

2030 Afton Place Farmington, NM 87401 (505) 325-6622

Analysis No: HM2021091 Cust No: 33700-10435

Well/Lease Information

Customer Name: HARVEST MIDSTREAM

Well Name: Laguna Seca CDP

County/State: Rio Arriba NM

Location: Lease/PA/CA: Formation:

Cust. Stn. No.: 02013-01

Heat Trace: N

Remarks: Calculated Molecular Weight: = 17.1334

Source: METER RUN

Well Flowing: Y

Pressure: 95 PSIG
Flow Temp: 65 DEG. F
Ambient Temp: 60 DEG. F
Flow Rate: 3.5 MCF/D

Sample Method:

Sample Date: 10/18/2021 Sample Time: 11.15 AM Sampled By: D. Valencia

Sampled by (CO): Harvest Mid

Analysis

CO2 3.6895 3.7147 0.6310 0.00 0.0561 Methane 95.9097 96.5652 16.2900 968.69 0.5312 Ethane 0.1228 0.1236 0.0330 2.17 0.0013 Propane 0.0000 0.0000 0.0000 0.000 0.000 Iso-Butane 0.0000 0.0000 0.0000 0.000 0.000 N-Butane 0.0000 0.0000 0.0000 0.000 0.000 N-Butane 0.0000 0.0000 0.0000 0.000 0.0000 N-Butane 0.0000 0.0000 0.0000 0.000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.0000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.0000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.0000 0.0000 N-Pentane 0.0000 N/R </th <th>Component:</th> <th>Mole%:</th> <th>Unormalized %:</th> <th>**GPM:</th> <th>*BTU:</th> <th>*SP Gravity:</th>	Component:	Mole%:	Unormalized %:	**GPM:	*BTU:	*SP Gravity:
Methane 95.9097 96.5652 16.2900 968.69 0.5312 Ethane 0.1228 0.1236 0.0330 2.17 0.0013 Propane 0.0000 0.0000 0.0000 0.000 0.000 Iso-Butane 0.0000 0.0000 0.0000 0.000 0.000 N-Butane 0.0000 0.0000 0.0000 0.000 0.000 N-Pottane 0.0000 0.0000 0.0000 0.000 0.000 N-Pentane 0.0000 0.0000 0.0000 0.000 0.0000 N-Pentane 0.0000 0.000	Nitrogen	0.2513	0.2530	0.0280	0.00	0.0024
Ethane 0.1228 0.1236 0.0330 2.17 0.0013 Propane 0.0000 0.0000 0.0000 0.000 0.000 0.0000 Iso-Butane 0.0000 0.0000 0.0000 0.000 0.000 0.0000 N-Butane 0.0000 0.0000 0.0000 0.000 0.000 0.0000 Neopentane 2,2 dmc3 0.0268 0.0270 0.0100 1.07 0.0007 I-Pentane 0.0000 0.0000 0.0000 0.000 0.00 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.000 0.000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.000 0.000 0.0000 N-Pentane 0.0000 N/R 0.0000 0.00 0.000 0.0000 Cyclopentane 0.0000 N/R 0.0000 0.00 0.000 Cyclopentane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 C6 0.0000 N/R 0.0000 0.000 0.000 0.0000 Methylcyclopentane 0.0000 N/R 0.0000 0.000 0.0000 Department 0.0000 N/R 0.0000 0.000 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.000 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.000 0.0000 Department 0.0000 N/R 0.0000 0.000 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.0000 0.0000	CO2	3.6895	3.7147	0.6310	0.00	0.0561
Propane 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000	Methane	95.9097	96.5652	16.2900	968.69	0.5312
So-Butane 0.0000 0.0000 0.0000 0.000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.00000 0.00000 0.00000 0.00000 0.00000 0.00000 0.00000 0.000000 0.00000 0.00000 0.00000 0.0	Ethane	0.1228	0.1236	0.0330	2.17	0.0013
N-Butane 0.0000 0.0000 0.0000 0.000 0.0000 0.0000 0.0000 1.077 0.0007 1-Pentane 2,2 dmc3 0.0268 0.0270 0.0100 1.077 0.0007 1-Pentane 0.0000 0.0000 0.0000 0.0000 0.000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.000	Propane	0.0000	0.0000	0.0000	0.00	0.0000
Neopentane 2,2 dmc3 0.0268 0.0270 0.0100 1.07 0.0007 I-Pentane 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 N/R 0.0000 0.000 0.000 0.000 0.0000 0.0000 2-3-Dimethylbutane 0.0000 N/R 0.0000 0.00 0.0000 Cyclopentane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 Methylcyclopentane 0.0000 N/R 0.0000 0.00 0.0000 Benzene 0.0000 N/R 0.0000 0.00 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.00 0.000	Iso-Butane	0.0000	0.0000	0.0000	0.00	0.0000
