



303-637-0150

EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PROJECT NO. :	201409145	ANALYSIS NO. :	12
COMPANY NAME :	NIGHTHAWK PRODUCTION	ANALYSIS DATE :	SEPTEMBER 29, 2014
ACCOUNT NO. :		SAMPLE DATE :	SEPTEMBER 23, 2014
PRODUCER :	ARIKAREE CREEK	CYLINDER NO. :	0690
LEASE NO. :		SAMPLED BY :	JOHN MOSER - EMPACT
NAME/DESCRIP :	SALES GAS 15:00 SILVERTON 16-10		
FIELD DATA		SAMPLE TEMP. :	96
SAMPLE PRES. :	27	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE; LENGTH OF H2S STAIN @ 2.0 PPM (1-7PPM) 15:05		

<u>COMPONENT</u>	<u>MOLE %</u>	<u>MASS %</u>	<u>GPM @ 14.650</u>	<u>GPM @ 14.730</u>
ALCOHOLS	0.0015	0.0038		
HELIUM	0.57	0.08	---	---
HYDROGEN	0.03	0.00	---	---
OXYGEN/ARGON	0.39	0.42	---	---
NITROGEN	50.49	47.95	---	---
CARBON DIOXIDE	2.68	4.00	---	---
METHANE	27.45800	14.93730	---	---
ETHANE	3.9687	4.0462	1.0583	1.0640
PROPANE	5.6764	8.4868	1.5589	1.5674
I-BUTANE	1.2194	2.4030	0.3977	0.3999
N-BUTANE	3.8390	7.5655	1.2062	1.2127
I-PENTANE	0.9276	2.2670	0.3358	0.3375
N-PENTANE	1.2420	3.0383	0.4477	0.4501
HEXANES PLUS	1.5074	4.8021	0.6348	0.6374
<u>TOTALS</u>	<u>100.00000</u>	<u>100.00000</u>	<u>5.6394</u>	<u>5.6690</u>

<u>BTEX COMPONENTS</u>	<u>MOLE%</u>	<u>WT%</u>		<u>BTU @ 14.650</u>	<u>BTU @ 14.730</u>
BENZENE	0.0031	0.0082	LOW NET DRY REAL :	748.2 /scf	752.3 /scf
TOLUENE	0.0033	0.0103	NET WET REAL :	735.1 /scf	739.2 /scf
ETHYLBENZENE	0.0013	0.0047	HIGH GROSS DRY REAL :	817.8 /scf	822.3 /scf
XYLENES	0.0032	0.0116	GROSS WET REAL :	803.5 /scf	808.0 /scf
<u>TOTAL BTEX</u>	<u>0.0109</u>	<u>0.0348</u>	NET DRY REAL :	9649.8 /lb	9702.5 /lb
			GROSS DRY REAL :	10549.1 /lb	10606.7 /lb

RELATIVE DENSITY (AIR=1): 1.0166
 COMPRESSIBILITY FACTOR : 0.99757

(CALC: GPA STD 2145 & TP-17 @14.696 & 60 F)

*(DETAILED HYDROCARBON ANALYSIS/NJ 1993) ; ASTM D6730

THIS DATA HAS BEEN ACQUIRED THROUGH APPLICATION OF CURRENT STATE-OF-THE-ART ANALYTICAL TECHNIQUES.
 THE USE OF THIS INFORMATION IS THE RESPONSIBILITY OF THE USER. EMPACT ANALYTICAL SYSTEMS, ASSUMES NO
 RESPONSIBILITY FOR ACCURACY OF THE REPORTED INFORMATION NOR ANY CONSEQUENCES OF ITS APPLICATION.



303-637-0150

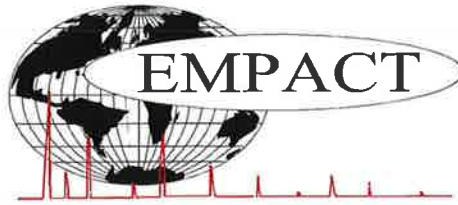
EXTENDED NATURAL GAS ANALYSIS (*DHA)

GLYCALC INFORMATION

PROJECT NO. :	201409145	ANALYSIS NO. :	12
COMPANY NAME :	NIGHTHAWK PRODUCTION	ANALYSIS DATE:	SEPTEMBER 29, 2014
ACCOUNT NO. :		SAMPLE DATE :	SEPTEMBER 23, 2014
PRODUCER :	ARIKAREE CREEK	CYLINDER NO. :	0690
LEASE NO. :		SAMPLED BY :	JOHN MOSER - EMPACT
NAME/DESCRIP :	SALES GAS 15:00 SILVERTON 16-10		
FIELD DATA		SAMPLE TEMP. :	96
SAMPLE PRES. :	27	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE; LENGTH OF H2S STAIN @ 2.0 PPM (1-7PPM) 15:05		

