

EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

LEASE #: NAME/DESCRIP : **JONES 18-14-10-59
SEPARATOR GAS**

PROJECT NO. : **201703033** ANALYSIS NO. : **01**

COMPANY NAME : **CARRIZO OIL & GAS, INC** ANALYSIS DATE: **MARCH 09, 2017 10:05**

OFFICE / BRANCH: **HOUSTON, TX** SAMPLE DATE : **MARCH 6, 2017 15:50**

CUSTOMER REF: TO:

PRODUCER : EFFECTIVE DATE:

FIELD DATA

SAMPLE CYCLE: SAMPLE TYPE: SPOT

SAMPLE PRES. : 23.5 psig CYLINDER NO. : 1590

LAB PRES: psig SAMPLED BY : JOHN MOSER

SAMPLE TEMP. : 58.4 °f SAMPLING COMPANY: EMPACT

AMBIENT TEMP.: °f H2S BY STAIN TUBE: **35.0** ppm

H2O BY STAIN TUBE: - #/mmcf CO2 BY STAIN TUBE: - Mol %

FIELD COMMENTS: NO PROBE

LAB COMMENTS:

COMPONENT	MOLE %	MASS %	GPM @ 14.730	GPM @ 14.650
HELIUM	0.02	0.00	---	---
HYDROGEN	0.00	0.00	---	---
OXYGEN/ARGON	0.01	0.01	---	---
NITROGEN	0.9900	1.0600	---	---
CARBON DIOXIDE	3.36	5.65	---	---
METHANE	65.00590	39.87080	---	---
ETHANE	10.5875	12.1715	2.8454	2.8299
PROPANE	10.7316	18.0922	2.9713	2.9552
I-BUTANE	1.1741	2.6090	0.3859	0.3838
N-BUTANE	4.5790	10.1752	1.4509	1.4430
I-PENTANE	1.1154	3.0670	0.4010	0.3989
N-PENTANE	1.4696	4.0537	0.5350	0.5321
HEXANES PLUS	0.9569	3.2406	0.3785	0.3766
TOTALS	100.00000	100.00000	8.9680	8.9195

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.730	14.650
BENZENE	0.0381	0.1138	LOW NET DRY REAL :	1331.7 /scf	1324.5 /scf
TOLUENE	0.0082	0.0289	NET WET REAL :	1308.5 /scf	1301.3 /scf
ETHYLBENZENE	0.0000	0.0000	HIGH GROSS DRY REAL :	1460.5 /scf	1452.6 /scf
XYLENES	0.0017	0.0068	GROSS WET REAL :	1435.1 /scf	1427.2 /scf
TOTAL BTEX	0.0480	0.1495	NET DRY REAL :	19337.2 /lb	19232.2 /lb
			GROSS DRY REAL :	21214.6 /lb	21099.3 /lb

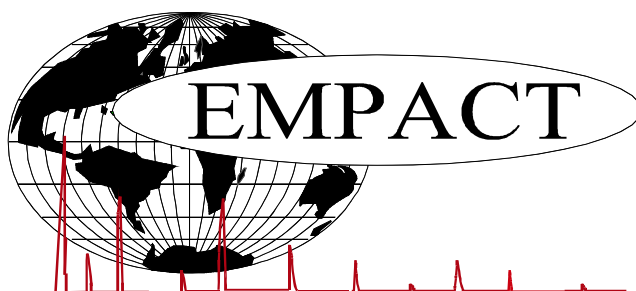
(CALC: GPA STD 2145 & TP-17 @14.696 & 60 F)

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

RELATIVE DENSITY (AIR=1): 0.9024

COMPRESSIBILITY FACTOR : 0.99478

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.



