

# CRUDE OIL ASSAY

PROJECT NO. :	201403015	ANALYSIS NO. :	01
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 @ 15:50		EMPACT
	BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	85
SAMPLE PRES. :		AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; TK# 52620		

<u>SPECIFICATION</u>	<u>TEST METHOD</u>	<u>UNITS</u>	<u>RESULTS</u>
API GRAVITY		API 60/60	36.1
RVP @100 DEG F	D323	PSIG	7.4
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&amp;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		
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

*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.*



303-637-0150

**EXTENDED NATURAL GAS LIQUID ANALYSIS (\*DHA)**

**MAIN PAGE**

PROJECT NO. :	201403015	ANALYSIS NO. :	02
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 5, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	13005
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:00		EMPACT
	BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	144
SAMPLE PRES. :	28 PSI	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	MOLE %	MASS %	VOL %
NITROGEN (AIR)	0.0382	0.0099	0.0090
CARBON DIOXIDE	0.0214	0.0087	0.0079
METHANE	0.0648	0.0096	0.0237
ETHANE	0.4273	0.1187	0.2463
PROPANE	1.7183	0.6998	1.0203
I-BUTANE	0.4589	0.2463	0.3234
N-BUTANE	2.3066	1.2380	1.5670
I-PENTANE	1.0620	0.7076	0.8382
N-PENTANE	1.7205	1.1464	1.3427
HEXANES PLUS	92.1820	95.8150	94.6215
TOTALS	100.0000	100.0000	100.0000

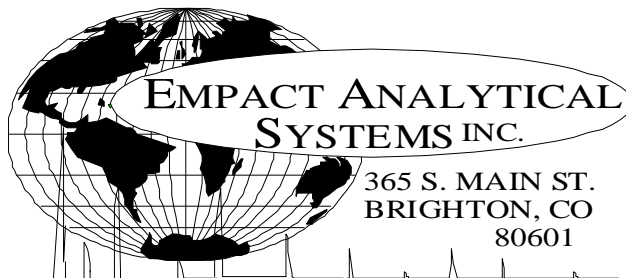
BTEX COMPONENTS	MOLE%	MASS%
BENZENE	1.4776	1.0659
TOLUENE	2.4294	2.0672
ETHYLBENZENE	0.6456	0.6330
XYLENE	2.0414	2.0016
TOTAL BTEX	6.5940	5.7677

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

	TOTAL SAMPLE	C6+ FRACTION
Specific Gravity (H2O=1) =	0.7398	0.7491 60/60
API Gravity =	59.77	57.39 60/60
Molecular Weight =	108.28	113.48
Absolute Density =	6.17	6.24 LBS/GAL
Heating Value Liq. Idl Gas=	125540	127077 BTU/GAL
Vapor/Liquid =	21.64	20.94 CUFT/GAL
Vapor Pressure =	13.75	1.93 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. :	02
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 5, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	13005
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:00		EMPACT
	BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	144
SAMPLE PRES. :	28 PSI	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	Mole %	Wt %	LV %			
CARBON DIOXIDE	0.0214	0.0087	0.0079			
NITROGEN (AIR)	0.0382	0.0099	0.0090			
METHANE	0.0648	0.0096	0.0237			
ETHANE	0.4273	0.1187	0.2463			
PROPANE	1.7183	0.6998	1.0203			
I-BUTANE	0.4589	0.2463	0.3234			
N-BUTANE	2.3066	1.2380	1.5670			
I-PENTANE	1.0620	0.7076	0.8382			
N-PENTANE	1.7205	1.1464	1.3427			
CYCLOPENTANE (N-C5)	1.9603	1.2696	1.2345			
N-HEXANE	8.2688	6.5800	7.3272			
CYCLOHEXANE (OTHER C6)	2.8147	2.1876	2.0640			
OTHER HEXANES	12.7470	10.0520	10.6971			
OTHER HEPTANES	12.4481	11.4414	11.9054			
METHYLCYCLOHEXANE (OTHER C7)	3.8285	3.4716	3.3124			
2,2,4 TRIMETHYLPENTANE	0.7740	0.7019	0.6886			
BENZENE	1.4776	1.0659	0.8924			
TOLUENE	2.4294	2.0672	1.7478			
ETHYLBENZENE	0.6456	0.6330	0.5351			
XYLENES	2.0414	2.0016	1.6941			
OTHER OCTANES	9.8303	10.3987	10.4335			
OCTANES PLUS	----	46.2076	----	57.6797	----	55.4407
NONANES	10.3292	12.0935	11.7819			
DECANES PLUS	22.5871	31.8510	30.3075			
SUB TOTAL	100.0000	100.0000	100.0000			
TOTAL	100.0000	100.0000	100.0000			

API Gravity	=	59.77	60/60
Vapor Pressure	=	13.75	PSIA & 100 F
Average Molecular Weight of Decanes plus	=	152.69	
Average Specific Gravity of Decanes plus	=	0.7770	

