



303-637-0150

EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

MAIN PAGE

PROJECT NO. :	201401066	ANALYSIS NO. :	01
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 17, 2014
ACCOUNT NO. :		SAMPLE DATE :	JANUARY 15, 2014
PRODUCER :		CYLINDER NO. :	2064
LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	TREATER @ 12:40		EMPACT
	KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	120
SAMPLE PRES. :	31	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	MOLE %	MASS %	VOL %
NITROGEN (AIR)	0.0549	0.0138	0.0127
CARBON DIOXIDE	0.0282	0.0111	0.0101
METHANE	0.2004	0.0288	0.0716
ETHANE	0.1878	0.0506	0.1059
PROPANE	0.5391	0.2131	0.3134
I-BUTANE	0.1723	0.0898	0.1189
N-BUTANE	0.8633	0.4497	0.5741
I-PENTANE	0.4683	0.3028	0.3614
N-PENTANE	0.8396	0.5429	0.6413
HEXANES PLUS	96.6461	98.2974	97.7906
TOTALS	100.0000	100.0000	100.0000

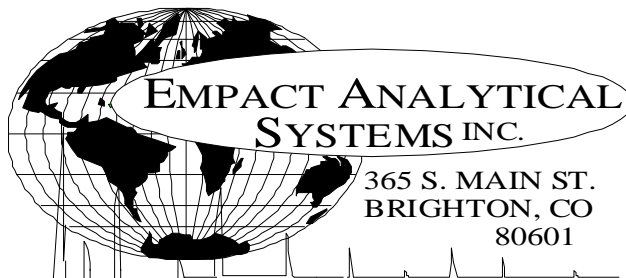
BTEX COMPONENTS	MOLE%	MASS%
BENZENE	1.6417	1.1493
TOLUENE	3.5125	2.9007
ETHYLBENZENE	0.6955	0.6618
XYLENE	2.2832	2.1726
TOTAL BTEX	8.1329	6.8844

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

	TOTAL SAMPLE	C6+ FRACTION
Specific Gravity (H2O=1) =	0.7457	0.7495 60/60
API Gravity =	58.25	57.29 60/60
Molecular Weight =	111.57	114.115
Absolute Density =	6.22	6.25 LBS/GAL
Heating Value Liq. Idl Gas=	126370	127287 BTU/GAL
Vapor/Liquid =	21.15	20.84 CUFT/GAL
Vapor Pressure =	15.05	1.72 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.
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 LIQUID ANALYSIS (*DHA)

E & P TANK / GLYCALC INFORMATION

PROJECT NO. :	201401066	ANALYSIS NO. :	01
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 17, 2014
ACCOUNT NO. :		SAMPLE DATE :	JANUARY 15, 2014
PRODUCER :		CYLINDER NO. :	2064
LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	TREATER @ 12:40		EMPACT
	KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	120
SAMPLE PRES. :	31	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	Mole %	Wt %	LV %
CARBON DIOXIDE	0.0282	0.0111	0.0101
NITROGEN (AIR)	0.0549	0.0138	0.0127
METHANE	0.2004	0.0288	0.0716
ETHANE	0.1878	0.0506	0.1059
PROPANE	0.5391	0.2131	0.3134
I-BUTANE	0.1723	0.0898	0.1189
N-BUTANE	0.8633	0.4497	0.5741
I-PENTANE	0.4683	0.3028	0.3614
N-PENTANE	0.8396	0.5429	0.6413
CYCLOPENTANE (N-C5)	1.3857	0.8710	0.8542
N-HEXANE	7.0308	5.4302	6.0995
CYCLOHEXANE (OTHER C6)	2.9961	2.2600	2.1506
OTHER HEXANES	10.4785	8.0081	8.5277
OTHER HEPTANES	14.8577	13.2514	13.8896
METHYLCYCLOHEXANE (OTHER C7)	4.2350	3.7270	3.5866
2,2,4 TRIMETHYLPENTANE	0.8218	0.7232	0.7156
BENZENE	1.6417	1.1493	0.9705
TOLUENE	3.5125	2.9007	2.4736
ETHYLBENZENE	0.6955	0.6618	0.5643
XYLENES	2.2832	2.1726	1.8556
OTHER OCTANES	12.1873	12.4759	12.5262
OCTANES PLUS	----	50.5081	----
NONANES	12.1205	13.7848	13.6304
DECANES PLUS	22.3998	30.8814	29.9462
SUB TOTAL	100.0000	100.0000	100.0000
TOTAL	100.0000	100.0000	100.0000

API Gravity	=	58.25	60/60
Vapor Pressure	=	15.05	PSIA & 100 F
Average Molecular Weight of Decanes plus	=	153.82	
Average Specific Gravity of Decanes plus	=	0.7700	

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 ITS APPLICATION.



303-637-0150

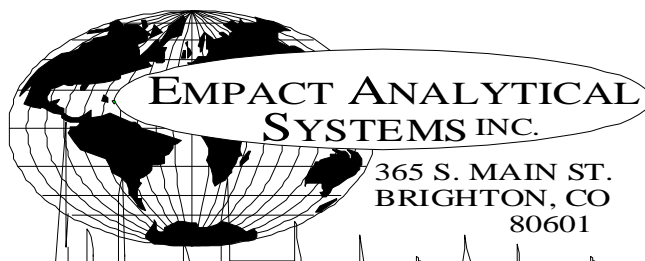
EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

BY CARBON NUMBER

PROJECT NO. :	201401066	ANALYSIS NO. :	01
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 17, 2014
ACCOUNT NO. :		SAMPLE DATE :	JANUARY 15, 2014
PRODUCER :		CYLINDER NO. :	2064
LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	TREATER @ 12:40		EMPACT
	KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	120
SAMPLE PRES. :	31	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT / CARBON NUMBER	MOLE%	MASS %	VOLUME %
NITROGEN	0.0549	0.0138	0.0127
CARBON DIOXIDE	0.0282	0.0111	0.0101
C1	0.2004	0.0288	0.0716
C2	0.1878	0.0506	0.1059
C3	0.5391	0.2131	0.3134
C4	1.0356	0.5395	0.6930
C5	2.6936	1.7167	1.8569
C6	22.1471	16.8476	17.7483
C7	22.6052	19.8791	19.9498
C8	15.9878	16.0335	15.6617
C9	12.1205	13.7848	13.6304
C10	11.2730	13.9502	13.5944
C11	4.9920	6.7500	6.4861
C12	2.6213	3.8174	3.6982
C13	1.1767	1.9169	1.8663
C14	0.8661	1.5400	1.5041
C15	0.8707	1.6577	1.6005
C16	0.3698	0.7505	0.7199
C17	0.2103	0.4533	0.4334
C18	0.0199	0.0454	0.0433
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

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 ITS APPLICATION.