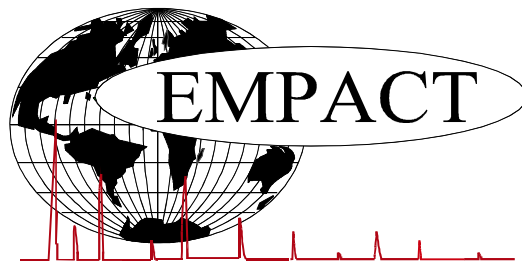
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GLYCALC INFORMATION

PROJECT NO. :	201703033	ANALYSIS NO. :	01
COMPANY NAME :	CARRIZO OIL & GAS, INC	ANALYSIS DATE:	MARCH 09, 2017 10:05
ACCOUNT NO. :		SAMPLE DATE :	MARCH 6, 2017 15:50
PRODUCER :		CYLINDER NO. :	1590
LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	JONES 18-14-10-59 SEPARATOR GAS		
FIELD DATA		SAMPLE TEMP. :	58.4
SAMPLE PRES. :	23.5	AMBIENT TEMP.:	
COMMENTS :	NO PROBE SPOT		

Componet	Mole %	Wt %
Helium	0.02	0.00
Hydrogen	0.00	0.00
Carbon Dioxide	3.36	5.65
Nitrogen	0.99	1.06
Methane	65.00590	39.87080
Ethane	10.5875	12.1715
Propane	10.7316	18.0922
Isobutane	1.1741	2.6090
n-Butane	4.5790	10.1752
Isopentane	0.9883	2.7262
n-Pentane	1.4696	4.0537
Cyclopentane	0.1271	0.3408
n-Hexane	0.2512	0.8276
Cyclohexane	0.0492	0.1583
Other Hexanes	0.4499	1.4719
Heptanes	0.1111	0.4225
Methycyclohexane	0.0098	0.0368
2,2,4 Trimethylpentane	0.0028	0.0122
Benzene	0.0381	0.1138
Toluene	0.0082	0.0289
Ethylbenzene	0.0000	0.0000
Xylenes	0.0017	0.0068
C8+ Heavies	0.0349	0.1618
<u>Subtotal</u>	<u>99.99000</u>	<u>99.99000</u>
<u>Oxygen/Argon</u>	<u>0.01</u>	<u>0.01</u>
Total	100.00000	100.00000

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DHA COMPONENT LIST

PROJECT NO. :	201703033	ANALYSIS NO. :	01
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LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	JONES 18-14-10-59		
	SEPARATOR GAS		
FIELD DATA		SAMPLE TEMP. :	58.4
SAMPLE PRES. :	23.5	AMBIENT TEMP.:	
COMMENTS :	NO PROBE		
	SPOT		

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.730	GPM @ 14.650
Helium	---	0.02	0.00	---	---
Oxygen/Argon	---	0.01	0.01	---	---
Nitrogen	---	0.99	1.06	---	---
Carbon Dioxide	---	3.36	5.65	---	---
Methane	P1	65.00590	39.87080	---	---
Ethane	P2	10.5875	12.1715	2.845	2.830
Propane	P3	10.7316	18.0922	2.971	2.955
i-Butane	I4	1.1741	2.6090	0.386	0.384
n-Butane	P4	4.5790	10.1752	1.451	1.443
2,2-Dimethylpropane	I5	0.0031	0.0086	0.001	0.001
i-Pentane	I5	0.9852	2.7176	0.362	0.360
n-Pentane	P5	1.4693	4.0529	0.535	0.532
2,2-Dimethylbutane	I6	0.0020	0.0066	0.001	0.001
Cyclopentane	N5	0.1271	0.3408	0.038	0.038
2,3-Dimethylbutane	I6	0.0150	0.0494	0.006	0.006
2-Methylpentane	I6	0.1999	0.6586	0.084	0.083
3-Methylpentane	I6	0.0992	0.3268	0.040	0.040
UnknownC5s	U5	0.0003	0.0008	0.000	0.000
n-Hexane	P6	0.2512	0.8276	0.104	0.103
2,2-Dimethylpentane	I7	0.0002	0.0008	0.000	0.000
Methylcyclopentane	N6	0.1334	0.4292	0.047	0.047
2,4-Dimethylpentane	I7	0.0035	0.0134	0.002	0.002
2,2,3-Trimethylbutane	I7	0.0009	0.0034	0.000	0.000
Benzene	A6	0.0381	0.1138	0.011	0.011
3,3-Dimethylpentane	I7	0.0002	0.0008	0.000	0.000
Cyclohexane	N6	0.0492	0.1583	0.017	0.017
2-Methylhexane	I7	0.0110	0.0421	0.005	0.005
2,3-Dimethylpentane	I7	0.0128	0.0491	0.006	0.006
1,1-Dimethylcyclopentane	N7	0.0043	0.0161	0.002	0.002
3-Methylhexane	I7	0.0195	0.0747	0.009	0.009
1c,3-Dimethylcyclopentane	N7	0.0041	0.0154	0.002	0.002
1t,3-Dimethylcyclopentane	N7	0.0072	0.0270	0.003	0.003
3-Ethylpentane	I7	0.0015	0.0057	0.001	0.001
1t,2-Dimethylcyclopentane	N7	0.0208	0.0781	0.010	0.010
2,2,4-Trimethylpentane	I8	0.0028	0.0122	0.001	0.001