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

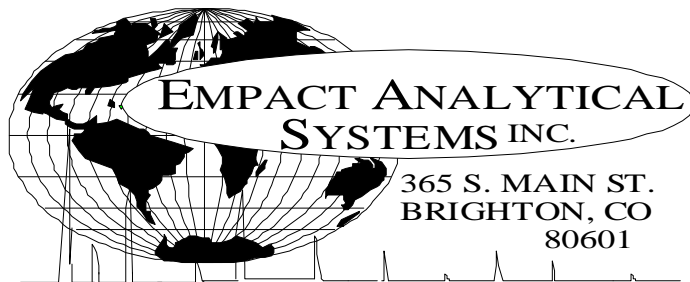
**EXTENDED NATURAL GAS LIQUID ANALYSIS (\*DHA)**

**BY CARBON NUMBER**

PROJECT NO. :	201403015	ANALYSIS NO. :	02
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 5, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	13005
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:00		EMPACT
	BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	144
SAMPLE PRES. :	28 PSI	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT / CARBON NUMBER	MOLE%	MASS %	VOLUME %
NITROGEN	0.0382	0.0099	0.0090
CARBON DIOXIDE	0.0214	0.0087	0.0079
C1	0.0648	0.0096	0.0237
C2	0.4273	0.1187	0.2463
C3	1.7183	0.6998	1.0203
C4	2.7655	1.4843	1.8904
C5	4.7428	3.1236	3.4154
C6	25.3081	19.8855	20.9807
C7	18.7060	16.9802	16.9656
C8	13.2913	13.7352	13.3513
C9	10.3292	12.0935	11.7819
C10	9.5509	12.1480	11.6187
C11	5.7652	7.8939	7.3318
C12	3.7828	5.6726	5.4487
C13	1.9153	3.1811	3.0588
C14	1.0043	1.8400	1.7818
C15	0.5686	1.1154	1.0677
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
<b>Total</b>	<b>100.0000</b>	<b>100.0000</b>	<b>100.0000</b>

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303-637-0150

**EXTENDED NATURAL GAS LIQUID ANALYSIS (\*DHA)**

**DHA COMPONENT LIST**

PROJECT NO. :	201403015	ANALYSIS NO. :	02
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 5, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	13005
LEASE NO. :		SAMPLED BY :	GALE MCENDREE
NAME/DESCRIP :	SEPARATOR @ 14:00		IMPACT
	BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	144
SAMPLE PRES. :	28 PSI	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	PIANO #	MOLE %	MASS %	VOL %
Nitrogen	NHC	0.0382	0.0099	0.0090
Carbon Dioxide	NHC	0.0214	0.0087	0.0079
Methane	P1	0.0648	0.0096	0.0237
Ethane	P2	0.4273	0.1187	0.2463
Propane	P3	1.7183	0.6998	1.0203
i-Butane	I4	0.4589	0.2463	0.3234
n-Butane	P4	2.3066	1.2380	1.5670
2,2-Dimethylpropane	I5	0.0219	0.0146	0.0181
i-Pentane	I5	1.0401	0.6930	0.8201
n-Pentane	P5	1.7205	1.1464	1.3427
2,2-Dimethylbutane	I6	0.0470	0.0374	0.0423
Cyclopentane	N5	1.9603	1.2696	1.2345
2,3-Dimethylbutane	I6	0.4563	0.3632	0.4025
2-Methylpentane	I6	4.6851	3.7287	4.1902
3-Methylpentane	I6	2.5754	2.0497	2.2650
n-Hexane	P6	8.2688	6.5800	7.3272
Methylcyclopentane	N6	4.9832	3.8730	3.7971
2,4-Dimethylpentane	I7	0.2515	0.2327	0.2543
Benzene	A6	1.4776	1.0659	0.8924
3,3-Dimethylpentane	I7	0.0148	0.0137	0.0145
Cyclohexane	N6	2.8147	2.1876	2.0640
2-Methylhexane	I7	1.2275	1.1359	1.2305
2,3-Dimethylpentane	I7	0.7221	0.6682	0.7032
1,1-Dimethylcyclopentane	N7	0.3587	0.3253	0.3167
3-Methylhexane	I7	1.6381	1.5158	1.6173
1c,3-Dimethylcyclopentane	N7	0.8465	0.7676	0.7570
1t,3-Dimethylcyclopentane	N7	0.7740	0.7019	0.6886
3-Ethylpentane	I7	0.1266	0.1171	0.1229
1t,2-Dimethylcyclopentane	N7	1.5188	1.3772	1.3464
2,2,4-Trimethylpentane	I8	0.0984	0.1038	0.1097
n-Heptane	P7	4.2579	3.9400	4.2317
1c,2-Dimethylcyclopentane	N7	0.1594	0.1445	0.1375
Methylcyclohexane	N7	3.8285	3.4716	3.3124
2,2-Dimethylhexane	I8	0.3744	0.3950	0.4171
Ethylcyclopentane	N7	0.5143	0.4664	0.4471