<u>Componet</u>	<u>Mole %</u>	<u>Wt %</u>
Helium	0.57	0.08
Hydrogen	0.03	0.00
Carbon Dioxide	2.68	4.00
Nitrogen	50.49	47.95
Methane	27.45800	14.93730
Ethane	3.9687	4.0462
Propane	5.6764	8.4868
Isobutane	1.2194	2.4030
n-Butane	3.8390	7.5655
Isopentane	0.8958	2.1914
n-Pentane	1.2420	3.0383
Cyclopentane	0.0318	0.0756
n-Hexane	0.3375	0.9861
Cyclohexane	0.0845	0.2411
Other Hexanes	0.4759	1.3838
Heptanes	0.3006	1.0155
Methycyclohexane	0.0869	0.2893
2,2,4 Trimethylpentane	0.0005	0.0019
Benzene	0.0031	0.0082
Toluene	0.0033	0.0103
Ethylbenzene	0.0013	0.0047
Xylenes	0.0032	0.0116
C8+ Heavies	0.2106	0.8496
Subtotal	99.60850	99.57620
Oxygen/Argon	0.39	0.42
Alcohols	0.0015	0.0038
Total	100.00000	100.00000

THE DATA PRESENTED HEREIN HAS BEEN ACQUIRED THROUGH JUDICIOUS APPLICATION OF CURRENT STATE-OF-THE ART ANALYTICAL TECHNIQUES. THE APPLICATIONS OF THIS INFORMATION IS THE RESPONSIBILITY OF THE USER. EMPACT ANALYTICAL SYSTEMS, INC. ASSUMES NO RESPONSIBILITY FOR ACCURACY OF THE REPORTED INFORMATION NOR ANY CONSEQUENCES OF IT'S APPLICATION.



**EXTENDED NATURAL GAS ANALYSIS (*DHA)
DHA COMPONENT LIST**

PROJECT NO. :	201409145	ANALYSIS NO. :	12
COMPANY NAME :	NIGHTHAWK PRODUCTION	ANALYSIS DATE:	SEPTEMBER 29, 2014
ACCOUNT NO. :		SAMPLE DATE :	SEPTEMBER 23, 2014
PRODUCER :	ARIKAREE CREEK	CYLINDER NO. :	0690
LEASE NO. :		SAMPLED BY :	JOHN MOSER - EMPACT
NAME/DESCRIP :	SALES GAS 15:00 SILVERTON 16-10		
FIELD DATA		SAMPLE TEMP. :	96
SAMPLE PRES. :	27	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE; LENGTH OF H2S STAIN @ 2.0 PPM (1-7PPM) 15:05		