UnknownC6s	U6	0.0004	0.0013	0.000	0.000
n-Heptane	P7	0.0116	0.0444	0.005	0.005
1c,2-Dimethylcyclopentane	N7	0.0004	0.0015	0.000	0.000
Methylcyclohexane	N7	0.0098	0.0368	0.004	0.004
2,2-Dimethylhexane	I8	0.0004	0.0018	0.000	0.000
1,1,3-Trimethylcyclopentane	N7	0.0001	0.0004	0.000	0.000
Ethylcyclopentane	N7	0.0025	0.0094	0.001	0.001
2,5-Dimethylhexane	I8	0.0002	0.0009	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0001	0.0004	0.000	0.000
2,4-Dimethylhexane	I8	0.0004	0.0018	0.000	0.000
1c,2t,4-Trimethylcyclopentane	N8	0.0008	0.0034	0.000	0.000
3,3-Dimethylhexane	I8	0.0001	0.0004	0.000	0.000
1t,2c,4-Trimethylcyclopentane	N8	0.0009	0.0039	0.000	0.000
2,3,4-Trimethylpentane	I8	0.0004	0.0018	0.000	0.000
2,3,3-Trimethylpentane	I8	0.0002	0.0009	0.000	0.000
Toluene	A7	0.0082	0.0289	0.003	0.003
2,3-Dimethylhexane	I8	0.0004	0.0018	0.000	0.000
2-Methyl-3-ethylpentane	I8	0.0002	0.0009	0.000	0.000
1,1,2-Trimethylcyclopentane	N8	0.0001	0.0004	0.000	0.000
2-Methylheptane	I8	0.0024	0.0105	0.001	0.001
4-Methylheptane	I8	0.0009	0.0039	0.000	0.000
3-Methyl-3-ethylpentane	I8	0.0002	0.0009	0.000	0.000
3,4-Dimethylhexane	I8	0.0001	0.0004	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0001	0.0004	0.000	0.000
3-Methylheptane	I8	0.0001	0.0004	0.000	0.000
1c,2t,3-Trimethylcyclopentane	N8	0.0024	0.0103	0.001	0.001
3-Ethylhexane	I8	0.0005	0.0022	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0006	0.0026	0.000	0.000
1,1-Dimethylcyclohexane	N8	0.0002	0.0008	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0006	0.0026	0.000	0.000
2t-Ethylmethylcyclopentane	N8	0.0005	0.0021	0.000	0.000
1,1-Methylethylcyclopentane	N8	0.0013	0.0056	0.001	0.001
1t,2-Dimethylcyclohexane	N8	0.0007	0.0030	0.000	0.000
1c,2c,3-Trimethylcyclopentane	N8	0.0002	0.0008	0.000	0.000
UnknownC7s	U7	0.0105	0.0402	0.005	0.005
n-Octane	P8	0.0043	0.0188	0.002	0.002
1c,4-Dimethylcyclohexane	N8	0.0006	0.0026	0.000	0.000
2,4,4-Trimethylhexane	I9	0.0002	0.0010	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0001	0.0005	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0002	0.0008	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0016	0.0077	0.001	0.001
2,2,3-Trimethylhexane	I9	0.0003	0.0015	0.000	0.000
2,4-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
4,4-Dimethylheptane	I9	0.0003	0.0015	0.000	0.000
Ethylcyclohexane	N8	0.0005	0.0021	0.000	0.000
n-Propylcyclopentane	N8	0.0002	0.0008	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0001	0.0005	0.000	0.000
2,5-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
3,3-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
2,6-Dimethylheptane	I9	0.0009	0.0044	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0003	0.0015	0.000	0.000
2,3-Dimethylheptane	I9	0.0002	0.0010	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0007	0.0028	0.000	0.000
1,4-Dimethylbenzene (p-Xylene)	A8	0.0003	0.0012	0.000	0.000
3,4-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0003	0.0015	0.000	0.000
4-Ethylheptane	I9	0.0001	0.0005	0.000	0.000
4-Methyloctane	I9	0.0002	0.0010	0.000	0.000
2-Methyloctane	I9	0.0003	0.0015	0.000	0.000