2,5-Dimethylhexane	I8	0.1033	0.1090	0.1154
2,2,3-Trimethylpentane	I8	0.0289	0.0305	0.0313
2,4-Dimethylhexane	I8	0.1931	0.2037	0.2146
1c,2t,4-Trimethylcyclopentane	N8	0.3607	0.3738	0.3598
3,3-Dimethylhexane	I8	0.0501	0.0529	0.0548
2,3,4-Trimethylpentane	I8	0.0851	0.0898	0.0917
2,3,3-Trimethylpentane	I8	0.0130	0.0137	0.0139
Toluene	A7	2.4294	2.0672	1.7478
2,3-Dimethylhexane	I8	0.1798	0.1897	0.1958
2-Methyl-3-ethylpentane	I8	0.1178	0.1243	0.1269
1,1,2-Trimethylcyclopentane	N8	0.0072	0.0075	0.0071
2-Methylheptane	I8	1.1758	1.2404	1.3032
4-Methylheptane	I8	0.3470	0.3661	0.3753
3-Methyl-3-ethylpentane	I8	0.0921	0.0972	0.0982
3,4-Dimethylhexane	I8	0.0631	0.0666	0.0680
1c,2c,4-Trimethylcyclopentane	N8	0.0335	0.0347	0.0331
1c,3-Dimethylcyclohexane	N8	0.0163	0.0169	0.0162
3-Methylheptane	I8	0.5298	0.5589	0.5821
1c,2t,3-Trimethylcyclopentane	N8	0.8649	0.8963	0.8549
3-Ethylhexane	I8	0.1875	0.1978	0.2038
1t,4-Dimethylcyclohexane	N8	0.3864	0.4004	0.3859
1,1-Dimethylcyclohexane	N8	0.1148	0.1190	0.1120
3t-Ethylmethylcyclopentane	N8	0.1797	0.1862	0.1785
2t-Ethylmethylcyclopentane	N8	0.1515	0.1570	0.1501
1,1-Methylethylcyclopentane	N8	0.5150	0.5337	0.5023
2,2,4-Trimethylhexane	I9	0.0439	0.0520	0.0534
1t,2-Dimethylcyclohexane	N8	0.5191	0.5379	0.5096
1t,3-Dimethylcyclohexane	N8	0.0148	0.0153	0.0143
UnknownC7s	U7	0.0379	0.0351	0.0377
n-Octane	P8	2.3071	2.4338	2.5448
1c,4-Dimethylcyclohexane	N8	0.4803	0.4977	0.4673
i-Propylcyclopentane	I8	0.0433	0.0449	0.0425
2,4,4-Trimethylhexane	I9	0.0260	0.0308	0.0314
2,2,3,4-Tetramethylpentane	I9	0.0145	0.0172	0.0176
2,3,4-Trimethylhexane	I9	0.0254	0.0301	0.0306
1c,2-Dimethylcyclohexane	N8	0.2346	0.2431	0.2244
2,3,5-Trimethylhexane	I9	0.0148	0.0175	0.0178
2,2-Dimethylheptane	I9	0.0181	0.0214	0.0221
1,1,4-Trimethylcyclohexane	N9	0.9046	1.0546	1.0043
2,2,3-Trimethylhexane	I9	0.3672	0.4349	0.4382
2,4-Dimethylheptane	I9	0.0700	0.0829	0.0852
4,4-Dimethylheptane	I9	0.0378	0.0448	0.0460
Ethylcyclohexane	N8	0.5116	0.5302	0.4948
n-Propylcyclopentane	N8	0.2024	0.2097	0.1985
1c,3c,5-Trimethylcyclohexane	N9	0.0397	0.0463	0.0441
2,5-Dimethylheptane	I9	0.0635	0.0752	0.0771
3,3-Dimethylheptane	I9	0.0710	0.0841	0.0862
3,5-Dimethylheptane	I9	0.0530	0.0628	0.0644
2,6-Dimethylheptane	I9	0.0468	0.0554	0.0574
1,1,3-Trimethylcyclohexane	N9	0.0942	0.1098	0.1046
Ethylbenzene	A8	0.6456	0.6330	0.5351
1c,2t,4t-Trimethylcyclohexane	N9	0.2169	0.2529	0.2362
2,3-Dimethylheptane	I9	0.0134	0.0159	0.0161
1,3-Dimethylbenzene (m-Xylene)	A8	0.7050	0.6912	0.5877
1,4-Dimethylbenzene (p-Xylene)	A8	0.6789	0.6657	0.5678
3,4-Dimethylheptane	I9	0.0691	0.0818	0.0822
3,4-Dimethylheptane (2)	I9	0.1204	0.1426	0.1433
4-Ethylheptane	I9	0.0234	0.0277	0.0285
4-Methyloctane	I9	0.2801	0.3318	0.3384
2-Methyloctane	I9	0.2956	0.3501	0.3606
1c,2t,4c-Trimethylcyclohexane	I9	0.0468	0.0554	0.0561
3-Ethylheptane	I9	0.0954	0.1130	0.1143
3-Methyloctane	I9	0.4452	0.5273	0.5377
3,3-Diethylpentane	I9	0.0513	0.0608	0.0592
1c,2t,3-Trimethylcyclohexane	N9	0.0691	0.0806	0.0753
1,1,2-Trimethylcyclohexane	N9	0.0220	0.0256	0.0239
1,2-Dimethylbenzene (o-Xylene)	A8	0.6575	0.6447	0.5386