303-637-0150

EXTENDED NATURAL GAS LIQUID ANALYSIS (*DHA)

DHA COMPONENT LIST

PROJECT NO. :	201401066	ANALYSIS NO. :	01
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 17, 2014
ACCOUNT NO. :		SAMPLE DATE :	JANUARY 15, 2014
PRODUCER :		CYLINDER NO.:	2064
LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	TREATER @ 12:40		EMPACT
	KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	120
SAMPLE PRES. :	31	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE		

COMPONENT	PIANO #	MOLE %	MASS %	VOL %
Nitrogen	NHC	0.0549	0.0138	0.0127
Carbon Dioxide	NHC	0.0282	0.0111	0.0101
Methane	P1	0.2004	0.0288	0.0716
Ethane	P2	0.1878	0.0506	0.1059
Propane	P3	0.5391	0.2131	0.3134
i-Butane	I4	0.1723	0.0898	0.1189
n-Butane	P4	0.8633	0.4497	0.5741
i-Pentane	I5	0.4683	0.3028	0.3614
n-Pentane	P5	0.8396	0.5429	0.6413
2,2-Dimethylbutane	I6	0.0348	0.0269	0.0307
Cyclopentane	N5	1.3857	0.8710	0.8542
2,3-Dimethylbutane	I6	0.3115	0.2406	0.2689
2-Methylpentane	I6	3.3689	2.6022	2.9494
3-Methylpentane	I6	2.0360	1.5726	1.7527
n-Hexane	P6	7.0308	5.4302	6.0995
2,2-Dimethylpentane	I7	0.0151	0.0136	0.0149
Methylcyclopentane	N6	4.7273	3.5658	3.5260
2,4-Dimethylpentane	I7	0.2815	0.2528	0.2787
2,2,3-Trimethylbutane	I7	0.0363	0.0326	0.0350
Benzene	A6	1.6417	1.1493	0.9705
3,3-Dimethylpentane	I7	0.0171	0.0154	0.0165
Cyclohexane	N6	2.9961	2.2600	2.1506
2-Methylhexane	I7	1.1123	0.9989	1.0914
2,3-Dimethylpentane	I7	0.9081	0.8155	0.8656
1,1-Dimethylcyclopentane	N7	0.4021	0.3539	0.3475
3-Methylhexane	I7	1.9079	1.7134	1.8438
1c,3-Dimethylcyclopentane	N7	0.8964	0.7889	0.7847
1t,3-Dimethylcyclopentane	N7	0.8218	0.7232	0.7156
3-Ethylpentane	I7	0.0405	0.0364	0.0385
1t,2-Dimethylcyclopentane	N7	1.8932	1.6661	1.6428
2,2,4-Trimethylpentane	I8	0.1892	0.1937	0.2065
n-Heptane	P7	5.4382	4.8839	5.2905
1c,2-Dimethylcyclopentane	N7	0.1419	0.1249	0.1198
Methylcyclohexane	N7	4.2350	3.7270	3.5866
2,2-Dimethylhexane	I8	0.4066	0.4163	0.4433
Ethylcyclopentane	N7	0.9453	0.8319	0.8043
2,5-Dimethylhexane	I8	0.1101	0.1127	0.1203
2,2,3-Trimethylpentane	I8	0.0460	0.0471	0.0488
2,4-Dimethylhexane	I8	0.2776	0.2842	0.3020
1c,2t,4-Trimethylcyclopentane	N8	0.4124	0.4148	0.4027
3,3-Dimethylhexane	I8	0.0645	0.0660	0.0689

