

CRUDE OIL ASSAY

PROJECT NO. :	201403015	ANALYSIS NO. :	04
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 6, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	1L GLASS JAR
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	TANK BATTERY @ 14:55		EMPACT
	BAILEY 7-26-8-60		
FIELD DATA		SAMPLE TEMP. :	58
SAMPLE PRES. :		AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; TK# 52116		

<u>SPECIFICATION</u>	<u>TEST METHOD</u>	<u>UNITS</u>	<u>RESULTS</u>
API GRAVITY		API 60/60	36.8
RVP @100 DEG F	D323	PSIG	9.4
TOTAL SULFUR	D2622	WT %	N/A
TOTAL CHLORIDE	D4929	ug/g	N/A
ORGANIC CHLORIDE	D4929	ug/g	N/A
FLASH POINT	D93	° F	N/A
HEATING VALUE	D4809	BTU/ LB	N/A
VISUAL APPEARANCE			RED/BROWN
<u>BS&W</u>	D96		
Crude Oil		VOL %	N/A
Water		VOL %	N/A
Emulsion		VOL %	N/A
Sediment		VOL %	N/A
<u>DISTILLATION:</u>	D86		
INITIAL POINT		DEG F	N/A
50%		DEG F	N/A
90%		DEG F	N/A
END POINT		DEG F	N/A
<u>DISTILLATION:</u>	@TEMP D445		
Average Centipoise	20°C		N/A
Average Centipoise	30°C		N/A
Average Centipoise	80°C		N/A
Kinetic Viscosity	20°C	cSt (mm2/s)	N/A
Kinetic Viscosity	30°C	cSt (mm2/s)	N/A
Kinetic Viscosity	80°C	cSt (mm2/s)	N/A

ND: NOT DETECTED

N/A: NO TEST PREFORMED FOR THIS PARAMETER

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303-637-0150

EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

MAIN PAGE

PROJECT NO. :	201403015	ANALYSIS NO. :	05
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 6, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	6843
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:30		EMPACT
	BAILEY 7-26-8-60		
FIELD DATA		SAMPLE TEMP. :	112
SAMPLE PRES. :	28	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	MOLE %	MASS %	VOL %
NITROGEN (AIR)	0.0361	0.0095	0.0086
CARBON DIOXIDE	0.0597	0.0247	0.0222
METHANE	0.2595	0.0391	0.0959
ETHANE	0.7802	0.2205	0.4554
PROPANE	2.2072	0.9147	1.3274
I-BUTANE	0.5136	0.2805	0.3666
N-BUTANE	2.4842	1.3568	1.7093
I-PENTANE	1.0747	0.7286	0.8592
N-PENTANE	1.7562	1.1907	1.3880
HEXANES PLUS	90.8286	95.2349	93.7674
TOTALS	100.0000	100.0000	100.0000

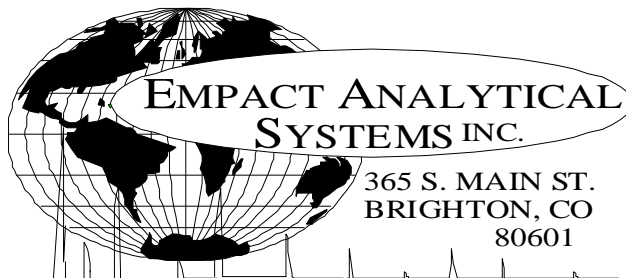
BTEX COMPONENTS	MOLE%	MASS%
BENZENE	1.5827	1.1617
TOLUENE	2.3294	2.0169
ETHYLBENZENE	0.6057	0.6043
XYLENE	1.8864	1.8821
TOTAL BTEX	6.4042	5.6650

(CALC: GPA STD 2145-94 & TP-17 @ 14.696 & 60 F)

	TOTAL SAMPLE	C6+ FRACTION
Specific Gravity (H2O=1) =	0.7357	0.7476 60/60
API Gravity =	60.83	57.77 60/60
Molecular Weight =	106.42	112.519
Absolute Density =	6.13	6.23 LBS/GAL
Heating Value Liq. Idl Gas=	125300	126856 BTU/GAL
Vapor/Liquid =	21.98	21.09 CUFT/GAL
Vapor Pressure =	27.42	2.01 PSIA @100 F

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

THIS DATA HAS BEEN ACQUIRED THROUGH APPLICATION OF CURRENT STATE-OF-THE-ART ANALYTICAL TECHNIQUES.
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303-637-0150

EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

E & P TANK / GLYCALC INFORMATION

PROJECT NO. :	201403015	ANALYSIS NO. :	05
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE :	MARCH 6, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	6843
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:30		EMPACT
	BAILEY 7-26-8-60		
FIELD DATA		SAMPLE TEMP. :	112
SAMPLE PRES. :	28	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	Mole %	Wt %	LV %			
CARBON DIOXIDE	0.0597	0.0247	0.0222			
NITROGEN (AIR)	0.0361	0.0095	0.0086			
METHANE	0.2595	0.0391	0.0959			
ETHANE	0.7802	0.2205	0.4554			
PROPANE	2.2072	0.9147	1.3274			
I-BUTANE	0.5136	0.2805	0.3666			
N-BUTANE	2.4842	1.3568	1.7093			
I-PENTANE	1.0747	0.7286	0.8592			
N-PENTANE	1.7562	1.1907	1.3880			
CYCLOPENTANE (N-C5)	2.0275	1.3362	1.2932			
N-HEXANE	8.5356	6.9131	7.6628			
CYCLOHEXANE (OTHER C6)	2.9675	2.3469	2.2039			
OTHER HEXANES	13.1322	10.5360	11.1509			
OTHER HEPTANES	12.9195	12.0813	12.5055			
METHYLCYCLOHEXANE (OTHER C7)	3.8127	3.5180	3.3410			
2,2,4 TRIMETHYLPENTANE	0.8121	0.7493	0.7316			
BENZENE	1.5827	1.1617	0.9681			
TOLUENE	2.3294	2.0169	1.6973			
ETHYLBENZENE	0.6057	0.6043	0.5085			
XYLENES	1.8864	1.8821	1.5861			
OTHER OCTANES	9.4248	10.1539	10.1411			
OCTANES PLUS	----	43.5215	----	55.3248	----	52.9447
NONANES	9.9108	11.8115	11.4512			
DECANES PLUS	20.8817	30.1237	28.5262			
SUB TOTAL	100.0000	100.0000	100.0000			
TOTAL	100.0000	100.0000	100.0000			

API Gravity	=	60.83	60/60
Vapor Pressure	=	27.42	PSIA & 100 F
Average Molecular Weight of Decanes plus	=	153.51	
Average Specific Gravity of Decanes plus	=	0.7800	

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.



303-637-0150

EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

BY CARBON NUMBER

PROJECT NO. : 201403015
 COMPANY NAME : CARRIZO OIL & GAS
 ACCOUNT NO. :
 PRODUCER :
 LEASE NO. :
 NAME/DESCRIP : SEPARATOR @ 14:30
 BAILEY 7-26-8-60

ANALYSIS NO. : 05
 ANALYSIS DATE: MARCH 6, 2014
 SAMPLE DATE : MARCH 4, 2014
 CYLINDER NO. : 6843
 SAMPLED BY : GALE MCENDREE
 EMPACT

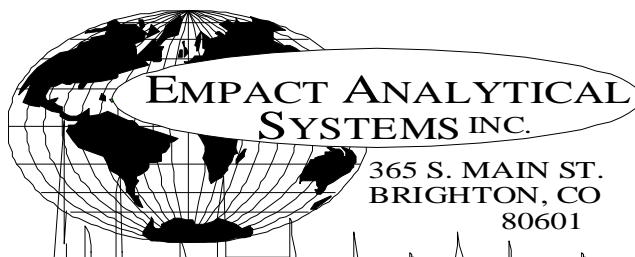
*****FIELD DATA*****

SAMPLE PRES.: 28
 VAPOR PRES. :
 COMMENTS : SPOT; NO PROBE

SAMPLE TEMP. : 112
 AMBIENT TEMP.:
 GRAVITY :

COMPONENT / CARBON NUMBER	MOLE%	MASS %	VOLUME %
NITROGEN	0.0361	0.0095	0.0086
CARBON DIOXIDE	0.0597	0.0247	0.0222
C1	0.2595	0.0391	0.0959
C2	0.7802	0.2205	0.4554
C3	2.2072	0.9147	1.3274
C4	2.9978	1.6373	2.0759
C5	4.8584	3.2555	3.5404
C6	26.2180	20.9577	21.9857
C7	19.0616	17.6162	17.5438
C8	12.7290	13.3896	12.9673
C9	9.9108	11.8115	11.4512
C10	8.5424	11.0277	10.4749
C11	5.2610	7.3483	6.8040
C12	3.5005	5.3418	5.1062
C13	1.8630	3.1451	3.0088
C14	1.2303	2.2937	2.2109
C15	0.4845	0.9671	0.9214
C16	0.0000	0.0000	0.0000
C17	0.0000	0.0000	0.0000
C18	0.0000	0.0000	0.0000
C19	0.0000	0.0000	0.0000
C20	0.0000	0.0000	0.0000
C21	0.0000	0.0000	0.0000
C22	0.0000	0.0000	0.0000
C23	0.0000	0.0000	0.0000
C24	0.0000	0.0000	0.0000
C25	0.0000	0.0000	0.0000
C26	0.0000	0.0000	0.0000
C27	0.0000	0.0000	0.0000
C28	0.0000	0.0000	0.0000
C29	0.0000	0.0000	0.0000
C30+	0.0000	0.0000	0.0000
Total	100.0000	100.0000	100.0000