i-Butylcyclopentane	N9	0.2264	0.2639	0.2484
UnknownC8s	U8	0.0219	0.0231	0.0242
n-Nonane	P9	1.8680	2.2126	2.2659
1,1-Methylethylcyclohexane	N9	0.2958	0.3504	0.3599
i-Propylbenzene	A9	0.2740	0.3041	0.2589
i-Propylcyclohexane	N9	0.1028	0.1198	0.1098
2,2-Dimethyloctane	I10	0.0731	0.0961	0.0955
2,4-Dimethyloctane	I10	0.0870	0.1143	0.1136
2,6-Dimethyloctane	I10	0.0126	0.0166	0.0171
2,5-Dimethyloctane	I10	0.0437	0.0574	0.0571
n-Butylcyclopentane	N9	0.3160	0.4093	0.3766
3,3-Dimethyloctane	I10	0.0251	0.0330	0.0328
n-Propylbenzene	A9	0.2564	0.2846	0.2423
3,6-Dimethyloctane	I10	0.2444	0.3211	0.3192
3-Methyl-5-ethylheptane	I10	0.3816	0.4520	0.4578
1,3-Methylethylbenzene	A9	0.3299	0.3662	0.3092
1,4-Methylethylbenzene	A9	0.1623	0.1802	0.1521
1,3,5-Trimethylbenzene	A9	0.1807	0.2006	0.1705
2,3-Dimethyloctane	I10	0.0743	0.0976	0.0970
5-Methylnonane	I10	0.2790	0.3666	0.3678
1,2-Methylethylbenzene	A9	0.5196	0.5768	0.4844
2-Methylnonane	I10	0.0763	0.1003	0.1015
3-Ethylheptane	I10	0.0712	0.0936	0.0930
3-Methylnonane	I10	0.1687	0.2217	0.2222
1,2,4-Trimethylbenzene	A9	0.0263	0.0292	0.0245
t-Butylbenzene	A10	0.2549	0.3160	0.2683
i-Butylcyclohexane	N10	0.2142	0.2775	0.2514
1t-Methyl-2-n-propylcyclohexane	I10	0.0366	0.0433	0.0439
i-Butylbenzene	A10	0.0791	0.0980	0.0845
sec-Butylbenzene	A10	0.0391	0.0485	0.0414
UnknownC9s	U9	1.9360	2.2932	2.3484
n-Decane	P10	1.2565	1.6510	1.6623
1,2,3-Trimethylbenzene	A9	0.1958	0.2173	0.1788
1,3-Methyl-i-propylbenzene	A10	0.1070	0.1188	0.0998
1,4-Methyl-i-propylbenzene	A10	0.0519	0.0576	0.0484
Sec-Butylcyclohexane	N10	0.3311	0.4289	0.3880
1,2-Methyl-i-propylbenzene	A10	0.1535	0.1903	0.1597
3-Ethylheptane	I10	0.0433	0.0569	0.0576
1,3-Diethylbenzene	A10	0.1413	0.1751	0.1491
1,3-Methyl-n-propylbenzene	A10	0.0369	0.0457	0.0390
1,4-Diethylbenzene	A10	0.1502	0.1862	0.1589
1,4-Methyl-n-propylbenzene	A10	0.0332	0.0412	0.0353
n-Butylbenzene	A10	0.0954	0.1183	0.1010
1,3-Dimethyl-5-ethylbenzene	A10	0.0669	0.0829	0.0705
1,2-Diethylbenzene	A10	0.1154	0.1430	0.1196
1,2-Methyl-n-propylbenzene	A10	0.1024	0.1269	0.1068
1,4-Dimethyl-2-ethylbenzene	A10	0.1352	0.1676	0.1405
1,3-Dimethyl-4-ethylbenzene	A10	0.0391	0.0485	0.0407
1,2-Dimethyl-4-ethylbenzene	A10	0.2072	0.2568	0.2160
1,3-Dimethyl-2-ethylbenzene	A10	0.1532	0.1899	0.1569
1t,2c,4-Trimethylcyclopentane	A10	0.4226	0.4379	0.4306
1,2-Dimethyl-3-ethylbenzene	A10	0.1147	0.1422	0.1172
1,2-Ethyl-i-propylbenzene	A10	0.1140	0.1413	0.1186
1,4-Methyl-t-butylbenzene	A11	0.1901	0.2356	0.1977
UnknownC10s	U10	3.1256	4.1069	4.1350
n-Undecane	P11	0.9529	1.3755	1.3657
1,4-Ethyl-i-propylbenzene	A11	0.0818	0.1014	0.0851
1,2,4,5-Tetramethylbenzene	A11	0.2225	0.2758	0.2290
1,2-Methyl-n-butylbenzene	A11	0.1165	0.1444	0.1212
1,2,3,5-Tetramethylbenzene	A11	0.1700	0.2107	0.1741
1,2-Methyl-t-butylbenzene	A11	0.1605	0.1989	0.1669
5-Methylindan	A11	0.0442	0.0695	0.0683
4-Methylindan	A11	0.0247	0.0389	0.0382
1,2-Ethyl-n-propylbenzene	A11	0.1773	0.2198	0.1844
2-Methylindan	A11	0.1029	0.1619	0.1590
1,3-Methyl-n-butylbenzene	A11	0.0847	0.1050	0.0881
1,3-Di-i-propylbenzene	A11	0.1161	0.1439	0.1207