2,3,4-Trimethylpentane	I8	0.0832	0.0852	0.0877
2,3,3-Trimethylpentane	I8	0.0414	0.0424	0.0433
Toluene	A7	3.5125	2.9007	2.4736
2,3-Dimethylhexane	I8	0.1338	0.1370	0.1426
2-Methyl-3-ethylpentane	I8	0.2352	0.2408	0.2480
1,1,2-Trimethylcyclopentane	N8	0.0517	0.0520	0.0499
2-Methylheptane	I8	1.3600	1.3924	1.4754
4-Methylheptane	I8	0.4640	0.4751	0.4913
3-Methyl-3-ethylpentane	I8	0.1669	0.1709	0.1742
3,4-Dimethylhexane	I8	0.1281	0.1312	0.1351
1c,2c,4-Trimethylcyclopentane	N8	0.0600	0.0603	0.0579
1c,3-Dimethylcyclohexane	N8	0.0459	0.0462	0.0447
3-Methylheptane	I8	0.0800	0.0819	0.0860
1c,2t,3-Trimethylcyclopentane	N8	1.1163	1.1227	1.0801
3-Ethylhexane	I8	0.3459	0.3541	0.3681
1t,4-Dimethylcyclohexane	N8	0.7434	0.7476	0.7266
1,1-Dimethylcyclohexane	N8	0.1400	0.1408	0.1337
3c-Ethylmethylcyclopentane	N8	0.0128	0.0129	0.0125
3t-Ethylmethylcyclopentane	N8	0.2842	0.2858	0.2763
2t-Ethylmethylcyclopentane	N8	0.2130	0.2142	0.2065
1,1-Methylethylcyclopentane	N8	0.8591	0.8640	0.8202
2,2,4-Trimethylhexane	I9	0.0703	0.0808	0.0837
1t,2-Dimethylcyclohexane	N8	0.6070	0.6105	0.5834
1t,3-Dimethylcyclohexane	N8	0.0022	0.0022	0.0021
n-Octane	P8	2.1943	2.2466	2.3692
1c,4-Dimethylcyclohexane	N8	0.8895	0.8946	0.8471
i-Propylcyclopentane	I8	0.1310	0.1318	0.1258
2,4,4-Trimethylhexane	I9	0.0233	0.0268	0.0275
2,2,3,4-Tetramethylpentane	I9	0.0843	0.0969	0.0998
2,3,4-Trimethylhexane	I9	0.0174	0.0200	0.0205
1c,2-Dimethylcyclohexane	N8	0.1733	0.1743	0.1623
2,3,5-Trimethylhexane	I9	0.0593	0.0682	0.0700
2,2-Dimethylheptane	I9	0.0220	0.0253	0.0264
1,1,4-Trimethylcyclohexane	N9	1.1061	1.2515	1.2020
2,2,3-Trimethylhexane	I9	0.3641	0.4186	0.4254
2,4-Dimethylheptane	I9	0.0743	0.0854	0.0885
4,4-Dimethylheptane	I9	0.1366	0.1570	0.1626
Ethylcyclohexane	N8	0.7117	0.7158	0.6737
n-Propylcyclopentane	N8	0.1796	0.1806	0.1724
1c,3c,5-Trimethylcyclohexane	N9	0.0588	0.0665	0.0639
2,5-Dimethylheptane	I9	0.1272	0.1462	0.1512
3,3-Dimethylheptane	I9	0.1065	0.1224	0.1266
3,5-Dimethylheptane	I9	0.0802	0.0922	0.0953
2,6-Dimethylheptane	I9	0.0437	0.0502	0.0525
1,1,3-Trimethylcyclohexane	N9	0.0842	0.0953	0.0915
Ethylbenzene	A8	0.6955	0.6618	0.5643
1c,2t,4t-Trimethylcyclohexane	N9	0.4765	0.5391	0.5079
2,3-Dimethylheptane	I9	0.0102	0.0117	0.0119
1,3-Dimethylbenzene (m-Xylene)	A8	0.7224	0.6874	0.5895
1,4-Dimethylbenzene (p-Xylene)	A8	0.8748	0.8324	0.7161
3,4-Dimethylheptane	I9	0.2313	0.2659	0.2695
3,4-Dimethylheptane (2)	I9	0.2505	0.2880	0.2919
4-Ethylheptane	I9	0.0973	0.1119	0.1159
4-Methyloctane	I9	0.2892	0.3325	0.3420
2-Methyloctane	I9	0.2621	0.3013	0.3130
1c,2t,4c-Trimethylcyclohexane	I9	0.0824	0.0947	0.0967
3-Ethylheptane	I9	0.0832	0.0956	0.0975
3-Methyloctane	I9	0.2919	0.3356	0.3451
3,3-Diethylpentane	I9	0.0947	0.1089	0.1070
1c,2t,3-Trimethylcyclohexane	N9	0.1421	0.1608	0.1515
1,1,2-Trimethylcyclohexane	N9	0.0672	0.0760	0.0716
1,2-Dimethylbenzene (o-Xylene)	A8	0.6860	0.6528	0.5500
i-Butylcyclopentane	N9	0.3050	0.3451	0.3276
UnknownC8s	U8	0.0492	0.0504	0.0532
n-Nonane	P9	1.0472	1.2038	1.2434
1,1-Methylethylcyclohexane	N9	0.8599	0.9885	1.0240
i-Propylbenzene	A9	0.2774	0.2988	0.2565
i-Propylcyclohexane	N9	0.0915	0.1035	0.0957
2,2-Dimethyloctane	I10	0.0430	0.0548	0.0549
2,4-Dimethyloctane	I10	0.0501	0.0639	0.0641
2,6-Dimethyloctane	I10	0.0175	0.0223	0.0231
2,5-Dimethyloctane	I10	0.0341	0.0435	0.0436
n-Butylcyclopentane	N9	0.1955	0.2458	0.2281