I-Pentane 0.0000 0.0000 0.0000 0.000 0.000 0.0000 0.0000 N-Pentane 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0	N-Butane	0.0000	0.0000	0.0000	0.00	0.0000
N-Pentane 0.0000 0.0000 0.0000 0.000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000	Neopentane 2,2 dmc3	0.0268	0.0270	0.0100	1.07	0.0007
Neohexane 0.0000 N/R 0.0000 0.00 0.0000 2-3-Dimethylbutane 0.0000 N/R 0.0000 0.00 0.0000 Cyclopentane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 C6 0.0000 N/R 0.0000 0.00 0.0000 Methylcyclopentane 0.0000 N/R 0.0000 0.00 0.0000 Benzene 0.0000 N/R 0.0000 0.00 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 1-heptanes 0.0000 N/R 0.0000 0.00 0.0000	I-Pentane	0.0000	0.0000	0.0000	0.00	0.0000
2-3-Dimethylbutane 0.0000 N/R 0.0000 0.00 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.00	N-Pentane	0.0000	0.0000	0.0000	0.00	0.0000
Cyclopentane 0.0000 N/R 0.0000 0.000 0.0000 2-Methylpentane 0.0000 N/R 0.0000 0.000 0.0000 3-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 C6 0.0000 0.0000 0.0000 0.000 0.000 0.0000 Methylcyclopentane 0.0000 N/R 0.0000 0.00 0.0000 Benzene 0.0000 N/R 0.0000 0.00 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.00 0.0000 Heptanes 0.0000 N/R 0.0000 0.00 0.0000	Neohexane	0.0000	N/R	0.0000	0.00	0.0000
2-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.000	2-3-Dimethylbutane	0.0000	N/R	0.0000	0.00	0.0000
3-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.0000 0.000	Cyclopentane	0.0000	N/R	0.0000	0.00	0.0000
3-Methylpentane 0.0000 N/R 0.0000 0.000 0.0000 C6 0.0000 0.0000 0.0000 0.000 0.000 0.000 Methylcyclopentane 0.0000 N/R 0.0000 0.00 0.000 Benzene 0.0000 N/R 0.0000 0.00 0.000 Cyclohexane 0.0000 N/R 0.0000 0.00 0.000 2-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylpentane 0.0000 N/R 0.0000 0.00 0.0000 i-heptanes 0.0000 N/R 0.0000 0.000 0.0000	2-Methylpentane	0.0000	N/R	0.0000	0.00	0.0000
C6 0.0000 0.0000 0.0000 0.0000 0.0000 Methylcyclopentane 0.0000 N/R 0.0000 0.00 0.0000 Benzene 0.0000 N/R 0.0000 0.00 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.000 0.0000 M/R 0.0000 0.0000 0.0000 0.0000 0.0000	3-Methylpentane	0.0000	N/R			0.0000
Methylcyclopentane 0.0000 N/R 0.0000 0.000 0.0000 Benzene 0.0000 N/R 0.0000 0.00 0.0000 Cyclohexane 0.0000 N/R 0.0000 0.00 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.00 0.0000 I-heptanes 0.0000 N/R 0.0000 0.00 0.0000	C6	0.0000	0.0000			0.0000
Cyclohexane 0.0000 N/R 0.0000 0.000 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.000 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.000 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.000 0.0000 i-heptanes 0.0000 N/R 0.0000 0.000 0.0000	Methylcyclopentane	0.0000	N/R			0.0000
Cyclohexane 0.0000 N/R 0.0000 0.000 0.0000 2-Methylhexane 0.0000 N/R 0.0000 0.000 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.00 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.000 0.0000 i-heptanes 0.0000 N/R 0.0000 0.000 0.0000	Benzene	0.0000	N/R	0.0000	0.00	0.0000
2-Methylhexane 0.0000 N/R 0.0000 0.000 0.0000 3-Methylhexane 0.0000 N/R 0.0000 0.000 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.000 0.0000 i-heptanes 0.0000 N/R 0.0000 0.0000 0.0000	Cyclohexane	0.0000	N/R			0.0000
3-Methylhexane 0.0000 N/R 0.0000 0.000 0.0000 2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.000 0.0000 i-heptanes 0.0000 N/R 0.0000 0.0000	2-Methylhexane	0.0000	N/R			0.0000
2-2-4-Trimethylpentane 0.0000 N/R 0.0000 0.000 0.0000 i-heptanes 0.0000 N/R 0.0000 0.000 0.0000	3-Methylhexane	0.0000	N/R			0.0000
i-heptanes 0.0000 N/R 0.0000 0.0000	2-2-4-Trimethylpentane	0.0000	N/R			0.0000
0.0000 N/D	i-heptanes	0.0000	N/R			0.0000
	Heptane	0.0000	N/R	0.0000	0.00	0.0000