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303-637-0150

EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

DHA COMPONENT LIST

PROJECT NO. :	201403015	ANALYSIS NO. :	05
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 6, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	6843
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:30		EMPACT
	BAILEY 7-26-8-60		
FIELD DATA		SAMPLE TEMP. :	112
SAMPLE PRES. :	28	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	PIANO #	MOLE %	MASS %	VOL %
Nitrogen	NHC	0.0361	0.0095	0.0086
Carbon Dioxide	NHC	0.0597	0.0247	0.0222
Methane	P1	0.2595	0.0391	0.0959
Ethane	P2	0.7802	0.2205	0.4554
Propane	P3	2.2072	0.9147	1.3274
i-Butane	I4	0.5136	0.2805	0.3666
n-Butane	P4	2.4842	1.3568	1.7093
2,2-Dimethylpropane	I5	0.0288	0.0195	0.0240
i-Pentane	I5	1.0459	0.7091	0.8352
n-Pentane	P5	1.7562	1.1907	1.3880
2,2-Dimethylbutane	I6	0.0438	0.0355	0.0399
Cyclopentane	N5	2.0275	1.3362	1.2932
2,3-Dimethylbutane	I6	0.4517	0.3658	0.4035
2-Methylpentane	I6	4.7813	3.8721	4.3311
3-Methylpentane	I6	2.6399	2.1379	2.3514
n-Hexane	P6	8.5356	6.9131	7.6628
2,2-Dimethylpentane	I7	0.0225	0.0212	0.0229
Methylcyclopentane	N6	5.2155	4.1247	4.0250
2,4-Dimethylpentane	I7	0.2524	0.2377	0.2586
Benzene	A6	1.5827	1.1617	0.9681
3,3-Dimethylpentane	I7	0.0247	0.0233	0.0246
Cyclohexane	N6	2.9675	2.3469	2.2039
2-Methylhexane	I7	1.2785	1.2038	1.2980
2,3-Dimethylpentane	I7	0.7137	0.6720	0.7039
1,1-Dimethylcyclopentane	N7	0.3772	0.3480	0.3372
3-Methylhexane	I7	1.7127	1.6127	1.7127
1c,3-Dimethylcyclopentane	N7	0.9015	0.8318	0.8165
1t,3-Dimethylcyclopentane	N7	0.8121	0.7493	0.7316
3-Ethylpentane	I7	0.1334	0.1256	0.1312
1t,2-Dimethylcyclopentane	N7	1.6695	1.5404	1.4989
2,2,4-Trimethylpentane	I8	0.0512	0.0550	0.0579
n-Heptane	P7	4.3561	4.1017	4.3848
1c,2-Dimethylcyclopentane	N7	0.1466	0.1353	0.1281
Methylcyclohexane	N7	3.8127	3.5180	3.3410
2,2-Dimethylhexane	I8	0.3738	0.4012	0.4216
Ethylcyclopentane	N7	0.5186	0.4785	0.4565
2,5-Dimethylhexane	I8	0.1018	0.1093	0.1152
2,2,3-Trimethylpentane	I8	0.0273	0.0293	0.0299
2,4-Dimethylhexane	I8	0.2096	0.2250	0.2359
1c,2t,4-Trimethylcyclopentane	N8	0.3606	0.3802	0.3642
3,3-Dimethylhexane	I8	0.0400	0.0429	0.0442
2,3,4-Trimethylpentane	I8	0.0837	0.0898	0.0913