COMPONENT	PIANO #	MOLE %	MASS %	GPM @	
				14.650	14.730
Helium	---	0.57	0.08	---	---
Hydrogen	---	0.03	0.00	---	---
Oxygen/Argon	---	0.39	0.42	---	---
Nitrogen	---	50.49	47.95	---	---
Carbon Dioxide	---	2.68	4.00	---	---
Methane	P1	27.45800	14.93730	---	---
Ethane	P2	3.9687	4.0462	1.058	1.064
Propane	P3	5.6764	8.4868	1.559	1.567
i-Butane	I4	1.2194	2.4030	0.398	0.400
n-Butane	P4	3.8388	7.5651	1.206	1.213
2,2-Dimethylpropane	I5	0.0070	0.0171	0.003	0.003
i-Pentane	I5	0.8888	2.1743	0.324	0.326
UnknownC4s	U4	0.0002	0.0004	0.000	0.000
n-Pentane	P5	1.2403	3.0341	0.448	0.450
t-Butanol	X4	0.0015	0.0038	0.001	0.001
2,2-Dimethylbutane	I6	0.0060	0.0175	0.002	0.002
Cyclopentane	N5	0.0318	0.0756	0.009	0.009
2,3-Dimethylbutane	I6	0.0251	0.0733	0.010	0.010
2-Methylpentane	I6	0.2300	0.6720	0.095	0.096
3-Methylpentane	I6	0.1170	0.3419	0.048	0.048
UnknownC5s	U5	0.0017	0.0042	0.001	0.001
n-Hexane	P6	0.3375	0.9861	0.138	0.139
2,2-Dimethylpentane	I7	0.0006	0.0020	0.000	0.000
Methylcyclopentane	N6	0.0977	0.2788	0.034	0.034
2,4-Dimethylpentane	I7	0.0067	0.0228	0.003	0.003
2,2,3-Trimethylbutane	I7	0.0005	0.0017	0.000	0.000
Benzene	A6	0.0031	0.0082	0.001	0.001
3,3-Dimethylpentane	I7	0.0001	0.0003	0.000	0.000
Cyclohexane	N6	0.0845	0.2411	0.029	0.029
2-Methylhexane	I7	0.0367	0.1247	0.017	0.017
2,3-Dimethylpentane	I7	0.0104	0.0353	0.005	0.005
1,1-Dimethylcyclopentane	N7	0.0115	0.0383	0.005	0.005
3-Methylhexane	I7	0.0471	0.1600	0.022	0.022
1c,3-Dimethylcyclopentane	N7	0.0183	0.0609	0.008	0.008
1t,3-Dimethylcyclopentane	N7	0.0154	0.0513	0.007	0.007
3-Ethylpentane	I7	0.0017	0.0058	0.001	0.001
1t,2-Dimethylcyclopentane	N7	0.0320	0.1065	0.015	0.015
2,2,4-Trimethylpentane	I8	0.0005	0.0019	0.000	0.000
UnknownC6s	U6	0.0001	0.0003	0.000	0.000
n-Heptane	P7	0.1096	0.3724	0.050	0.050
1c,2-Dimethylcyclopentane	N7	0.0024	0.0080	0.001	0.001

Methylcyclohexane	N7	0.0869	0.2893	0.035	0.035
2,2-Dimethylhexane	I8	0.0099	0.0384	0.005	0.005
Ethylcyclopentane	N7	0.0046	0.0153	0.002	0.002
2,5-Dimethylhexane	I8	0.0012	0.0046	0.001	0.001
2,2,3-Trimethylpentane	I8	0.0009	0.0035	0.000	0.000
2,4-Dimethylhexane	I8	0.0033	0.0128	0.002	0.002
1c,2t,4-Trimethylcyclopentane	N8	0.0068	0.0259	0.003	0.003
3,3-Dimethylhexane	I8	0.0007	0.0027	0.000	0.000
1t,2c,4-Trimethylcyclopentane	N8	0.0100	0.0380	0.005	0.005
2,3,4-Trimethylpentane	I8	0.0003	0.0012	0.000	0.000
Toluene	A7	0.0033	0.0103	0.001	0.001
2,3-Dimethylhexane	I8	0.0030	0.0116	0.002	0.002
2-Methyl-3-ethylpentane	I8	0.0010	0.0039	0.000	0.000
1,1,2-Trimethylcyclopentane	N8	0.0001	0.0004	0.000	0.000
2-Methylheptane	I8	0.0193	0.0748	0.010	0.010
4-Methylheptane	I8	0.0052	0.0201	0.003	0.003
3-Methyl-3-ethylpentane	I8	0.0010	0.0039	0.000	0.000
3,4-Dimethylhexane	I8	0.0007	0.0027	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0003	0.0012	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0004	0.0015	0.000	0.000
3-Methylheptane	I8	0.0060	0.0232	0.003	0.003
1c,2t,3-Trimethylcyclopentane	N8	0.0167	0.0635	0.009	0.009
3-Ethylhexane	I8	0.0019	0.0074	0.001	0.001
1t,4-Dimethylcyclohexane	N8	0.0045	0.0171	0.002	0.002
1,1-Dimethylcyclohexane	N8	0.0013	0.0050	0.001	0.001
3t-Ethylmethylcyclopentane	N8	0.0013	0.0050	0.001	0.001
2t-Ethylmethylcyclopentane	N8	0.0010	0.0038	0.001	0.001
1,1-Methylethylcyclopentane	N8	0.0030	0.0114	0.002	0.002
2,2,4-Trimethylhexane	I9	0.0003	0.0013	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0049	0.0187	0.003	0.003
UnknownC7s	U7	0.0030	0.0102	0.001	0.001
n-Octane	P8	0.0172	0.0666	0.009	0.009
1c,4-Dimethylcyclohexane	N8	0.0163	0.0620	0.008	0.008
i-Propylcyclopentane	I8	0.0005	0.0019	0.000	0.000
2,4,4-Trimethylhexane	I9	0.0003	0.0013	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0005	0.0022	0.000	0.000
2,2,3,4-Tetramethylpentane	I9	0.0004	0.0017	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0002	0.0009	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0010	0.0038	0.001	0.001
2,2-Dimethylheptane	I9	0.0002	0.0009	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0060	0.0257	0.003	0.003
2,2,3-Trimethylhexane	I9	0.0031	0.0135	0.002	0.002
2,4-Dimethylheptane	I9	0.0002	0.0009	0.000	0.000
4,4-Dimethylheptane	I9	0.0009	0.0039	0.000	0.000
Ethylcyclohexane	N8	0.0025	0.0095	0.001	0.001
n-Propylcyclopentane	N8	0.0016	0.0061	0.001	0.001
1c,3c,5-Trimethylcyclohexane	N9	0.0003	0.0013	0.000	0.000
2,5-Dimethylheptane	I9	0.0006	0.0026	0.000	0.000
3,3-Dimethylheptane	I9	0.0006	0.0026	0.000	0.000
3,5-Dimethylheptane	I9	0.0004	0.0017	0.000	0.000
2,6-Dimethylheptane	I9	0.0005	0.0022	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0008	0.0034	0.000	0.000
Ethylbenzene	I8	0.0013	0.0047	0.001	0.001
1c,2t,4t-Trimethylcyclohexane	N9	0.0010	0.0043	0.001	0.001
2,3-Dimethylheptane	I9	0.0003	0.0013	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0015	0.0054	0.001	0.001
1,4-Dimethylbenzene (p-Xylene)	A8	0.0009	0.0033	0.000	0.000
3,4-Dimethylheptane	I9	0.0004	0.0017	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0005	0.0022	0.000	0.000
4-Ethylheptane	I9	0.0003	0.0013	0.000	0.000
4-Methyloctane	I9	0.0017	0.0074	0.001	0.001
2-Methyloctane	I9	0.0021	0.0091	0.001	0.001
1c,2t,3-Trimethylcyclohexane	N9	0.0004	0.0017	0.000	0.000
3-Ethylheptane	I9	0.0006	0.0026	0.000	0.000