C10 i-C11	0.0000 0.0000	N/R N/R	0.0000	0.00	0.0000 0.0000
i-C10	0.0000	N/R	0.0000	0.00	0.0000
C9	0.0000	N/R	0.0000	0.00	0.0000
i-C9	0.0000	N/R	0.0000	0.00	0.0000
o Xylene (& 2,2,4 tmc7)	0.0000	N/R	0.0000	0.00 0.00	0.0000 0.0000
m, p Xylene	0.0000	N/R	0.0000	0.00	0.0000
Octane Ethylbenzene	0.0000 0.0000	N/R N/R	0.0000	0.00	0.0000
i-Octanes	0.0000	N/R	0.0000	0.00	0.0000
4-Methylheptane	0.0000	N/R	0.0000	0.00	0.0000
2-Methylheptane	0.0000	N/R	0.0000	0.00	0.0000
Toluene	0.0000	N/R	0.0000	0.00	0.0000
Received by OCD: 9/23/2022 1: Methylcyclohexane	0.0000	N/R	0.0000	0.00	Page 2 of 10 0.0000

^{* @ 14.730} PSIA DRY & UNCORRECTED FOR COMPRESSIBILITY

^{**@ 14.730} PSIA & 60 DEG. F.

COMPRESSIBLITY FACTOR	(1/Z):	1.0021	CYLINDER #:	0163
BTU/CU.FT IDEAL:		974.2	CYLINDER PRESSURE:	90 PSIG
BTU/CU.FT (DRY) CORRECTED FO	PR (1/Z):	976.2	ANALYSIS DATE:	10/20/2021
BTU/CU.FT (WET) CORRECTED FO	OR (1/Z):	959.2	ANALYIS TIME:	12:42:01 AM
DRY BTU @ 15.025:		995.8	ANALYSIS RUN BY:	ELAINE MORRISON
REAL SPECIFIC GRAVITY:		0.5927		

GPM, BTU, and SPG calculations as shown above are based on current GPA constants.

GPA Standard: GPA 2286-14

GC: SRI Instruments 8610 Last Cal/Verify: 10/26/2021

GC Method: C12+BTEX Gas



HARVEST MIDSTREAM WELL ANALYSIS COMPARISON

 Lease:
 Laguna Seca CDP
 METER RUN
 10/26/2021

 Stn. No.:
 02013-01
 33700-10435

Mtr. No.:

Smpl Date: 10/18/2021 04/16/2021 02/06/202 Test Date: 10/20/2021 04/21/2021 02/13/202 Run No: HM2021091 HM2021025 HM20001	20
TWIZOZ 1031 TWIZOZ 1020	
Nitrogen: 0.2513 0.0701 0.0654	
CO2: 3.6895 4.1404 5.1009	
Methane: 95.9097 95.5707 93.9967	
Ethane: 0.1228 0.2188 0.7593	
Propane: 0.0000 0.0000 0.0641	
I-Butane: 0.0000 0.0000 0.0075	
N-Butane: 0.0000 0.0000 0.0024	
2.2 dmc3: 0.0268 0.0000 0.0000	
I-Pentane: 0.0000 0.0000 0.0000	
N-Pentane: 0.0000 0.0000 0.0000	
Neohexane: 0.0000 0.0000 0.0001	
2-3- 0.0000 0.0000 0.0000	
Cyclopentane: 0.0000 0.0000 0.0000	
2-Methylpentane: 0.0000 0.0000 0.0003	
3-Methylpentane: 0.0000 0.0000 0.0001	
C6: 0.0000 0.0000 0.0003	
Methylcyclopentane: 0.0000 0.0000 0.0003	
Benzene: 0.0000 0.0000 0.0004	
Cyclohexane: 0.0000 0.0000 0.0001	
2-Methylhexane: 0.0000 0.0000 0.0001	
3-Methylhexane: 0.0000 0.0000 0.0000	
0.0000 0.0000 0.0000	
U.0000 U.0000 U.0000	
Methylovolohevane: 0.0000 0.0000 0.0002	
0.0000 0.0000 0.0005 Toluene:	
0.0000 0.0000 0.0005	
2-Methylheptane: 0.0000 0.0000 0.0000	
4-Methylheptane: 0.0000 0.0000 0.0001	
i-Octanes: 0.0000 0.0000 0.0001	
Octane: 0.0000 0.0000 0.0002	
Ethylbenzene: 0.0000 0.0000 0.0000	
m, p Xylene: 0.0000 0.0000 0.0002	
o Xylene (& 2,2,4 0.0000 0.0000 0.0000	
i-C9: 0.0000 0.0000 0.0001	
C9: 0.0000 0.0000 0.0001	
i-C10: 0.0000 0.0000 0.0000	
C10: 0.0000 0.0000 0.0000	
i-C11: 0.0000 0.0000 0.0000	
C11: 0.0000 0.0000 0.0000	
C12P: 0.0000 0.0000 0.0000	
BTU: 976.2 973.4 969.3	
GPM: 16.9820 17.0070 17.0710	
SPG: 0.5927 0.5962 0.6091	

2030 Afton Place, Farmington, NM 87401 - (5	705) 325-6622 GN#
C6+ □ C9+ □ C12+	BTEX 🗆 Helium 🗆
NALYSIS N2 Flowback Sulfu	rs 🗌 Ext. Liquid 🗌
SERVICE Other Extended gas Analy	SDate 10/18/2021
Sampled By:(Co.) Howest Midstreum	_Time 11/5
Sampled by:(Person) D. Vallucs	_Well Flowing: Wes No
Company: Homest Midstream	
Well Name: Asy lanner seca	Flow Pressure (PSIG): 98
Lease#: Werrers	_ Flow Temp (°F):
	_ Ambient Temp (°F):
State: NM Location: A59 Layre Sca	_Flow Rate (MCF/D):_3.5
Source: Meter Run Tubing Casing Bradenhead Other_	
Sample Type: Spot Composite Sample Method: Purge & Fill	Other
Meter Number: 02 29-01	Cylinder Number: 0163
Contact: D. Valuna 505-419-8357	33700-10435
Remarks: Extrade gns Analysis	Hm 2021091

LINE LEAK OR CONTINUOUS PSV RELEASE CALCULATOR AND REPORTING

Rosa 19D

ASSUMES NO PRESSURE LOSS AS RESULT OF LEAK

	DATE	PSI	PORT DIAMETER IN INCHES	TIME IN MINUTES BLOWN	MCF LOST	COMMENTS	
ſ		40.0	0.03125	17100.00	14.39		
		40.0	0.0625	17100.00	58.03		
		40.0	0.149	17100.00	329.83		
		40.0	0.344	17100.00	1,758.06		

2,160.31

District I
1625 N. French Dr., Hobbs, NM 88240
Phone: (575) 393-6161 Fax: (575) 393-0720

District II 811 S. First St., Artesia, NM 88210 Phone:(575) 748-1283 Fax:(575) 748-9720

District III 1000 Rio Brazos Rd., Aztec, NM 87410 Phone:(505) 334-6178 Fax:(505) 334-6170