sec-Pentylbenzene	A11	0.1129	0.1399	0.1174
n-Pentylbenzene	A11	0.0695	0.0951	0.0815
1t-M-2-(4MP)cyclopentane	P12	0.1086	0.1708	0.1677
1,2-Di-n-propylbenzene	A11	0.1507	0.1868	0.1567
1,4-Di-i-propylbenzene	A11	0.1706	0.2115	0.1775
Tetrahydronaphthalene	A10	0.1292	0.1601	0.1343
t-Decahydronaphthalene	A10	0.1374	0.1703	0.1429
Naphthalene	A10	0.1268	0.1501	0.1259
1-t-Butyl-3,5-dimethylbenzene	A12	0.0837	0.1037	0.0870
1,4-Ethyl-t-butylbenzene	A11	0.1522	0.1887	0.1583
UnknownC11s	U11	2.2312	3.2208	3.1979
n-Dodecane	P12	0.8024	1.2623	1.2395
1,3-Di-n-propylbenzene	A12	0.0946	0.1173	0.0984
1,3,5-Triethylbenzene	A12	0.0774	0.0859	0.0730
1,2,4-Triethylbenzene	A12	0.3268	0.3627	0.3044
1,4-Methyl-n-pentylbenzene	A12	0.0786	0.0974	0.0817
n-Hexylbenzene	A12	0.0690	0.1034	0.0887
1,2,3,4,5-Pentamethylbenzene	A13	0.1724	0.2137	0.1793
2-Methylnaphthalene	A11	0.2140	0.2810	0.2358
1-Methylnaphthalene	A11	0.2199	0.2888	0.2083
UnknownC12s	U12	2.1417	3.3691	3.3083
n-Tridecane	P13	0.5436	0.9255	0.8981
UnknownC13s	U13	1.1993	2.0419	1.9814
n-Tetradecane	P14	0.2727	0.4996	0.4838
UnknownC14s	U14	0.7316	1.3404	1.2980
UnknownC15s	U15	0.5686	1.1154	1.0677
<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. :	03
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 10, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	0710
LEASE NO. :		SAMPLED BY :	GALE MCENDREE-EMPACT
NAME/DESCRIP :	SALES GAS @ 14:10 BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	90
SAMPLE PRES. :	119	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; PROBE; LENGTH OF H2S STAIN @ 5PPM (1-7PPM) @ 14:15		

COMPONENT	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
ALCOHOLS	0.0001	0.0003		
HELIUM	0.02	0.00	---	---
HYDROGEN	0.00	0.00	---	---
OXYGEN/ARGON	0.01	0.01	---	---
NITROGEN	0.91	1.06	---	---
CARBON DIOXIDE	2.47	4.50	---	---
METHANE	68.80020	45.69970	---	---
ETHANE	12.6380	15.7346	3.3753	3.3938
PROPANE	9.2753	16.9348	2.5523	2.5662
I-BUTANE	0.8964	2.1573	0.2934	0.2950
N-BUTANE	2.9011	6.9817	0.9132	0.9182
I-PENTANE	0.5810	1.7305	0.2082	0.2094
N-PENTANE	0.6662	1.9902	0.2413	0.2426
HEXANES PLUS	0.8317	3.2009	0.3372	0.3389
TOTALS	100.00000	100.00000	7.9209	7.9641

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.650	14.730
BENZENE	0.0333	0.1077	LOW NET DRY REAL :	1245.0 /scf	1251.8 /scf
TOLUENE	0.0227	0.0866	NET WET REAL :	1223.2 /scf	1230.0 /scf
ETHYLBENZENE	0.0025	0.0110	HIGH GROSS DRY REAL :	1368.2 /scf	1375.6 /scf
XYLENES	0.0064	0.0282	GROSS WET REAL :	1344.3 /scf	1351.7 /scf
TOTAL BTEX	0.0649	0.2335	NET DRY REAL :	19581.9 /lb	19688.9 /lb
			GROSS DRY REAL :	21520.9 /lb	21638.5 /lb

RELATIVE DENSITY (AIR=1): 0.8327  
COMPRESSIBILITY FACTOR : 0.99559

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



303-637-0150

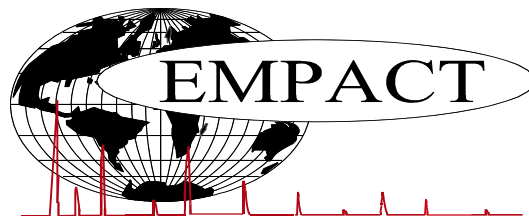
**EXTENDED NATURAL GAS ANALYSIS (\*DHA)**

**GLYCALC INFORMATION**

PROJECT NO. :	201403015	ANALYSIS NO. :	03
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 10, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	0710
LEASE NO. :		SAMPLED BY :	GALE MCENDREE-EMPACT
NAME/DESCRIP :	SALES GAS @ 14:10		
	BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	90
SAMPLE PRES. :	119	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; PROBE; LENGTH OF H2S STAIN @ 5PPM (1-7PPM) @ 14:15		