2,3,3-Trimethylpentane	I8	0.0077	0.0083	0.0084
Toluene	A7	2.3294	2.0169	1.6973
2,3-Dimethylhexane	I8	0.1879	0.2017	0.2072
2-Methyl-3-ethylpentane	I8	0.0975	0.1047	0.1064
1,1,2-Trimethylcyclopentane	N8	0.0119	0.0125	0.0118
2-Methylheptane	I8	1.1391	1.2227	1.2786
4-Methylheptane	I8	0.3266	0.3506	0.3578
3-Methyl-3-ethylpentane	I8	0.0952	0.1022	0.1028
3,4-Dimethylhexane	I8	0.0764	0.0820	0.0833
1c,2c,4-Trimethylcyclopentane	N8	0.0324	0.0342	0.0324
1c,3-Dimethylcyclohexane	N8	0.0223	0.0235	0.0224
3-Methylheptane	I8	0.5078	0.5451	0.5651
1c,2t,3-Trimethylcyclopentane	N8	0.8212	0.8659	0.8221
3-Ethylhexane	I8	0.1922	0.2063	0.2116
1t,4-Dimethylcyclohexane	N8	0.3900	0.4112	0.3944
1,1-Dimethylcyclohexane	N8	0.1014	0.1069	0.1001
3t-Ethylmethylcyclopentane	N8	0.1735	0.1829	0.1745
2t-Ethylmethylcyclopentane	N8	0.1438	0.1516	0.1442
1,1-Methylethylcyclopentane	N8	0.4955	0.5225	0.4895
2,2,4-Trimethylhexane	I9	0.0422	0.0509	0.0520
1t,2-Dimethylcyclohexane	N8	0.5157	0.5438	0.5128
1t,3-Dimethylcyclohexane	N8	0.0124	0.0131	0.0122
n-Octane	P8	2.1806	2.3407	2.4360
1c,4-Dimethylcyclohexane	N8	0.4556	0.4804	0.4489
i-Propylcyclopentane	I8	0.0601	0.0634	0.0597
2,4,4-Trimethylhexane	I9	0.0161	0.0194	0.0197
2,2,3,4-Tetramethylpentane	I9	0.0166	0.0200	0.0203
2,3,4-Trimethylhexane	I9	0.0193	0.0233	0.0236
1c,2-Dimethylcyclohexane	N8	0.2174	0.2292	0.2106
2,3,5-Trimethylhexane	I9	0.0599	0.0722	0.0732
2,2-Dimethylheptane	I9	0.0146	0.0176	0.0181
1,1,4-Trimethylcyclohexane	N9	0.8495	1.0078	0.9552
2,2,3-Trimethylhexane	I9	0.3598	0.4337	0.4350
2,4-Dimethylheptane	I9	0.0807	0.0973	0.0995
4,4-Dimethylheptane	I9	0.0426	0.0513	0.0524
Ethylcyclohexane	N8	0.4800	0.5061	0.4701
n-Propylcyclopentane	N8	0.1919	0.2023	0.1906
1c,3c,5-Trimethylcyclohexane	N9	0.0425	0.0504	0.0478
2,5-Dimethylheptane	I9	0.0700	0.0844	0.0861
3,3-Dimethylheptane	I9	0.0769	0.0927	0.0946
3,5-Dimethylheptane	I9	0.0521	0.0628	0.0641
2,6-Dimethylheptane	I9	0.0478	0.0576	0.0594
1,1,3-Trimethylcyclohexane	N9	0.0874	0.1037	0.0983
Ethylbenzene	A8	0.6057	0.6043	0.5085
1c,2t,4t-Trimethylcyclohexane	N9	0.2191	0.2599	0.2416
2,3-Dimethylheptane	I9	0.0082	0.0099	0.0100
1,3-Dimethylbenzene (m-Xylene)	A8	0.6110	0.6096	0.5159
1,4-Dimethylbenzene (p-Xylene)	A8	0.6936	0.6920	0.5875
3,4-Dimethylheptane	I9	0.0469	0.0565	0.0565
3,4-Dimethylheptane (2)	I9	0.1459	0.1758	0.1758
4-Ethylheptane	I9	0.0356	0.0429	0.0439
4-Methyloctane	I9	0.2661	0.3207	0.3255
2-Methyloctane	I9	0.2966	0.3575	0.3665
1c,2t,4c-Trimethylcyclohexane	I9	0.0481	0.0580	0.0585
3-Ethylheptane	I9	0.0823	0.0992	0.0999
3-Methyloctane	I9	0.3265	0.3935	0.3994
3,3-Diethylpentane	I9	0.0528	0.0636	0.0617
1c,2t,3-Trimethylcyclohexane	N9	0.0728	0.0864	0.0803
1,1,2-Trimethylcyclohexane	N9	0.0310	0.0368	0.0342
1,2-Dimethylbenzene (o-Xylene)	A8	0.5818	0.5805	0.4827
i-Butylcyclopentane	N9	0.2154	0.2555	0.2394
UnknownC8s	U8	0.0528	0.0567	0.0590
n-Nonane	P9	1.6874	2.0338	2.0730
1,1-Methylethylcyclohexane	N9	0.2946	0.3551	0.3630
i-Propylbenzene	A9	0.2628	0.2968	0.2515
i-Propylcyclohexane	N9	0.0951	0.1128	0.1029
2,2-Dimethyloctane	I10	0.0698	0.0933	0.0923
2,4-Dimethyloctane	I10	0.0780	0.1043	0.1032
2,6-Dimethyloctane	I10	0.0140	0.0187	0.0191
2,5-Dimethyloctane	I10	0.0389	0.0520	0.0514
n-Butylcyclopentane	N9	0.3315	0.4370	0.4002
3,3-Dimethyloctane	I10	0.0204	0.0273	0.0270
n-Propylbenzene	A9	0.2537	0.2865	0.2428
3,6-Dimethyloctane	I10	0.2167	0.2897	0.2866