3,3-Dimethyloctane	I10	0.1365	0.1741	0.1747
n-Propylbenzene	A9	0.2439	0.2627	0.2256
3,6-Dimethyloctane	I10	0.1905	0.2429	0.2435
3-Methyl-5-ethylheptane	I10	0.3950	0.4541	0.4639
1,3-Methylethylbenzene	A9	0.2615	0.2817	0.2399
1,4-Methylethylbenzene	A9	0.2948	0.3176	0.2704
1,3,5-Trimethylbenzene	A9	0.0985	0.1061	0.0910
2,3-Dimethyloctane	I10	0.0568	0.0724	0.0726
5-Methylnonane	I10	0.3872	0.4938	0.4997
1,2-Methylethylbenzene	A9	0.2641	0.2845	0.2410
2-Methylnonane	I10	0.1172	0.1495	0.1526
3-Ethylheptane	I10	0.1067	0.1361	0.1364
3-Methylnonane	I10	0.1766	0.2252	0.2276
1,2,4-Trimethylbenzene	A9	0.0741	0.0798	0.0676
t-Butylbenzene	A10	0.1930	0.2322	0.1988
i-Butylcyclohexane	N10	0.1670	0.2100	0.1918
1t-Methyl-2-n-propylcyclohexane	I10	0.1057	0.1215	0.1241
i-Butylbenzene	A10	0.0656	0.0789	0.0686
sec-Butylbenzene	A10	0.1192	0.1434	0.1234
UnknownC9s	U9	3.0244	3.4768	3.5910
n-Decane	P10	1.1011	1.4042	1.4259
1,2,3-Trimethylbenzene	A9	0.2458	0.2648	0.2197
1,3-Methyl-i-propylbenzene	A10	0.0964	0.1039	0.0880
1,4-Methyl-i-propylbenzene	A10	0.1005	0.1083	0.0917
Sec-Butylcyclohexane	N10	0.2412	0.3032	0.2766
1,2-Methyl-i-propylbenzene	A10	0.1872	0.2252	0.1906
3-Ethylheptane	I10	0.0705	0.0899	0.0917
1,3-Diethylbenzene	A10	0.1408	0.1694	0.1455
1,3-Methyl-n-propylbenzene	A10	0.0565	0.0680	0.0586
1,4-Diethylbenzene	A10	0.1886	0.2269	0.1953
1,4-Methyl-n-propylbenzene	A10	0.0682	0.0820	0.0708
n-Butylbenzene	A10	0.0775	0.0932	0.0802
1,3-Dimethyl-5-ethylbenzene	A10	0.1155	0.1389	0.1191
1,2-Diethylbenzene	A10	0.0707	0.0850	0.0717
1,2-Methyl-n-propylbenzene	A10	0.0745	0.0896	0.0761
1,4-Dimethyl-2-ethylbenzene	A10	0.0801	0.0964	0.0815
1,3-Dimethyl-4-ethylbenzene	A10	0.1593	0.1916	0.1622
1,2-Dimethyl-4-ethylbenzene	A10	0.1405	0.1690	0.1434
1,3-Dimethyl-2-ethylbenzene	A10	0.0886	0.1066	0.0888
1t,2c,4-Trimethylcyclopentane	A10	0.5291	0.5321	0.5277
1,2-Dimethyl-3-ethylbenzene	A10	0.1594	0.1918	0.1595
1,2-Ethyl-i-propylbenzene	A10	0.1213	0.1459	0.1235
1,4-Methyl-t-butylbenzene	A11	0.0851	0.1024	0.0867
UnknownC10s	U10	4.8087	6.1322	6.2271
n-Undecane	P11	0.5866	0.8218	0.8230
1,4-Ethyl-i-propylbenzene	A11	0.3126	0.3761	0.3183
1,2,4,5-Tetramethylbenzene	A11	0.0874	0.1051	0.0880
1,2-Methyl-n-butylbenzene	A11	0.0662	0.0796	0.0674
1,2,3,5-Tetramethylbenzene	A11	0.0493	0.0593	0.0494
1,2-Methyl-t-butylbenzene	A11	0.0415	0.0499	0.0422
5-Methylindan	A11	0.0135	0.0206	0.0204
4-Methylindan	A11	0.0242	0.0369	0.0365
1,2-Ethyl-n-propylbenzene	A11	0.1616	0.1944	0.1645
2-Methylindan	A11	0.0895	0.1366	0.1353
1,3-Methyl-n-butylbenzene	A11	0.0714	0.0859	0.0727
1,3-Di-i-propylbenzene	A11	0.0334	0.0402	0.0340
sec-Pentylbenzene	A11	0.0520	0.0626	0.0530
n-Pentylbenzene	A11	0.0713	0.0947	0.0818
1t-M-2-(4MP)cyclopentane	P12	0.0776	0.1185	0.1174
1,2-Di-n-propylbenzene	A11	0.1114	0.1340	0.1134
1,4-Di-i-propylbenzene	A11	0.0762	0.0917	0.0776
Tetrahydronaphthalene	A10	0.0436	0.0525	0.0444
t-Decahydronaphthalene	A10	0.1069	0.1286	0.1088
Naphthalene	A10	0.0846	0.0972	0.0823
1-t-Butyl-3,5-dimethylbenzene	A12	0.0395	0.0475	0.0402
1,4-Ethyl-t-butylbenzene	A11	0.0618	0.0743	0.0629
UnknownC11s	U11	2.8799	4.0347	4.0404
n-Dodecane	P12	0.5395	0.8237	0.8158
1,3-Di-n-propylbenzene	A12	0.0594	0.0715	0.0605
1,3,5-Triethylbenzene	A12	0.0571	0.0615	0.0527
1,2,4-Triethylbenzene	A12	0.2266	0.2441	0.2067
1,4-Methyl-n-pentylbenzene	A12	0.0641	0.0771	0.0652
n-Hexylbenzene	A12	0.0604	0.0878	0.0759

1,2,3,4,5-Pentamethylbenzene	A13	0.0610	0.0734	0.0621
2-Methylnaphthalene	A11	0.0663	0.0845	0.0715
1-Methylnaphthalene	A11	0.0508	0.0647	0.0471
UnknownC12s	U12	1.4971	2.2857	2.2638
n-Tridecane	P13	0.1642	0.2713	0.2655
UnknownC13s	U13	0.9515	1.5722	1.5387
n-Tetradecane	P14	0.4051	0.7203	0.7035
UnknownC14s	U14	0.4610	0.8197	0.8006
n-Pentadecane	P15	0.2112	0.4021	0.3882
UnknownC15s	U15	0.6595	1.2556	1.2123
n-Hexadecane	P16	0.0921	0.1869	0.1793
UnknownC16s	U16	0.2777	0.5636	0.5406
n-Heptadecane	P17	0.0219	0.0472	0.0451
UnknownC17s	U17	0.1884	0.4061	0.3883
n-Octadecane	P18	0.0040	0.0091	0.0087
UnknownC18s	U18	0.0159	0.0363	0.0346
<u>TOTAL</u>		<u>100.0000</u>	<u>100.0000</u>	<u>100.0000</u>

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 ITS APPLICATION.