Componet	Mole %	Wt %
Helium	0.02	0.00
Hydrogen	0.00	0.00
Carbon Dioxide	2.47	4.50
Nitrogen	0.91	1.06
Methane	68.80020	45.69970
Ethane	12.6380	15.7346
Propane	9.2753	16.9348
Isobutane	0.8964	2.1573
n-Butane	2.9011	6.9817
Isopentane	0.5183	1.5484
n-Pentane	0.6662	1.9902
Cyclopentane	0.0627	0.1821
n-Hexane	0.1538	0.5488
Cyclohexane	0.0462	0.1610
Other Hexanes	0.2749	0.9729
Heptanes	0.1499	0.6175
Methycyclohexane	0.0395	0.1606
2,2,4 Trimethylpentane	0.0001	0.0005
Benzene	0.0333	0.1077
Toluene	0.0227	0.0866
Ethylbenzene	0.0025	0.0110
Xylenes	0.0064	0.0282
C8+ Heavies	0.1024	0.5061
<b>Subtotal</b>	<b>99.98990</b>	<b>99.98970</b>
Oxygen/Argon	0.01	0.01
Alcohols	0.0001	0.0003
<b>Total</b>	<b>100.00000</b>	<b>100.00000</b>

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

PROJECT NO. :	201403015	ANALYSIS NO. :	03
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	MARCH 10, 2014
ACCOUNT NO. :		SAMPLE DATE :	MARCH 4, 2014
PRODUCER :		CYLINDER NO. :	0710
LEASE NO. :		SAMPLED BY :	GALE MCENDREE-EMPACT
NAME/DESCRIP :	SALES GAS @ 14:10 BAILEY 8-26-8-60		
***FIELD DATA***		SAMPLE TEMP. :	90
SAMPLE PRES. :	119	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; PROBE; LENGTH OF H2S STAIN @ 5PPM (1-7PPM) @ 14:15		

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
Helium	---	0.02	0.00	---	---
Hydrogen	---	0.00	0.00	---	---
Oxygen/Argon	---	0.01	0.01	---	---
Nitrogen	---	0.91	1.06	---	---
Carbon Dioxide	---	2.47	4.50	---	---
Methane	P1	68.80020	45.69970	---	---
Ethane	P2	12.6380	15.7346	3.375	3.394
Propane	P3	9.2753	16.9348	2.552	2.566
i-Butane	I4	0.8964	2.1573	0.293	0.295
n-Butane	P4	2.9011	6.9817	0.913	0.918
2,2-Dimethylpropane	I5	0.0025	0.0075	0.001	0.001
i-Pentane	I5	0.5158	1.5409	0.188	0.189
n-Pentane	P5	0.6662	1.9902	0.241	0.243
t-Butanol	X4	0.0001	0.0003	0.000	0.000
2,2-Dimethylbutane	I6	0.0017	0.0061	0.001	0.001
Cyclopentane	N5	0.0627	0.1821	0.019	0.019
2,3-Dimethylbutane	I6	0.0102	0.0364	0.004	0.004
2-Methylpentane	I6	0.1106	0.3946	0.046	0.046
3-Methylpentane	I6	0.0566	0.2020	0.023	0.023
n-Hexane	P6	0.1538	0.5488	0.063	0.063
2,2-Dimethylpentane	I7	0.0001	0.0004	0.000	0.000
Methylcyclopentane	N6	0.0958	0.3338	0.034	0.034
2,4-Dimethylpentane	I7	0.0040	0.0166	0.002	0.002
2,2,3-Trimethylbutane	I7	0.0001	0.0004	0.000	0.000
Benzene	A6	0.0333	0.1077	0.009	0.009
3,3-Dimethylpentane	I7	0.0004	0.0017	0.000	0.000
Cyclohexane	N6	0.0462	0.1610	0.016	0.016
2-Methylhexane	I7	0.0175	0.0726	0.008	0.008
2,3-Dimethylpentane	I7	0.0080	0.0332	0.004	0.004
1,1-Dimethylcyclopentane	N7	0.0052	0.0212	0.002	0.002
3-Methylhexane	I7	0.0202	0.0838	0.009	0.009
1c,3-Dimethylcyclopentane	N7	0.0115	0.0468	0.005	0.005
1t,3-Dimethylcyclopentane	N7	0.0103	0.0419	0.005	0.005
3-Ethylpentane	I7	0.0015	0.0062	0.001	0.001
1t,2-Dimethylcyclopentane	N7	0.0204	0.0829	0.009	0.009
2,2,4-Trimethylpentane	I8	0.0001	0.0005	0.000	0.000
n-Heptane	P7	0.0438	0.1817	0.020	0.020
1c,2-Dimethylcyclopentane	N7	0.0017	0.0069	0.001	0.001
Methylcyclohexane	N7	0.0395	0.1606	0.016	0.016
2,2-Dimethylhexane	I8	0.0032	0.0152	0.001	0.001
Ethylcyclopentane	N7	0.0052	0.0212	0.002	0.002
2,5-Dimethylhexane	I8	0.0009	0.0043	0.000	0.000
2,2,3-Trimethylpentane	I8	0.0002	0.0010	0.000	0.000
2,4-Dimethylhexane	I8	0.0017	0.0080	0.001	0.001
1c,2t,4-Trimethylcyclopentane	N8	0.0033	0.0153	0.002	0.002
3,3-Dimethylhexane	I8	0.0002	0.0010	0.000	0.000
1t,2c,4-Trimethylcyclopentane	N8	0.0036	0.0167	0.002	0.002
2,3,4-Trimethylpentane	I8	0.0007	0.0033	0.000	0.000