3-Methyl-5-ethylheptane	I10	0.3887	0.4685	0.4723
1,3-Methylethylbenzene	A9	0.3372	0.3809	0.3201
1,4-Methylethylbenzene	A9	0.1371	0.1549	0.1302
1,3,5-Trimethylbenzene	A9	0.1631	0.1842	0.1559
2,3-Dimethyloctane	I10	0.0648	0.0866	0.0857
5-Methylnonane	I10	0.2304	0.3080	0.3076
1,2-Methylethylbenzene	A9	0.4746	0.5361	0.4481
2-Methylnonane	I10	0.0738	0.0987	0.0994
3-Ethyl-octane	I10	0.0726	0.0971	0.0961
3-Methylnonane	I10	0.1474	0.1971	0.1966
1,2,4-Trimethylbenzene	A9	0.0327	0.0369	0.0308
t-Butylbenzene	A10	0.2636	0.3325	0.2810
i-Butylcyclohexane	N10	0.1994	0.2628	0.2369
1t-Methyl-2-n-propylcyclohexane	I10	0.0344	0.0415	0.0418
i-Butylbenzene	A10	0.0651	0.0821	0.0704
sec-Butylbenzene	A10	0.0272	0.0343	0.0291
UnknownC9s	U9	1.9215	2.3159	2.3606
n-Decane	P10	1.1453	1.5313	1.5346
1,2,3-Trimethylbenzene	A9	0.1942	0.2193	0.1796
1,3-Methyl-i-propylbenzene	A10	0.0994	0.1123	0.0939
1,4-Methyl-i-propylbenzene	A10	0.0437	0.0494	0.0413
Sec-Butylcyclohexane	N10	0.3212	0.4234	0.3812
1,2-Methyl-i-propylbenzene	A10	0.1396	0.1761	0.1471
3-Ethyl-nonane	I10	0.0446	0.0596	0.0600
1,3-Diethylbenzene	A10	0.1343	0.1694	0.1435
1,3-Methyl-n-propylbenzene	A10	0.0258	0.0325	0.0276
1,4-Diethylbenzene	A10	0.1435	0.1810	0.1537
1,4-Methyl-n-propylbenzene	A10	0.0318	0.0401	0.0342
n-Butylbenzene	A10	0.0806	0.1017	0.0864
1,3-Dimethyl-5-ethylbenzene	A10	0.0595	0.0750	0.0635
1,2-Diethylbenzene	A10	0.0925	0.1167	0.0971
1,2-Methyl-n-propylbenzene	A10	0.0886	0.1118	0.0937
1,4-Dimethyl-2-ethylbenzene	A10	0.1182	0.1491	0.1244
1,3-Dimethyl-4-ethylbenzene	A10	0.0424	0.0535	0.0447
1,2-Dimethyl-4-ethylbenzene	A10	0.1938	0.2444	0.2046
1,3-Dimethyl-2-ethylbenzene	A10	0.1447	0.1825	0.1501
1t,2c,4-Trimethylcyclopentane	A10	0.4305	0.4539	0.4442
1,2-Dimethyl-3-ethylbenzene	A10	0.1133	0.1429	0.1173
1,2-Ethyl-i-propylbenzene	A10	0.1116	0.1408	0.1176
1,4-Methyl-t-butylbenzene	A11	0.1724	0.2174	0.1815
UnknownC10s	U10	2.5644	3.4287	3.4360
n-Undecane	P11	0.9511	1.3970	1.3806
1,4-Ethyl-i-propylbenzene	A11	0.0690	0.0870	0.0727
1,2,4,5-Tetramethylbenzene	A11	0.2065	0.2605	0.2153
1,2-Methyl-n-butylbenzene	A11	0.1099	0.1386	0.1157
1,2,3,5-Tetramethylbenzene	A11	0.1525	0.1923	0.1582
1,2-Methyl-t-butylbenzene	A11	0.1505	0.1898	0.1585
5-Methylindan	A11	0.0317	0.0507	0.0496
4-Methylindan	A11	0.0261	0.0418	0.0409
1,2-Ethyl-n-propylbenzene	A11	0.1657	0.2090	0.1745
2-Methylindan	A11	0.0949	0.1519	0.1485
1,3-Methyl-n-butylbenzene	A11	0.0859	0.1083	0.0904
sec-Pentylbenzene	A11	0.0972	0.1226	0.1024
n-Pentylbenzene	A11	0.0730	0.1017	0.0867
1t-M-2-(4MP)cyclopentane	P12	0.0597	0.0956	0.0935
1,2-Di-n-propylbenzene	A11	0.1361	0.1717	0.1434
1,4-Di-i-propylbenzene	A11	0.1639	0.2067	0.1726
Tetrahydronaphthalene	A10	0.1212	0.1529	0.1277
t-Decahydronaphthalene	A10	0.1238	0.1562	0.1304
Naphthalene	A10	0.1229	0.1480	0.1236
1-t-Butyl-3,5-dimethylbenzene	A12	0.0793	0.1000	0.0835
1,4-Ethyl-t-butylbenzene	A11	0.1113	0.1404	0.1172
UnknownC11s	U11	2.0300	2.9818	2.9467
n-Dodecane	P12	0.7512	1.2024	1.1752
1,3-Di-n-propylbenzene	A12	0.0782	0.0986	0.0823
1,3,5-Triethylbenzene	A12	0.0719	0.0812	0.0687
1,2,4-Triethylbenzene	A12	0.2979	0.3365	0.2811
1,4-Methyl-n-pentylbenzene	A12	0.0836	0.1054	0.0880
n-Hexylbenzene	A12	0.0696	0.1061	0.0906
1,2,3,4,5-Pentamethylbenzene	A13	0.1750	0.2207	0.1843
2-Methylnaphthalene	A11	0.2098	0.2804	0.2342
1-Methylnaphthalene	A11	0.2235	0.2987	0.2144
UnknownC12s	U12	2.0091	3.2160	3.1433
n-Tridecane	P13	0.5611	0.9721	0.9389

UnknownC13s	U13	1.1269	1.9523	1.8856
n-Tetradecane	P14	0.3701	0.6900	0.6651
UnknownC14s	U14	0.8602	1.6037	1.5458
UnknownC15s	U15	<u>0.4845</u>	<u>0.9671</u>	<u>0.9214</u>
<u>TOTAL</u>		<u>100.0000</u>	<u>100.0000</u>	<u>100.0000</u>

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303-637-0150

EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PROJECT NO. :	201403015	ANALYSIS NO. :	06
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 12, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 11, 2014
PRODUCER :		CYLINDER NO. :	0958
LEASE NO. :		SAMPLED BY :	GALE MCENDREE-EMPACT
NAME/DESCRIP :	SALES GAS @ 10:15 BAILEY 7-26-8-60		
FIELD DATA		SAMPLE TEMP. :	77
SAMPLE PRES. :	118	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; PROBE; LENGTH OF H2S STAIN @ 1-7=5 PPM @ 14:45 RESAMPLE-EMPACT		