303-637-0150

EXTENDED NATURAL GAS ANALYSIS (*DHA)

MAIN PAGE

PROJECT NO. :	201401066	ANALYSIS NO. :	02
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 20, 2014
ACCOUNT NO. :		SAMPLE DATE :	JANUARY 15, 2014
PRODUCER :		CYLINDER NO. :	1173
LEASE NO. :		SAMPLED BY :	JOHN MOSER-EMPACT
NAME/DESCRIP :	SALES GAS @ 12:50 KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	105
SAMPLE PRES. :	31	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE; LENGTH OF H2S STAIN @ 6.5 PPM (1-7 PPM) @ 13:05		

COMPONENT	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
ALCOHOLS	0.0090	0.0192		
HELIUM	0.02	0.00	---	---
HYDROGEN	0.01	0.00	---	---
OXYGEN/ARGON	0.01	0.01	---	---
NITROGEN	1.21	1.23	---	---
CARBON DIOXIDE	2.60	4.15	---	---
METHANE	64.25210	37.42210	---	---
ETHANE	10.0749	10.9985	2.6939	2.7086
PROPANE	10.0799	16.1371	2.7760	2.7912
I-BUTANE	1.1803	2.4906	0.3858	0.3879
N-BUTANE	4.6663	9.8467	1.4702	1.4782
I-PENTANE	1.2934	3.3742	0.4600	0.4625
N-PENTANE	1.7007	4.4548	0.6163	0.6197
HEXANES PLUS	2.8934	9.8668	1.2013	1.2075
TOTALS	100.00000	100.00000	9.6035	9.6556

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.650	14.730
BENZENE	0.1074	0.3046	LOW NET DRY REAL :	1405.5 /scf	1413.1 /scf
TOLUENE	0.0843	0.2820	NET WET REAL :	1380.9 /scf	1388.6 /scf
ETHYLBENZENE	0.0092	0.0355	HIGH GROSS DRY REAL :	1540.6 /scf	1549.0 /scf
XYLENES	0.0224	0.0863	GROSS WET REAL :	1513.7 /scf	1522.0 /scf
TOTAL BTEX	0.2233	0.7084	NET DRY REAL :	19414.1 /lb	19520.1 /lb
			GROSS DRY REAL :	21267.0 /lb	21383.1 /lb

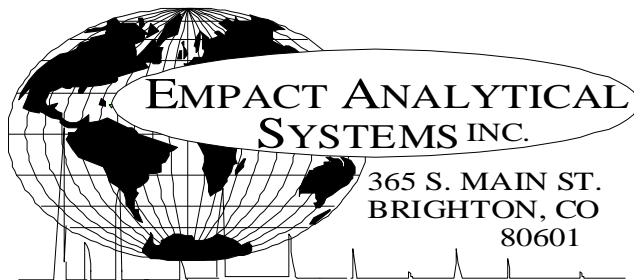
RELATIVE DENSITY (AIR=1): 0.9498
COMPRESSIBILITY FACTOR : 0.99470

(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

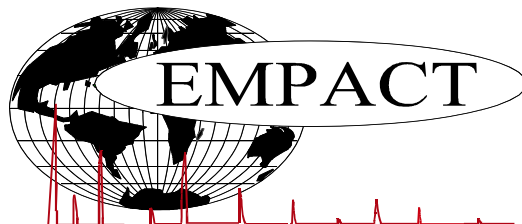
EXTENDED NATURAL GAS ANALYSIS (*DHA)

GLYCALC INFORMATION

PROJECT NO. :	201401066	ANALYSIS NO. :	02
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 20, 2014
ACCOUNT NO. :		SAMPLE DATE :	JANUARY 15, 2014
PRODUCER :		CYLINDER NO. :	1173
LEASE NO. :		SAMPLED BY :	JOHN MOSER-EMPACT
NAME/DESCRIP :	SALES GAS @ 12:50 KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	105
SAMPLE PRES. :	31	AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT; NO PROBE; LENGTH OF H2S STAIN @ 6.5 PPM (1-7 PPM) @ 13:05		

Componet	Mole %	Wt %
Helium	0.02	0.00
Hydrogen	0.01	0.00
Carbon Dioxide	2.60	4.15
Nitrogen	1.21	1.23
Methane	64.25210	37.42210
Ethane	10.0749	10.9985
Propane	10.0799	16.1371
Isobutane	1.1803	2.4906
n-Butane	4.6663	9.8467
Isopentane	1.1051	2.8947
n-Pentane	1.7007	4.4548
Cyclopentane	0.1883	0.4795
n-Hexane	0.5326	1.6663
Cyclohexane	0.1535	0.4690
Other Hexanes	0.8780	2.7232
Heptanes	0.5743	2.0753
Methycyclohexane	0.1259	0.4488
2,2,4 Trimethylpentane	0.0002	0.0008
Benzene	0.1074	0.3046
Toluene	0.0843	0.2820
Ethylbenzene	0.0092	0.0355
Xylenes	0.0224	0.0863
C8+ Heavies	0.4056	1.7750
Subtotal	99.98100	99.97080
Oxygen/Argon	0.01	0.01
Alcohols	0.0090	0.0192
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. : 201401066	ANALYSIS NO. : 02
COMPANY NAME : CARRIZO OIL & GAS	ANALYSIS DATE: JANUARY 20, 2014
ACCOUNT NO. :	SAMPLE DATE : JANUARY 15, 2014
PRODUCER :	CYLINDER NO. : 1173
LEASE NO. :	SAMPLED BY : JOHN MOSER-EMPACT
NAME/DESCRIP : SALES GAS @ 12:50	
KONIG 2-31-11-59	
FIELD DATA	SAMPLE TEMP. : 105
SAMPLE PRES. : 31	AMBIENT TEMP.:
VAPOR PRES. :	GRAVITY :
COMMENTS : SPOT; NO PROBE; LENGTH OF H2S STAIN @ 6.5 PPM (1-7 PPM) @ 13:05	