1220 S. St Francis Dr., Santa Fe, NM 87505 Phone:(505) 476-3470 Fax:(505) 476-3462

State of New Mexico Energy, Minerals and Natural Resources Oil Conservation Division 1220 S. St Francis Dr. **Santa Fe, NM 87505**

DEFINITIONS

Action 145865

DEFINITIONS

Operator:	OGRID:
Harvest Four Corners, LLC	373888
1111 Travis Street	Action Number:
Houston, TX 77002	145865
	Action Type:
	[C-129] Amend Venting and/or Flaring (C-129A)

DEFINITIONS

For the sake of brevity and completeness, please allow for the following in all groups of questions and for the rest of this application:

- this application's operator, hereinafter "this operator";
- · venting and/or flaring, hereinafter "vent or flare";
- any notification or report(s) of the C-129 form family, hereinafter "any C-129 forms";
- the statements in (and/or attached to) this, hereinafter "the statements in this";
- and the past tense will be used in lieu of mixed past/present tense questions and statements.

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State of New Mexico Energy, Minerals and Natural Resources Oil Conservation Division 1220 S. St Francis Dr. **Santa Fe, NM 87505**

QUESTIONS

Action 145865

QUESTIONS

Operator.		JONE.	
Harvest Four Corners, LLC		373888	
1111 Travis Street		Action Number:	
Houston, TX 77002		145865	
		ction Type: [C-129] Amend Venting and/or Flaring (C-129A)	
QUESTIONS	·		
Prerequisites			
Any messages presented in this section, will prevent submission of this application. Please resolve	these issues before continui	ng with the rest of the questions.	
Incident Operator	[373888] Harvest Fou	r Corners, LLC	
Incident Type	Flare		
Incident Status	Closure Not Approved	1	
Incident Well	Not answered.		
Incident Facility	[fAPP2123052765] HAI	RVEST FOUR CORNERS GATHER SYSTEM	
Only valid Vent, Flare or Vent with Flaring incidents (selected above in the Application Details section	on) that are assigned to your	current operator can be amended with this C-129A application.	
Determination of Reporting Requirements			
	nd may provide addianal avia	danaa	
Answer all questions that apply. The Reason(s) statements are calculated based on your answers and may provide addionate the control of the c		nance.	
Was this vent or flare caused by an emergency or malfunction	No		
Did this vent or flare last eight hours or more cumulatively within any 24-hour period from a single event			
Is this considered a submission for a vent or flare event Yes, major		d/or flaring of natural gas.	
An operator shall file a form C-141 instead of a form C-129 for a release that, includes liquid during v	enting and/or flaring that is o	r may be a major or minor release under 19.15.29.7 NMAC.	
Was there at least 50 MCF of natural gas vented and/or flared during this event	Yes		
Did this vent or flare result in the release of ANY liquids (not fully and/or completely			
flared) that reached (or has a chance of reaching) the ground, a surface, a			
watercourse, or otherwise, with reasonable probability, endanger public health, the	No		
environment or fresh water			
Was the vent or flare within an incorporated municipal boundary or withing 300 feet			
from an occupied permanent residence, school, hospital, institution or church in	No		
existence			
	•		
Equipment Involved			
Primary Equipment Involved	Pipeline (Any)	v)	
	1 ()		
	I		

Representative Compositional Analysis of Vented or Flared Natural Gas		
Please provide the mole percent for the percentage questions in this group.		
Methane (CH4) percentage	96	
Nitrogen (N2) percentage, if greater than one percent	0	
Hydrogen Sulfide (H2S) PPM, rounded up	0	
Carbon Dioxide (C02) percentage, if greater than one percent	4	
Oxygen (02) percentage, if greater than one percent 0		
If you are venting and/or flaring because of Pipeline Specification, please provide the required specifications for each gas.		
Methane (CH4) percentage quality requirement	Not answered.	
Nitrogen (N2) percentage quality requirement	Not answered.	
Hydrogen Sufide (H2S) PPM quality requirement	Not answered.	
Carbon Dioxide (C02) percentage quality requirement	Not answered.	
Oxygen (02) percentage quality requirement	Not answered.	

Not answered.