COMPONENT	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
ALCOHOLS	0.0002	0.0005		
HELIUM	0.01	0.00	---	---
HYDROGEN	0.00	0.00	---	---
OXYGEN/ARGON	0.01	0.01	---	---
NITROGEN	0.90	1.10	---	---
CARBON DIOXIDE	2.60	4.98	---	---
METHANE	71.95970	50.20510	---	---
ETHANE	12.0172	15.7146	3.2089	3.2264
PROPANE	7.9896	15.3215	2.1970	2.2090
I-BUTANE	0.7396	1.8695	0.2412	0.2425
N-BUTANE	2.3553	5.9535	0.7417	0.7457
I-PENTANE	0.4429	1.3861	0.1591	0.1600
N-PENTANE	0.5027	1.5773	0.1822	0.1832
HEXANES PLUS	0.4728	1.8819	0.1880	0.1890
TOTALS	100.00000	100.00000	6.9181	6.9558

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.650	14.730
BENZENE	0.0162	0.0550	LOW NET DRY REAL :	1184.0 /scf	1190.4 /scf
TOLUENE	0.0095	0.0380	NET WET REAL :	1163.3 /scf	1169.7 /scf
ETHYLBENZENE	0.0009	0.0042	HIGH GROSS DRY REAL :	1302.7 /scf	1309.8 /scf
XYLENES	0.0022	0.0101	GROSS WET REAL :	1279.9 /scf	1287.0 /scf
TOTAL BTEX	0.0288	0.1073	NET DRY REAL :	19557.4 /lb	19664.2 /lb
			GROSS DRY REAL :	21520.9 /lb	21638.4 /lb

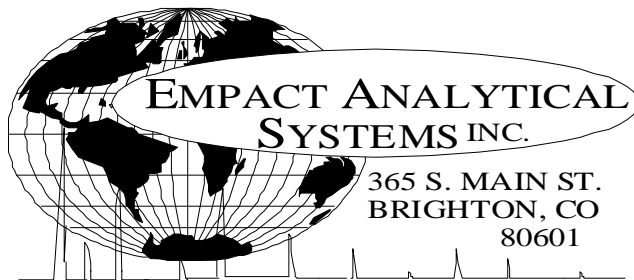
RELATIVE DENSITY (AIR=1):	0.7931
COMPRESSIBILITY FACTOR :	0.99598

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

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

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303-637-0150

EXTENDED NATURAL GAS ANALYSIS (*DHA)

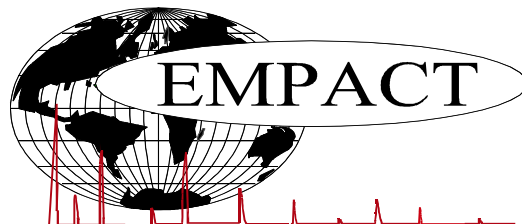
GLYCALC INFORMATION

PROJECT NO. : 201403015 ANALYSIS NO. : 06
COMPANY NAME : CARRIZO OIL & GAS ANALYSIS DATE: MARCH 12, 2014
ACCOUNT NO. : SAMPLE DATE : MARCH 11, 2014
PRODUCER : CYLINDER NO. : 0958
LEASE NO. : SAMPLED BY : GALE MCENDREE-EMPACT
NAME/DESCRIP : SALES GAS @ 10:15
BAILEY 7-26-8-60

FIELD DATA
SAMPLE PRES. : 118 SAMPLE TEMP. : 77
VAPOR PRES. : AMBIENT TEMP.:
COMMENTS : GRAVITY :
SPOT; PROBE; LENGTH OF H2S STAIN @ 1-7=5 PPM @ 14:45
RESAMPLE-EMPACT

Componet	Mole %	Wt %
Helium	0.01	0.00
Carbon Dioxide	2.60	4.98
Carbon Monoxide	0.00	0.00
Nitrogen	0.90	1.10
Methane	71.95970	50.20510
Ethane	12.0172	15.7146
Propane	7.9896	15.3215
Isobutane	0.7396	1.8695
n-Butane	2.3553	5.9535
Isopentane	0.4021	1.2617
n-Pentane	0.5027	1.5773
Cyclopentane	0.0408	0.1244
n-Hexane	0.0992	0.3718
Cyclohexane	0.0252	0.0922
Other Hexanes	0.1763	0.6558
Heptanes	0.0825	0.3578
Methycyclohexane	0.0192	0.0820
2,2,4 Trimethylpentane	0.0002	0.0010
Benzene	0.0162	0.0550
Toluene	0.0095	0.0380
Ethylbenzene	0.0009	0.0042
Xylenes	0.0022	0.0101
C8+ Heavies	0.0414	0.2140
Subtotal	99.98980	99.98950
Oxygen/Argon	0.01	0.01
Alcohols	0.0002	0.0005
Total	100.00000	100.00000

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EXTENDED NATURAL GAS ANALYSIS (*DHA)

DHA COMPONENT LIST

PROJECT NO. :	201403015	ANALYSIS NO. :	06
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 12, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 11, 2014
PRODUCER :		CYLINDER NO. :	0958
LEASE NO. :		SAMPLED BY :	GALE MCENDREE-EMPACT
NAME/DESCRIP :	SALES GAS @ 10:15 BAILEY 7-26-8-60		
FIELD DATA		SAMPLE TEMP. :	77
SAMPLE PRES. :	118	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; PROBE; LENGTH OF H2S STAIN @ 1-7=5 PPM @ 14:45 RESAMPLE-EMPACT		