COMPONENT	PIANO #	MOLE %	MASS %	GPM @ 14.650	GPM @ 14.730
Helium	---	0.02	0.00	---	---
Hydrogen	---	0.01	0.00	---	---
Oxygen/Argon	---	0.01	0.01	---	---
Nitrogen	---	1.21	1.23	---	---
Carbon Dioxide	---	2.60	4.15	---	---
Methane	P1	64.25210	37.42210	---	---
Ethane	P2	10.0749	10.9985	2.694	2.709
Propane	P3	10.0799	16.1371	2.776	2.791
i-Butane	I4	1.1803	2.4906	0.386	0.388
n-Butane	P4	4.6663	9.8467	1.470	1.478
2,2-Dimethylpropane	I5	0.0035	0.0092	0.001	0.001
Ethanol	X2	0.0004	0.0007	0.000	0.000
i-Pentane	I5	1.1016	2.8855	0.403	0.405
Acetone	X3	0.0067	0.0141	0.002	0.002
i-Propanol	X3	0.0013	0.0028	0.000	0.000
n-Pentane	P5	1.7007	4.4548	0.616	0.620
t-Butanol	X4	0.0006	0.0016	0.000	0.000
2,2-Dimethylbutane	I6	0.0036	0.0113	0.001	0.001
Cyclopentane	N5	0.1883	0.4795	0.056	0.056
2,3-Dimethylbutane	I6	0.0283	0.0885	0.012	0.012
2-Methylpentane	I6	0.3395	1.0622	0.141	0.142
3-Methylpentane	I6	0.1814	0.5676	0.074	0.075
n-Hexane	P6	0.5326	1.6663	0.219	0.220
2,2-Dimethylpentane	I7	0.0012	0.0044	0.001	0.001
Methylcyclopentane	N6	0.3249	0.9927	0.115	0.116
2,4-Dimethylpentane	I7	0.0130	0.0473	0.006	0.006
2,2,3-Trimethylbutane	I7	0.0003	0.0011	0.000	0.000
Benzene	A6	0.1074	0.3046	0.030	0.030
3,3-Dimethylpentane	I7	0.0010	0.0036	0.000	0.000
Cyclohexane	N6	0.1535	0.4690	0.052	0.052
2-Methylhexane	I7	0.0675	0.2456	0.031	0.031
2,3-Dimethylpentane	I7	0.0321	0.1168	0.015	0.015
1,1-Dimethylcyclopentane	N7	0.0096	0.0342	0.004	0.004
3-Methylhexane	I7	0.0800	0.2910	0.037	0.037
1c,3-Dimethylcyclopentane	N7	0.0378	0.1348	0.017	0.017
1t,3-Dimethylcyclopentane	N7	0.0342	0.1219	0.016	0.016

3-Ethylpentane	I7	0.0058	0.0211	0.003	0.003
1t,2-Dimethylcyclopentane	N7	0.0761	0.2713	0.035	0.035
2,2,4-Trimethylpentane	I8	0.0002	0.0008	0.000	0.000
UnknownC6s	U6	0.0003	0.0009	0.000	0.000
n-Heptane	P7	0.1809	0.6581	0.083	0.084
1c,2-Dimethylcyclopentane	N7	0.0053	0.0189	0.002	0.002
Methylcyclohexane	N7	0.1259	0.4488	0.050	0.050
2,2-Dimethylhexane	I8	0.0062	0.0257	0.003	0.003
Ethylcyclopentane	N7	0.0293	0.1045	0.012	0.012
2,5-Dimethylhexane	I8	0.0039	0.0162	0.002	0.002
2,2,3-Trimethylpentane	I8	0.0006	0.0025	0.000	0.000
2,4-Dimethylhexane	I8	0.0068	0.0282	0.003	0.003
1c,2t,4-Trimethylcyclopentane	N8	0.0117	0.0477	0.005	0.005
3,3-Dimethylhexane	I8	0.0006	0.0025	0.000	0.000
1t,2c,4-Trimethylcyclopentane	N8	0.0143	0.0583	0.007	0.007
2,3,4-Trimethylpentane	I8	0.0030	0.0125	0.001	0.001
2,3,3-Trimethylpentane	I8	0.0001	0.0004	0.000	0.000
Toluene	A7	0.0843	0.2820	0.028	0.028
2,3-Dimethylhexane	I8	0.0055	0.0228	0.003	0.003
2-Methyl-3-ethylpentane	I8	0.0038	0.0158	0.002	0.002
2-Methylheptane	I8	0.0332	0.1377	0.017	0.017
4-Methylheptane	I8	0.0094	0.0390	0.005	0.005
3-Methyl-3-ethylpentane	I8	0.0009	0.0037	0.000	0.000
3,4-Dimethylhexane	I8	0.0011	0.0046	0.001	0.001
1c,2c,4-Trimethylcyclopentane	N8	0.0006	0.0024	0.000	0.000
1c,3-Dimethylcyclohexane	N8	0.0006	0.0024	0.000	0.000
3-Methylheptane	I8	0.0157	0.0651	0.008	0.008
1c,2t,3-Trimethylcyclopentane	N8	0.0180	0.0733	0.009	0.009
3-Ethylhexane	I8	0.0019	0.0079	0.001	0.001
1t,4-Dimethylcyclohexane	N8	0.0059	0.0240	0.003	0.003
1,1-Dimethylcyclohexane	N8	0.0023	0.0094	0.001	0.001
3t-Ethylmethylcyclopentane	N8	0.0058	0.0236	0.003	0.003
2t-Ethylmethylcyclopentane	N8	0.0051	0.0208	0.003	0.003
1,1-Methylethylcyclopentane	N8	0.0163	0.0664	0.008	0.008
2,2,4-Trimethylhexane	I9	0.0006	0.0028	0.000	0.000
1t,2-Dimethylcyclohexane	N8	0.0093	0.0379	0.005	0.005
1t,3-Dimethylcyclohexane	N8	0.0001	0.0004	0.000	0.000
UnknownC7s	U7	0.0002	0.0007	0.000	0.000
n-Octane	P8	0.0525	0.2177	0.027	0.027
1c,4-Dimethylcyclohexane	N8	0.0038	0.0155	0.002	0.002
i-Propylcyclopentane	I8	0.0015	0.0061	0.001	0.001
2,4,4-Trimethylhexane	I9	0.0003	0.0014	0.000	0.000
2,3,5-Trimethylhexane	I9	0.0006	0.0028	0.000	0.000
2,2,3,4-Tetramethylpentane	I9	0.0003	0.0014	0.000	0.000
2,3,4-Trimethylhexane	I9	0.0002	0.0009	0.000	0.000
1c,2-Dimethylcyclohexane	N8	0.0036	0.0147	0.002	0.002
2,2-Dimethylheptane	I9	0.0002	0.0009	0.000	0.000
1,1,4-Trimethylcyclohexane	N9	0.0142	0.0651	0.007	0.007
2,2,3-Trimethylhexane	I9	0.0057	0.0265	0.003	0.003
2,4-Dimethylheptane	I9	0.0015	0.0070	0.001	0.001
4,4-Dimethylheptane	I9	0.0004	0.0019	0.000	0.000
Ethylcyclohexane	N8	0.0079	0.0322	0.004	0.004
n-Propylcyclopentane	N8	0.0030	0.0122	0.001	0.001
1c,3c,5-Trimethylcyclohexane	N9	0.0004	0.0018	0.000	0.000
2,5-Dimethylheptane	I9	0.0011	0.0051	0.001	0.001
3,3-Dimethylheptane	I9	0.0012	0.0056	0.001	0.001
3,5-Dimethylheptane	I9	0.0009	0.0042	0.001	0.001
2,6-Dimethylheptane	I9	0.0010	0.0047	0.001	0.001
1,1,3-Trimethylcyclohexane	N9	0.0020	0.0092	0.001	0.001
Ethylbenzene	I8	0.0092	0.0355	0.004	0.004

