	0.0018	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0021	0.0166	0.001	0.001
1,3-Dimethyl-4-ethylbenzene	A10	0.0048	0.0378	0.003	0.003
1,2-Dimethyl-4-ethylbenzene	A10	0.0015	0.0118	0.001	0.001

1,3-Dimethyl-2-ethylbenzene	A10	0.0032	0.0253	0.002	0.002
1,2-Dimethyl-3-ethylbenzene	A10	0.0007	0.0055	0.000	0.000
1,2-Ethyl-i-propylbenzene	A10	0.0020	0.0174	0.001	0.001
1,4-Methyl-t-butylbenzene	A11	0.0017	0.0148	0.001	0.001
UnknownC10s	U10	0.0061	0.0510	0.004	0.004
n-Undecane	P11	0.0278	0.2552	0.018	0.018
1,4-Ethyl-i-propylbenzene	A11	0.0015	0.0130	0.001	0.001
1,2,4,5-Tetramethylbenzene	A11	0.0019	0.0150	0.001	0.001
1,2-Methyl-n-butylbenzene	A11	0.0014	0.0122	0.001	0.001
1,2,3,5-Tetramethylbenzene	A11	0.0021	0.0166	0.001	0.001
1,2-Methyl-t-butylbenzene	A11	0.0018	0.0157	0.001	0.001
5-Methylindan	A11	0.0013	0.0101	0.001	0.001
4-Methylindan	A11	0.0010	0.0078	0.001	0.001
1,2-Ethyl-n-propylbenzene	A11	0.0009	0.0078	0.001	0.001
2-Methylindan	A11	0.0005	0.0039	0.000	0.000
1,3-Methyl-n-butylbenzene	A11	0.0009	0.0078	0.001	0.001
1,3-Di-i-propylbenzene	A11	0.0019	0.0181	0.001	0.001
sec-Pentylbenzene	A11	0.0030	0.0261	0.002	0.002
n-Pentylbenzene	A11	0.0005	0.0044	0.000	0.000
1,2-Di-n-propylbenzene	A11	0.0031	0.0295	0.002	0.002
1,4-Di-i-propylbenzene	A11	0.0011	0.0105	0.001	0.001
Tetrahydronaphthalene	A10	0.0045	0.0350	0.003	0.003
Naphthalene	A10	0.0007	0.0053	0.001	0.001
1-t-Butyl-3,5-dimethylbenzene	A12	0.0010	0.0095	0.001	0.001
1,4-Ethyl-t-butylbenzene	A11	0.0010	0.0095	0.001	0.001
1,3-Di-n-propylbenzene	A12	0.0017	0.0162	0.001	0.001
UnknownC11s	U11	0.0117	0.1074	0.008	0.008
n-Dodecane	P12	0.0314	0.3142	0.022	0.022
1,3,5-Triethylbenzene	A12	0.0056	0.0534	0.003	0.003
1,2,4-Triethylbenzene	A12	0.0020	0.0191	0.001	0.001
1,4-Methyl-n-pentylbenzene	A12	0.0037	0.0352	0.003	0.003
n-Hexylbenzene	A12	0.0042	0.0401	0.002	0.002
1,2,3,4,5-Pentamethylbenzene	A13	0.0054	0.0470	0.004	0.004
2-Methylnaphthalene	A11	0.0013	0.0109	0.001	0.001
1-Methylnaphthalene	A11	0.0012	0.0100	0.001	0.001
UnknownC12s	U12	0.0142	0.1304	0.009	0.009
n-Tridecane	P13	0.0179	0.1938	0.014	0.014
UnknownC13s	U13	0.0231	0.2501	0.018	0.018
n-Tetradecane	P14	0.0032	0.0373	0.003	0.003
UnknownC14s	U14	0.0117	0.1363	0.010	0.010
n-Pentadecane	P15	0.0004	0.0050	0.000	0.000
UnknownC15s	U15	0.0104	0.1297	0.009	0.009
n-Hexadecane	P16	0.0001	0.0014	0.000	0.000
UnknownC16s	U16	0.0012	0.0160	0.001	0.001
TOTAL		100.00000	100.00000	0.8674	0.8711

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BTEX COMPONENTS	MOLE%	WT%
BENZENE	0.0182	0.0835
TOLUENE	0.0347	0.1878
ETHYLBENZENE	0.0029	0.0181
XYLENES	0.0283	0.1766
TOTAL BTEX	0.0841	0.4660

**(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 :	954.5 /scf	959.7 /scf
	NET WET REAL :	937.8 /scf	943.0 /scf
HHV	GROSS DRY REAL :	1058.1 /scf	1063.9 /scf
	GROSS WET REAL :	1039.6 /scf	1045.4 /scf
	NET HEATING VALUE (60 °F ideal reaction):		21297.6 Btu/lbm
	GROSS HEATING VALUE (60°F ideal reaction):		23613.1 Btu/lbm
	RELATIVE DENSITY (AIR=1):		0.5870
	DENSITY		0.04486 lb/scf
	COMPRESSIBILITY FACTOR :		0.9979
	REGULAR WOBBE INDEX		1382.5

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

Net Dry Ideal BTU	6686.1 /scf	Relative Density - SG (Air=1)	4.6204	***C6+factors***
Gross Dry Ideal BTU	7201.7 /scf	**Z** Compressiblity Factor	0.99882	0.99795
Net Dry Ideal BTU	18506.7 /lb	Density Factor	352.686 lbm/1000 ft3	
Gross Dry Ideal BTU	19929.6 /lb	Molar Mass or MW	133.833 g/mol	
		Volume Liquid Ideal gas	0.287 scf/gal	18.2

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.