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
Helium	---	0.01	0.00	---	---
Hydrogen	---	0.00	0.00	---	---
Oxygen/Argon	---	0.01	0.01	---	---
Nitrogen	---	0.90	1.10	---	---
Carbon Dioxide	---	2.60	4.98	---	---
Methane	P1	71.95970	50.20510	---	---
Ethane	P2	12.0172	15.7146	3.209	3.226
Propane	P3	7.9896	15.3215	2.197	2.209
i-Butane	I4	0.7396	1.8695	0.241	0.243
n-Butane	P4	2.3553	5.9535	0.742	0.746
2,2-Dimethylpropane	I5	0.0019	0.0060	0.001	0.001
Ethanol	X2	0.0001	0.0002	0.000	0.000
i-Pentane	I5	0.4002	1.2557	0.146	0.147
n-Pentane	P5	0.5027	1.5773	0.182	0.183
t-Butanol	X4	0.0001	0.0003	0.000	0.000
2,2-Dimethylbutane	I6	0.0012	0.0045	0.000	0.000
Cyclopentane	N5	0.0408	0.1244	0.012	0.012
2,3-Dimethylbutane	I6	0.0067	0.0251	0.003	0.003
2-Methylpentane	I6	0.0748	0.2803	0.031	0.031
3-Methylpentane	I6	0.0379	0.1420	0.015	0.015
n-Hexane	P6	0.0992	0.3718	0.041	0.041
2,2-Dimethylpentane	I7	0.0004	0.0017	0.000	0.000
Methylcyclopentane	N6	0.0557	0.2039	0.020	0.020
2,4-Dimethylpentane	I7	0.0024	0.0104	0.001	0.001
2,2,3-Trimethylbutane	I7	0.0001	0.0004	0.000	0.000
Benzene	A6	0.0162	0.0550	0.005	0.005
3,3-Dimethylpentane	I7	0.0002	0.0009	0.000	0.000
Cyclohexane	N6	0.0252	0.0922	0.009	0.009
2-Methylhexane	I7	0.0098	0.0427	0.005	0.005
2,3-Dimethylpentane	I7	0.0046	0.0200	0.002	0.002
1,1-Dimethylcyclopentane	N7	0.0027	0.0115	0.001	0.001
3-Methylhexane	I7	0.0112	0.0488	0.005	0.005
1c,3-Dimethylcyclopentane	N7	0.0062	0.0265	0.003	0.003
1t,3-Dimethylcyclopentane	N7	0.0055	0.0235	0.003	0.003
3-Ethylpentane	I7	0.0008	0.0035	0.000	0.000
1t,2-Dimethylcyclopentane	N7	0.0107	0.0457	0.005	0.005

2,2,4-Trimethylpentane	I8	0.0002	0.0010	0.000	0.000
n-Heptane	P7	0.0231	0.1007	0.011	0.011
1c,2-Dimethylcyclopentane	N7	0.0008	0.0034	0.000	0.000
Methylcyclohexane	N7	0.0192	0.0820	0.008	0.008
1,1,3-Trimethylcyclopentane	N7	0.0016	0.0078	0.001	0.001
Ethylcyclopentane	N7	0.0024	0.0103	0.001	0.001
2,5-Dimethylhexane	I8	0.0004	0.0020	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0001	0.0005	0.000	0.000
2,4-Dimethylhexane	I8	0.0008	0.0040	0.000	0.000
1c,2t,4-Trimethylcyclopentane	N8	0.0016	0.0078	0.001	0.001
3,3-Dimethylhexane	I8	0.0001	0.0005	0.000	0.000
1t,2c,4-Trimethylcyclopentane	N8	0.0017	0.0083	0.001	0.001
2,3,4-Trimethylpentane	I8	0.0003	0.0015	0.000	0.000
2,3,3-Trimethylpentane	I8	0.0001	0.0005	0.000	0.000
Toluene	A7	0.0095	0.0380	0.003	0.003
2,3-Dimethylhexane	I8	0.0006	0.0030	0.000	0.000
2-Methyl-3-ethylpentane	I8	0.0005	0.0025	0.000	0.000
2-Methylheptane	I8	0.0037	0.0184	0.002	0.002
4-Methylheptane	I8	0.0010	0.0050	0.001	0.001
3-Methyl-3-ethylpentane	I8	0.0001	0.0005	0.000	0.000
3,4-Dimethylhexane	I8	0.0002	0.0010	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0001	0.0005	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0001	0.0005	0.000	0.000
3-Methylheptane	I8	0.0018	0.0090	0.001	0.001
1c,2t,3-Trimethylcyclopentane	N8	0.0025	0.0122	0.001	0.001
3-Ethylhexane	I8	0.0002	0.0010	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0009	0.0044	0.000	0.000
1,1-Dimethylcyclohexane	N8	0.0003	0.0015	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0005	0.0024	0.000	0.000
2t-Ethylmethylcyclopentane	N8	0.0004	0.0020	0.000	0.000
1,1-Methylethylcyclopentane	N8	0.0013	0.0063	0.001	0.001
2,2,4-Trimethylhexane	I9	0.0001	0.0006	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0013	0.0063	0.001	0.001
n-Octane	P8	0.0054	0.0268	0.003	0.003
1c,4-Dimethylcyclohexane	N8	0.0005	0.0024	0.000	0.000
i-Propylcyclopentane	I8	0.0001	0.0005	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0003	0.0016	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0003	0.0015	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0014	0.0077	0.001	0.001
2,2,3-Trimethylhexane	I9	0.0006	0.0033	0.000	0.000
2,4-Dimethylheptane	I9	0.0002	0.0011	0.000	0.000
Ethylcyclohexane	N8	0.0008	0.0039	0.000	0.000
n-Propylcyclopentane	N8	0.0003	0.0015	0.000	0.000
2,5-Dimethylheptane	I9	0.0001	0.0006	0.000	0.000
3,3-Dimethylheptane	I9	0.0001	0.0006	0.000	0.000
3,5-Dimethylheptane	I9	0.0001	0.0006	0.000	0.000
2,6-Dimethylheptane	I9	0.0001	0.0006	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0003	0.0016	0.000	0.000
Ethylbenzene	I8	0.0009	0.0042	0.000	0.000
1c,2t,4t-Trimethylcyclohexane	N9	0.0003	0.0016	0.000	0.000
2,3-Dimethylheptane	I9	0.0007	0.0039	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0012	0.0055	0.000	0.000
1,4-Dimethylbenzene (p-Xylene)	A8	0.0003	0.0014	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0002	0.0011	0.000	0.000
4-Methyloctane	I9	0.0003	0.0016	0.000	0.000
2-Methyloctane	I9	0.0004	0.0022	0.000	0.000
1c,2t,3-Trimethylcyclohexane	N9	0.0001	0.0006	0.000	0.000
3-Ethylheptane	I9	0.0001	0.0006	0.000	0.000
3-Methyloctane	I9	0.0004	0.0022	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0007	0.0032	0.000	0.000