1c,2t,4t-Trimethylcyclohexane	N9	0.0038	0.0174	0.002	0.002
2,3-Dimethylheptane	I9	0.0035	0.0163	0.002	0.002
1,3-Dimethylbenzene (m-Xylene)	A8	0.0128	0.0493	0.005	0.005
1,4-Dimethylbenzene (p-Xylene)	A8	0.0026	0.0100	0.001	0.001
3,4-Dimethylheptane	I9	0.0006	0.0028	0.000	0.000
3,4-Dimethylheptane (2)	I9	0.0017	0.0079	0.001	0.001
4-Ethylheptane	I9	0.0005	0.0023	0.000	0.000
4-Methyloctane	I9	0.0034	0.0158	0.002	0.002
2-Methyloctane	I9	0.0040	0.0186	0.002	0.002
1c,2t,3-Trimethylcyclohexane	N9	0.0010	0.0046	0.001	0.001
3-Ethylheptane	I9	0.0008	0.0037	0.000	0.000
3-Methyloctane	I9	0.0049	0.0228	0.003	0.003
1c,2t,4c-Trimethylcyclohexane	I9	0.0003	0.0014	0.000	0.000
1,1,2-Trimethylcyclohexane	N9	0.0002	0.0009	0.000	0.000
3,3-Diethylpentane	I9	0.0007	0.0033	0.000	0.000
1,2-Dimethylbenzene (o-Xylene)	A8	0.0070	0.0270	0.003	0.003
i-Butylcyclopentane	N9	0.0028	0.0128	0.001	0.001
UnknownC8s	U8	0.0002	0.0008	0.000	0.000
n-Nonane	P9	0.0142	0.0661	0.008	0.008
1,1-Methylethylcyclohexane	N9	0.0015	0.0069	0.001	0.001
i-Propylbenzene	A9	0.0028	0.0122	0.001	0.001
i-Propylcyclohexane	N9	0.0007	0.0032	0.000	0.000
2,2-Dimethyloctane	I10	0.0007	0.0036	0.000	0.000
2,4-Dimethyloctane	I10	0.0005	0.0026	0.000	0.000
2,5-Dimethyloctane	I10	0.0003	0.0016	0.000	0.000
n-Butylcyclopentane	N9	0.0027	0.0124	0.002	0.002
3,3-Dimethyloctane	I10	0.0007	0.0036	0.000	0.000
n-Propylbenzene	A9	0.0027	0.0118	0.001	0.001
3,6-Dimethyloctane	I10	0.0024	0.0124	0.001	0.001
3-Methyl-5-ethylheptane	I10	0.0027	0.0139	0.002	0.002
1,3-Methylethylbenzene	A9	0.0021	0.0092	0.001	0.001
1,4-Methylethylbenzene	A9	0.0006	0.0026	0.000	0.000
1,3,5-Trimethylbenzene	A9	0.0005	0.0022	0.000	0.000
2,3-Dimethyloctane	I10	0.0005	0.0026	0.000	0.000
5-Methylnonane	I10	0.0012	0.0062	0.001	0.001
1,2-Methylethylbenzene	A9	0.0016	0.0070	0.001	0.001
2-Methylnonane	I10	0.0001	0.0005	0.000	0.000
3-Ethyloctane	I10	0.0004	0.0021	0.000	0.000
3-Methylnonane	I10	0.0011	0.0057	0.001	0.001
1,2,4-Trimethylbenzene	A9	0.0003	0.0013	0.000	0.000
t-Butylbenzene	A10	0.0016	0.0078	0.001	0.001
i-Butylcyclohexane	N10	0.0009	0.0046	0.000	0.000
1t-Methyl-2-n-propylcyclohexane	I10	0.0003	0.0015	0.000	0.000
i-Butylbenzene	A10	0.0002	0.0010	0.000	0.000
sec-Butylbenzene	A10	0.0006	0.0029	0.000	0.000
UnknownC9s	U9	0.0125	0.0582	0.007	0.007
n-Decane	P10	0.0045	0.0232	0.003	0.003
1,2,3-Trimethylbenzene	A9	0.0009	0.0039	0.001	0.001
1,3-Methyl-i-propylbenzene	A10	0.0003	0.0015	0.000	0.000
1,4-Methyl-i-propylbenzene	A10	0.0003	0.0015	0.000	0.000
Sec-Butylcyclohexane	A10	0.0010	0.0051	0.001	0.001
1,2-Methyl-i-propylbenzene	A10	0.0005	0.0024	0.000	0.000
3-Ethylnonane	I10	0.0001	0.0006	0.000	0.000
1,3-Diethylbenzene	A10	0.0005	0.0024	0.000	0.000
1,3-Methyl-n-propylbenzene	A10	0.0001	0.0005	0.000	0.000
1,4-Diethylbenzene	A10	0.0006	0.0029	0.000	0.000
n-Butylbenzene	A10	0.0004	0.0020	0.000	0.000
1,3-Dimethyl-5-ethylbenzene	A10	0.0002	0.0010	0.000	0.000
1,2-Diethylbenzene	A10	0.0003	0.0015	0.000	0.000
t-Decahydronaphthalene	A9	0.0004	0.0023	0.000	0.000