Toluene	A7	0.0227	0.0866	0.008	0.008
2,3-Dimethylhexane	I8	0.0016	0.0076	0.001	0.001
2-Methyl-3-ethylpentane	I8	0.0009	0.0043	0.000	0.000
2-Methylheptane	I8	0.0082	0.0388	0.004	0.004
4-Methylheptane	I8	0.0023	0.0109	0.001	0.001
3-Methyl-3-ethylpentane	I8	0.0002	0.0010	0.000	0.000
3,4-Dimethylhexane	I8	0.0003	0.0014	0.000	0.000
1c,2c,4-Trimethylcyclopentane	N8	0.0002	0.0009	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0001	0.0005	0.000	0.000
3-Methylheptane	I8	0.0038	0.0180	0.002	0.002
1c,2t,3-Trimethylcyclopentane	N8	0.0057	0.0265	0.003	0.003
3-Ethylhexane	I8	0.0007	0.0033	0.000	0.000
1t,4-Dimethylcyclohexane	N8	0.0020	0.0093	0.001	0.001
1,1-Dimethylcyclohexane	N8	0.0007	0.0033	0.000	0.000
3t-Ethylmethylcyclopentane	N8	0.0011	0.0051	0.001	0.001
2t-Ethylmethylcyclopentane	N8	0.0010	0.0046	0.001	0.001
1,1-Methylethylcyclopentane	N8	0.0031	0.0144	0.002	0.002
2,2,4-Trimethylhexane	I9	0.0002	0.0011	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0031	0.0144	0.002	0.002
1t,3-Dimethylcyclohexane	N8	0.0001	0.0005	0.000	0.000
n-Octane	P8	0.0124	0.0586	0.006	0.006
1c,4-Dimethylcyclohexane	N8	0.0014	0.0065	0.001	0.001
i-Propylcyclopentane	I8	0.0003	0.0014	0.000	0.000
2,4,4-Trimethylhexane	I9	0.0001	0.0005	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0002	0.0011	0.000	0.000
2,2,3,4-Tetramethylpentane	I9	0.0001	0.0005	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0001	0.0005	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0010	0.0046	0.001	0.001
1,1,4-Trimethylcyclohexane	N9	0.0037	0.0193	0.002	0.002
2,2,3-Trimethylhexane	I9	0.0014	0.0075	0.001	0.001
2,4-Dimethylheptane	I9	0.0006	0.0032	0.000	0.000
4,4-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
Ethylcyclohexane	N8	0.0020	0.0093	0.001	0.001
n-Propylcyclopentane	N8	0.0008	0.0037	0.000	0.000
1c,3c,5-Trimethylcyclohexane	N9	0.0001	0.0005	0.000	0.000
2,5-Dimethylheptane	I9	0.0002	0.0011	0.000	0.000
3,3-Dimethylheptane	I9	0.0003	0.0016	0.000	0.000
3,5-Dimethylheptane	I9	0.0002	0.0011	0.000	0.000
2,6-Dimethylheptane	I9	0.0002	0.0011	0.000	0.000
1,1,3-Trimethylcyclohexane	N9	0.0004	0.0021	0.000	0.000
Ethylbenzene	I8	0.0025	0.0110	0.001	0.001
1c,2t,4t-Trimethylcyclohexane	N9	0.0008	0.0042	0.000	0.000
2,3-Dimethylheptane	I9	0.0001	0.0005	0.000	0.000
1,3-Dimethylbenzene (m-Xylene)	A8	0.0026	0.0114	0.001	0.001
1,4-Dimethylbenzene (p-Xylene)	A8	0.0019	0.0084	0.001	0.001
3,4-Dimethylheptane	I9	0.0014	0.0075	0.001	0.001
3,4-Dimethylheptane (2)	I9	0.0005	0.0027	0.000	0.000
4-Ethylheptane	I9	0.0002	0.0011	0.000	0.000
4-Methyloctane	I9	0.0008	0.0043	0.000	0.000
2-Methyloctane	I9	0.0009	0.0048	0.001	0.001
1c,2t,3-Trimethylcyclohexane	N9	0.0002	0.0010	0.000	0.000
3-Ethylheptane	I9	0.0004	0.0021	0.000	0.000
3-Methyloctane	I9	0.0011	0.0058	0.001	0.001
3,3-Diethylpentane	I9	0.0001	0.0005	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0019	0.0084	0.001	0.001
i-Butylcyclopentane	N9	0.0009	0.0047	0.000	0.000
n-Nonane	P9	0.0036	0.0191	0.002	0.002
1,1-Methylethylcyclohexane	N9	0.0006	0.0032	0.000	0.000
i-Propylbenzene	A9	0.0007	0.0035	0.000	0.000
i-Propylcyclohexane	N9	0.0002	0.0010	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
2,5-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
n-Butylcyclopentane	N9	0.0007	0.0036	0.000	0.000
3,3-Dimethyloctane	I10	0.0002	0.0012	0.000	0.000
n-Propylbenzene	A9	0.0007	0.0035	0.000	0.000
3,6-Dimethyloctane	I10	0.0005	0.0029	0.000	0.000
3-Methyl-5-ethylheptane	I10	0.0007	0.0041	0.000	0.000
1,3-Methylethylbenzene	A9	0.0005	0.0025	0.000	0.000
1,4-Methylethylbenzene	A9	0.0002	0.0010	0.000	0.000
1,3,5-Trimethylbenzene	A9	0.0003	0.0015	0.000	0.000
2,3-Dimethyloctane	I10	0.0001	0.0006	0.000	0.000
5-Methylnonane	I10	0.0003	0.0018	0.000	0.000