i-Butylcyclopentane	N9	0.0003	0.0016	0.000	0.000
UnknownC8s	U8	0.0001	0.0005	0.000	0.000
n-Nonane	P9	0.0013	0.0073	0.001	0.001
1,1-Methylethylcyclohexane	N9	0.0002	0.0011	0.000	0.000
i-Propylbenzene	A9	0.0002	0.0010	0.000	0.000
i-Propylcyclohexane	N9	0.0001	0.0006	0.000	0.000
2,2-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
2,4-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
n-Butylcyclopentane	N9	0.0003	0.0016	0.000	0.000
3,3-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
n-Propylbenzene	A9	0.0002	0.0010	0.000	0.000
3,6-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
3-Methyl-5-ethylheptane	I10	0.0004	0.0025	0.000	0.000
1,3-Methylethylbenzene	A9	0.0002	0.0010	0.000	0.000
1,4-Methylethylbenzene	A9	0.0001	0.0005	0.000	0.000
1,3,5-Trimethylbenzene	A9	0.0001	0.0005	0.000	0.000
5-Methylnonane	I10	0.0001	0.0006	0.000	0.000
1,2-Methylethylbenzene	A9	0.0001	0.0005	0.000	0.000
2-Methylnonane	I10	0.0001	0.0006	0.000	0.000
3-Ethyloctane	I10	0.0001	0.0006	0.000	0.000
3-Methylnonane	I10	0.0001	0.0006	0.000	0.000
1,2,4-Trimethylbenzene	A9	0.0001	0.0005	0.000	0.000
t-Butylbenzene	A10	0.0002	0.0012	0.000	0.000
i-Butylcyclohexane	N10	0.0001	0.0006	0.000	0.000
UnknownC9s	U9	0.0010	0.0056	0.001	0.001
n-Decane	P10	0.0004	0.0025	0.000	0.000
1,2,3-Trimethylbenzene	A9	0.0001	0.0005	0.000	0.000
Sec-Butylcyclohexane	A10	0.0001	0.0006	0.000	0.000
UnknownC10s	U10	0.0008	0.0050	0.000	0.000
n-Undecane	P11	0.0002	0.0013	0.000	0.000
UnknownC11s	U11	0.0001	0.0007	0.000	0.000
n-Dodecane	P12	0.0001	0.0007	0.000	0.000
TOTAL		100.00000	100.00000	6.9181	6.9558

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.650	14.730
BENZENE	0.0162	0.0550	LOW NET DRY REAL :	1184.0 /scf	1190.4 /scf
TOLUENE	0.0095	0.0380	NET WET REAL :	1163.3 /scf	1169.7 /scf
ETHYLBENZENE	0.0009	0.0042	HIGH GROSS DRY REAL :	1302.7 /scf	1309.8 /scf
XYLENES	0.0022	0.0101	GROSS WET REAL :	1279.9 /scf	1287.0 /scf
TOTAL BTEX	0.0288	0.1073	NET DRY REAL :	19557.4 /lb	19664.2 /lb
			GROSS DRY REAL :	21520.9 /lb	21638.4 /lb

RELATIVE DENSITY (AIR=1): 0.7931
COMPRESSIBILITY FACTOR : 0.99598

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

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

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