1,2-Methyl-n-propylbenzene	A10	0.0003	0.0015	0.000	0.000
1,4-Dimethyl-2-ethylbenzene	A10	0.0003	0.0015	0.000	0.000
1,2-Dimethyl-4-ethylbenzene	A10	0.0004	0.0020	0.000	0.000
1,3-Dimethyl-2-ethylbenzene	A10	0.0002	0.0010	0.000	0.000
1,2-Dimethyl-3-ethylbenzene	A10	0.0002	0.0010	0.000	0.000
1,2-Ethyl-i-propylbenzene	A10	0.0001	0.0005	0.000	0.000
1,4-Methyl-t-butylbenzene	A11	0.0003	0.0016	0.000	0.000
UnknownC10s	U10	0.0112	0.0579	0.007	0.007
n-Undecane	P11	0.0017	0.0097	0.001	0.001
1,4-Ethyl-i-propylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2,4,5-Tetramethylbenzene	A11	0.0002	0.0010	0.000	0.000
1,2-Methyl-n-butylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2,3,5-Tetramethylbenzene	A11	0.0002	0.0010	0.000	0.000
1,2-Methyl-t-butylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2-Ethyl-n-propylbenzene	A11	0.0002	0.0011	0.000	0.000
2-Methylindan	A11	0.0001	0.0005	0.000	0.000
1,3-Methyl-n-butylbenzene	A11	0.0001	0.0005	0.000	0.000
1,3-Di-i-propylbenzene	A11	0.0001	0.0006	0.000	0.000
sec-Pentylbenzene	A11	0.0002	0.0011	0.000	0.000
n-Pentylbenzene	A11	0.0001	0.0005	0.000	0.000
1,2-Di-n-propylbenzene	A11	0.0001	0.0006	0.000	0.000
1,4-Di-i-propylbenzene	A11	0.0003	0.0018	0.000	0.000
Naphthalene	A10	0.0001	0.0005	0.000	0.000
1-t-Butyl-3,5-dimethylbenzene	A12	0.0001	0.0006	0.000	0.000
1,4-Ethyl-t-butylbenzene	A11	0.0001	0.0006	0.000	0.000
1,3-Di-n-propylbenzene	A12	0.0001	0.0006	0.000	0.000
UnknownC11s	U11	0.0043	0.0244	0.003	0.003
n-Dodecane	P12	0.0007	0.0043	0.001	0.001
1,2,4-Triethylbenzene	A12	0.0001	0.0006	0.000	0.000
2-Methylnaphthalene	A11	0.0001	0.0005	0.000	0.000
UnknownC12s	U12	0.0022	0.0125	0.001	0.001
n-Tridecane	P13	0.0003	0.0020	0.000	0.000
UnknownC13s	U13	0.0002	0.0013	0.000	0.000
n-Tetradecane	P14	0.0001	0.0007	0.000	0.000
UnknownC14s	U14	0.0001	0.0007	0.000	0.000
TOTAL		100.00000	100.00000	9.6055	9.6576

BTEX COMPONENTS	MOLE%	WT%	BTU @	14.650	14.730
BENZENE	0.1074	0.3046	LOW NET DRY REAL :	1405.5 /scf	1413.1 /scf
TOLUENE	0.0843	0.2820	NET WET REAL :	1380.9 /scf	1388.6 /scf
ETHYLBENZENE	0.0092	0.0355	HIGH GROSS DRY REAL :	1540.6 /scf	1549.0 /scf
XYLENES	0.0224	0.0863	GROSS WET REAL :	1513.7 /scf	1522.0 /scf
TOTAL BTEX	0.2233	0.7084	NET DRY REAL :	19414.1 /lb	19520.1 /lb
			GROSS DRY REAL :	21267.0 /lb	21383.1 /lb

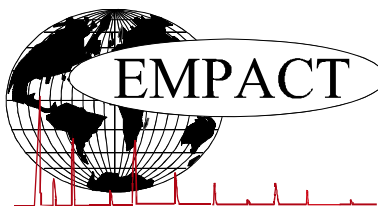
(CALC: GPA STD 2145 & TP-17 @14.696 & 60 F)

*(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.

RELATIVE DENSITY (AIR=1): 0.9498
COMPRESSIBILITY FACTOR : 0.99470



CRUDE OIL ASSAY

PROJECT NO. :	201401066	ANALYSIS NO. :	03
COMPANY NAME :	CARRIZO OIL & GAS	ANALYSIS DATE:	JANUARY 15, 2014
ACCOUNT NO. :		SAMPLE DATE :	
PRODUCER :		CYLINDER NO. :	JANUARY 17, 2014
LEASE NO. :		SAMPLED BY :	JOHN MOSER
NAME/DESCRIP :	PRODUCTION TANK @ 13:15		EMPACT
	KONIG 2-31-11-59		
FIELD DATA		SAMPLE TEMP. :	74
SAMPLE PRES. :		AMBIENT TEMP.:	
VAPOR PRES. :		GRAVITY :	
COMMENTS :	SPOT		

<u>SPECIFICATION</u>	<u>TEST METHOD</u>	<u>UNITS</u>	<u>RESULTS</u>
API GRAVITY		API 60/60	30.3
RVP @100 DEG F	D323	PSIG	4.6
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			DARK BROWN/BLACK
<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

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.