1,2-Methylethylbenzene	A9	0.0006	0.0030	0.000	0.000
2-Methylnonane	I10	0.0001	0.0006	0.000	0.000
3-Ethyl-octane	I10	0.0001	0.0006	0.000	0.000
3-Methylnonane	I10	0.0002	0.0012	0.000	0.000
t-Butylbenzene	A10	0.0006	0.0034	0.000	0.000
i-Butylcyclohexane	N10	0.0002	0.0012	0.000	0.000
i-Butylbenzene	A10	0.0001	0.0005	0.000	0.000
UnknownC9s	U9	0.0026	0.0138	0.001	0.001
n-Decane	P10	0.0009	0.0053	0.001	0.001
1,2,3-Trimethylbenzene	A9	0.0002	0.0010	0.000	0.000
1,3-Methyl-i-propylbenzene	A10	0.0001	0.0005	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0001	0.0005	0.000	0.000
Sec-Butylcyclohexane	A10	0.0002	0.0012	0.000	0.000
1,2-Methyl-i-propylbenzene	A10	0.0001	0.0005	0.000	0.000
1,3-Diethylbenzene	A10	0.0001	0.0005	0.000	0.000
1,4-Diethylbenzene	A10	0.0001	0.0005	0.000	0.000
n-Butylbenzene	A10	0.0001	0.0005	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0001	0.0005	0.000	0.000
1,2-Diethylbenzene	A10	0.0001	0.0005	0.000	0.000
t-Decahydronaphthalene	A9	0.0001	0.0006	0.000	0.000
1,2-Methyl-n-propylbenzene	A10	0.0001	0.0005	0.000	0.000
1,3-Dimethyl-4-ethylbenzene	A10	0.0001	0.0005	0.000	0.000
1,3-Dimethyl-2-ethylbenzene	A10	0.0001	0.0005	0.000	0.000
1,2-Dimethyl-3-ethylbenzene	A10	0.0001	0.0005	0.000	0.000
UnknownC10s	U10	0.0026	0.0153	0.002	0.002
n-Undecane	P11	0.0003	0.0020	0.000	0.000
1,2,4,5-Tetramethylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2,3,5-Tetramethylbenzene	A11	0.0001	0.0005	0.000	0.000
UnknownC11s	U11	0.0002	0.0013	0.000	0.000
n-Dodecane	P12	0.0001	0.0007	0.000	0.000
UnknownC12s	U12	0.0001	0.0007	0.000	0.000
n-Tridecane	P13	0.0001	0.0008	0.000	0.000
<b>TOTAL</b>		<b>100.00000</b>	<b>100.00000</b>	<b>7.9209</b>	<b>7.9641</b>

<b>BTEX COMPONENTS</b>	<b>MOLE%</b>	<b>WT%</b>	<b>BTU @</b>	<b>14.650</b>	<b>14.730</b>
BENZENE	0.0333	0.1077	<b>LOW</b> NET DRY REAL :	1245.0 /scf	1251.8 /scf
TOLUENE	0.0227	0.0866	NET WET REAL :	1223.2 /scf	1230.0 /scf
ETHYLBENZENE	0.0025	0.0110	<b>HIGH</b> GROSS DRY REAL :	1368.2 /scf	1375.6 /scf
XYLENES	0.0064	0.0282	GROSS WET REAL :	1344.3 /scf	1351.7 /scf
<b>TOTAL BTEX</b>	<b>0.0649</b>	<b>0.2335</b>	NET DRY REAL :	19581.9 /lb	19688.9 /lb
			GROSS DRY REAL :	21520.9 /lb	21638.5 /lb

RELATIVE DENSITY (AIR=1): 0.8327  
COMPRESSIBILITY FACTOR : 0.99559

(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.

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