3-Methyloctane	I9	0.0032	0.0139	0.002	0.002
1c,2t,4c-Trimethylcyclohexane	I9	0.0003	0.0013	0.000	0.000
1,1,2-Trimethylcyclohexane	N9	0.0002	0.0009	0.000	0.000
3,3-Diethylpentane	I9	0.0003	0.0013	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0008	0.0029	0.000	0.000
i-Butylcyclopentane	N9	0.0011	0.0047	0.001	0.001
UnknownC8s	U8	0.0005	0.0019	0.000	0.000
n-Nonane	P9	0.0068	0.0296	0.004	0.004
1,1-Methylethylcyclohexane	N9	0.0029	0.0124	0.002	0.002
i-Propylbenzene	A9	0.0004	0.0016	0.000	0.000
i-Propylcyclohexane	N9	0.0004	0.0017	0.000	0.000
2,2-Dimethyloctane	I10	0.0001	0.0005	0.000	0.000
2,4-Dimethyloctane	I10	0.0002	0.0010	0.000	0.000
2,6-Dimethyloctane	I10	0.0002	0.0010	0.000	0.000
n-Butylcyclopentane	N9	0.0011	0.0047	0.001	0.001
3,3-Dimethyloctane	I10	0.0002	0.0010	0.000	0.000
n-Propylbenzene	A9	0.0011	0.0045	0.000	0.000
3,6-Dimethyloctane	I10	0.0004	0.0019	0.000	0.000
3-Methyl-5-ethylheptane	I10	0.0005	0.0024	0.000	0.000
1,3-Methylethylbenzene	A9	0.0004	0.0016	0.000	0.000
1,4-Methylethylbenzene	A9	0.0003	0.0012	0.000	0.000
1,3,5-Trimethylbenzene	A9	0.0005	0.0020	0.000	0.000
2,3-Dimethyloctane	I10	0.0003	0.0015	0.000	0.000
5-Methylnonane	I10	0.0008	0.0039	0.000	0.000
1,2-Methylethylbenzene	A9	0.0005	0.0020	0.000	0.000
2-Methylnonane	I10	0.0001	0.0005	0.000	0.000
3-Ethyloctane	I10	0.0001	0.0005	0.000	0.000
3-Methylnonane	I10	0.0005	0.0024	0.000	0.000
1,2,4-Trimethylbenzene	A9	0.0001	0.0004	0.000	0.000
t-Butylbenzene	A10	0.0006	0.0028	0.000	0.000
i-Butylcyclohexane	N10	0.0003	0.0014	0.000	0.000
1t-Methyl-2-n-propylcyclohexane	I10	0.0001	0.0005	0.000	0.000
i-Butylbenzene	A10	0.0002	0.0009	0.000	0.000
sec-Butylbenzene	A10	0.0001	0.0004	0.000	0.000
UnknownC9s	U9	0.0065	0.0283	0.004	0.004
n-Decane	P10	0.0026	0.0125	0.002	0.002
1,2,3-Trimethylbenzene	A9	0.0002	0.0008	0.000	0.000
1,3-Methyl-i-propylbenzene	A10	0.0001	0.0004	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0001	0.0004	0.000	0.000
Sec-Butylcyclohexane	A10	0.0007	0.0033	0.000	0.000
1,2-Methyl-i-propylbenzene	A10	0.0002	0.0009	0.000	0.000
3-Ethylnonane	I10	0.0001	0.0005	0.000	0.000
1,3-Diethylbenzene	A10	0.0001	0.0004	0.000	0.000
1,3-Methyl-n-propylbenzene	A10	0.0001	0.0004	0.000	0.000
1,4-Diethylbenzene	A10	0.0002	0.0009	0.000	0.000
1,4-Methyl-n-propylbenzene	A10	0.0002	0.0009	0.000	0.000
n-Butylbenzene	A10	0.0001	0.0004	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0001	0.0004	0.000	0.000
1,2-Diethylbenzene	A10	0.0001	0.0004	0.000	0.000
t-Decahydronaphthalene	A9	0.0001	0.0005	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0001	0.0004	0.000	0.000
1,4-Dimethyl-2-ethylbenzene	A10	0.0001	0.0004	0.000	0.000
1,2-Dimethyl-4-ethylbenzene	A10	0.0002	0.0009	0.000	0.000
1,4-Methyl-t-butylbenzene	A11	0.0001	0.0005	0.000	0.000
UnknownC10s	U10	0.0045	0.0217	0.003	0.003
n-Undecane	P11	0.0008	0.0042	0.001	0.001
1,2-Methyl-t-butylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2-Ethyl-n-propylbenzene	A11	0.0001	0.0005	0.000	0.000
1,4-Di-i-propylbenzene	A11	0.0001	0.0005	0.000	0.000
UnknownC11s	U11	0.0010	0.0053	0.001	0.001
n-Dodecane	P12	0.0002	0.0012	0.000	0.000
1,2,4-Triethylbenzene	A12	0.0001	0.0005	0.000	0.000
UnknownC12s	U12	0.0002	0.0011	0.000	0.000
n-Tridecane	P13	0.0001	0.0006	0.000	0.000