Additional details for Equipment Involved. Please specify

QUESTIONS, Page 2

Action 145865

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District III

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State of New Mexico Energy, Minerals and Natural Resources Oil Conservation Division 1220 S. St Francis Dr. **Santa Fe, NM 87505**

QUESTI	ONS (continued)	
Operator:		OGRID:
Harvest Four Corners, LLC 1111 Travis Street		373888 Action Number:
Houston, TX 77002		145865
		Action Type: [C-129] Amend Venting and/or Flaring (C-129A)
QUESTIONS		
Date(s) and Time(s)		
Date vent or flare was discovered or commenced	09/15/2022	
Time vent or flare was discovered or commenced	12:32 PM	
Time vent or flare was terminated	01:15 PM	
Cumulative hours during this event	285	
Measured or Estimated Volume of Vented or Flared Natural Gas		
Network Con Virginia (AMA) Data il	Cause: Corrosion I	Pipeline (Any) Natural Gas Vented Released: 2,160 Mcf Recovered: 0
Natural Gas Vented (Mcf) Details	Mcf Lost: 2,160 Mc	
Natural Gas Flared (Mcf) Details	Not answered.	
Other Released Details	Cause: Other (Sp	pecify) Released: 0 (Unknown Released Amount) Recovered: 0 Lost: 0
Additional details for Measured or Estimated Volume(s). Please specify	Not answered.	
Is this a gas only submission (i.e. only significant Mcf values reported)	Yes, according to s	supplied volumes this appears to be a "gas only" report.
Venting or Flaring Resulting from Downstream Activity		
	T	
Was this vent or flare a result of downstream activity	No	
Was notification of downstream activity received by this operator	Not answered.	
Downstream OGRID that should have notified this operator Date notified of downstream activity requiring this vent or flare	Not answered.	
Time notified of downstream activity requiring this vent or flare	Not answered.	
Time notined of downstream activity requiring this vent of hare	Not answered.	
Steps and Actions to Prevent Waste		
For this event, this operator could not have reasonably anticipated the current event and it was beyond this operator's control	True	
Please explain reason for why this event was beyond this operator's control	Corrosion	
Steps taken to limit the duration and magnitude of vent or flare	Pipeline was shut i	n
Corrective actions taken to eliminate the cause and reoccurrence of vent or flare	Pipeline was repair	red

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ACKNOWLEDGMENTS

Action 145865

ACKNOWLEDGMENTS

Operator:	OGRID:
Harvest Four Corners, LLC	373888
1111 Travis Street	Action Number:
Houston, TX 77002	145865
	Action Type:
]	[C-129] Amend Venting and/or Flaring (C-129A)

ACKNOWLEDGMENTS

V	I acknowledge that with this application I will be amending an existing incident file (assigned to this operator) for a vent or flare event, pursuant to 19.15.27 and 19.15.28 NMAC	
V	I acknowledge that amending an incident file does not replace original submitted application(s) or information and understand that any C-129 forms submitted to the be logged and stored as public record.	
V	I hereby certify the statements in this amending report are true and correct to the best of my knowledge and acknowledge that any false statement may be subject to civil and criminal penalties under the Oil and Gas Act.	
V	I acknowledge that the acceptance of any C-129 forms by the OCD does not relieve this operator of liability should their operations have failed to adequately investigate, report, and remediate contamination that poses a threat to groundwater, surface water, human health, or the environment.	
V	I acknowledge that OCD acceptance of any C-129 forms does not relieve this operator of responsibility for compliance with any other applicable federal, state, or local laws and/or regulations.	

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State of New Mexico Energy, Minerals and Natural Resources Oil Conservation Division 1220 S. St Francis Dr. **Santa Fe, NM 87505**

CONDITIONS

Action 145865

CONDITIONS

Operator:	OGRID:
Harvest Four Corners, LLC	373888
1111 Travis Street	Action Number:
Houston, TX 77002	145865
	Action Type:
	[C-129] Amend Venting and/or Flaring (C-129A)

CONDITIONS

Created	d Condition	Condition
Ву		Date
jdeal	If the information provided in this report requires further amendment(s), submit a [C-129] Amend Venting and/or Flaring Incident (C-129A), utilizing your incident number from this event.	9/23/2022