1c,2t,3-Trimethylcyclohexane	N9	0.0001	0.0005	0.000	0.000
3-Methyloctane	I9	0.0002	0.0010	0.000	0.000
1c,2t,4c-Trimethylcyclohexane	I9	0.0001	0.0005	0.000	0.000
1,1,2-Trimethylcyclohexane	N9	0.0001	0.0005	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0007	0.0028	0.000	0.000
i-Butylcyclopentane	N9	0.0003	0.0015	0.000	0.000
UnknownC8s	U8	0.0005	0.0022	0.000	0.000
n-Nonane	P9	0.0012	0.0059	0.001	0.001
1,1-Methylethylcyclohexane	N9	0.0001	0.0005	0.000	0.000
i-Propylbenzene	A9	0.0002	0.0009	0.000	0.000
i-Propylcyclohexane	N9	0.0002	0.0010	0.000	0.000
n-Butylcyclopentane	N9	0.0002	0.0010	0.000	0.000
3,3-Dimethyloctane	I10	0.0002	0.0011	0.000	0.000
n-Propylbenzene	A9	0.0002	0.0009	0.000	0.000
3-Methyl-5-ethylheptane	I10	0.0002	0.0011	0.000	0.000
1,3-Methylethylbenzene	A9	0.0002	0.0009	0.000	0.000
5-Methylnonane	I10	0.0002	0.0011	0.000	0.000
1,2-Methylethylbenzene	A9	0.0001	0.0005	0.000	0.000
t-Butylbenzene	A10	0.0001	0.0005	0.000	0.000
UnknownC9s	U9	0.0028	0.0137	0.002	0.002
n-Decane	P10	0.0002	0.0011	0.000	0.000
Sec-Butylcyclohexane	A10	0.0001	0.0005	0.000	0.000
1,2-Methyl-i-propylbenzene	A10	0.0001	0.0005	0.000	0.000
UnknownC10s	U10	0.0005	0.0027	0.000	0.000
n-Undecane	P11	0.0001	0.0006	0.000	0.000
UnknownC11s	U11	0.0001	0.0006	0.000	0.000
n-Tetradecane	P14	0.0001	0.0008	0.000	0.000
n-Pentadecane	P15	0.0001	0.0008	0.000	0.000
n-Hexadecane	P16	0.0001	0.0009	0.000	0.000
n-Heptadecane	P17	0.0001	0.0009	0.000	0.000
UnknownC18s	U18	0.0001	0.0010	0.000	0.000
TOTAL		100.00000	100.00000	8.9680	8.9195

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.730	14.650
BENZENE	0.0381	0.1138	LOW NET DRY REAL :	1331.7 /scf	1324.5 /scf
TOLUENE	0.0082	0.0289	NET WET REAL :	1308.5 /scf	1301.3 /scf
ETHYLBENZENE	0.0000	0.0000	HIGH GROSS DRY REAL :	1460.5 /scf	1452.6 /scf
XYLENES	0.0017	0.0068	GROSS WET REAL :	1435.1 /scf	1427.2 /scf
TOTAL BTEX	0.0480	0.1495	NET DRY REAL :	19337.2 /lb	19232.2 /lb
			GROSS DRY REAL :	21214.6 /lb	21099.3 /lb

RELATIVE DENSITY (AIR=1): 0.9024

COMPRESSIBILITY FACTOR : 0.99478

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

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

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