n-Heptadecane	P17	0.0001	0.0008	0.000	0.000
n-Octadecane	P18	0.0001	0.0009	0.000	0.000
<u>TOTAL</u>		<u>100.00000</u>	<u>100.00000</u>	<u>5.6414</u>	<u>5.6710</u>

<u>BTEX COMPONENTS</u>	<u>MOLE%</u>	<u>WT%</u>	<u>BTU @</u>	<u>14.650</u>	<u>14.730</u>
BENZENE	0.0031	0.0082	LOW NET DRY REAL :	748.2 /scf	752.3 /scf
TOLUENE	0.0033	0.0103	NET WET REAL :	735.1 /scf	739.2 /scf
ETHYLBENZENE	0.0013	0.0047	HIGH GROSS DRY REAL :	817.8 /scf	822.3 /scf
<u>XYLENES</u>	<u>0.0032</u>	<u>0.0116</u>	GROSS WET REAL :	803.5 /scf	808.0 /scf
<u>TOTAL BTEX</u>	<u>0.0109</u>	<u>0.0348</u>	NET DRY REAL :	9649.8 /lb	9702.5 /lb
			GROSS DRY REAL :	10549.1 /lb	10606.7 /lb

RELATIVE DENSITY (AIR=1): 1.0166
 COMPRESSIBILITY FACTOR : 0.99757

(CALC: GPA STD 2145 & TP-17 @14.696 & 60 F)

*(DETAILED HYDROCARBON ANALYSIS/NJ 1993) ; ASTM D6730

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 RESPONSIBILITY FOR ACCURACY OF THE REPORTED INFORMATION NOR ANY CONSEQUENCES OF ITS APPLICATION.