Submit within 45 days of well completion		State of New Mexico Energy, Minerals and Natural			Revised February 15, 2012 1. WELL API NO.				
			nerais and Na esources	turai		30-039-26827			
		Oil Conservation Division			2. Well Name: JICARILLA 96 #002C				
1220 S. St Francis Dr.					3. Well Number: 002C				
Santa Fe, NM 87505  HYDRAULIC FRACTURING FLUID DISCLOSURE						Surface Hole Location:     Unit:G Lot:G Section:2 Township:26N Range:03			
O NET CARROLL STATE OF STATE O	IC FRAC	TURING FLU	ID DISCLOSURE		Feet fro	m:1750	N/S Li	ne:N	
						Hole Location			
_ Amendme	III.				Feet from	m:1750	N/S Li		
					6. latitud	de:	E/W Li	gitude:	
					7. County			0	
8. Operator Name an	d Address:				9. OGRID:	Rio Arriba	10 Pho	ne Number:	
	N RESOU	IRCES CORPORA	ATION			162928	10.1110	505-324-4154	
	on 87401	2012 Frac Perform	ed by: HALLIBURTO	J.	12 Produc	tion Type:			
13. Pool Code(s):	- 0/11/	2013 11001 2110111	- TALLIBORTO	•		G ractured Inter	val:		
77360 15. True Vertical Dep	oth (TVD):					3,722 ft to 3		d:	
6,252 ft	entre de la constitue	POSITION AND CO	NCENTRATION:			23,562 gals			
Trade Name	Supplier	Purpose	Ingredients	Abstrac	Chemical t Service	Maximum I Concentrati	on in	Maximum Ingredient	
IIVDDOOLII ODIO	I I a II i b	_	I hadaa ahlaaia a aid	#	14.0	Additive (%	, 14	Fluid (% by mass)	
HYDROCHLORIC ACID 10-30%	200000000000000000000000000000000000000	80000	Hydrochloric acid	7047-0	47-01-0		30% 0.69861		
2% KCL Water SAND - PREMIUM	Operator Halliburto		Crystalline silica,	14808-	-60-7		100%	51.62319 28.666949	
WHITE SSA-2	Halliburto	on Proppant	quartz Crystalline silica,	14808-60-7		100%		1.29959	
LGC-36 UC	Halliburto	on Liquid Gel	quartz Guar gum	9000-30-0 64742-48-9		60% 60%		0.115729	
		Concentrate						0.115729	
BC-140	Halliburto	on Crosslinker	Ethylene glycol Monoethanolamine	26038-			30% 0.014 60% 0.028		
AQF-2 FOAMING	Halliburto	on Foaming	borate Ethylene glycol	111-76				0.028077	
AGENT	riallibult	Agent	monobutyl ether Diethylene glycol	111-46-6		10%		0.037777	
GasPerm 1100	Halliburto		Ethanol	64-17-	5		30%	0.02438%	
		Surfactant	Isopropanol Methanol	67-63-0 67-56-1			10% 10%	0.00813% 0.00813%	
			Terpenes and Terpenoids, sweet	68647-	-72-3		5%	0.00406%	
HAI-OS ACID	Halliburto	on Corrosion	orange-oil Methanol	67-56-1			60%	0.000589	
INHIBITOR GBW-30	Halliburto	Inhibitor on Breaker	Propargyl alcohol Hemicellulase	107-19-7 9012-54-8			10% 30%	0.0001% 0.00117%	
BREAKER			enzyme Carbohydrates	Confid	\$2000 W		100%	0.0039%	
			Carbonyarates	Busine	ess		10070	0.00007	
OPTIFLO-HTE	Halliburto	on Breaker	Crystalline silica,	Information 14808-60-7			30%	0.00195%	
			quartz Walnut hulls	Mixture			100%	0.0065%	
NITROGEN LIQUEFIED	Halliburto	on Fluid	Nitrogen	7727-3			100%	15.62359%	
Ingredients Listed Below This Line Are Part of the			Alcohols, C14-C15, ethoxylated				0%	0%	
			Alkyl sulfonate	Confid Busine	ess		0%	0%	
			C.I. Pigment Red 5	6410-4			0%	0%	
			Crystalline silica, quartz	14808-	-60-7		0%	0%	
			Cured acrylic resin	Confid			0%	09	
			Cured acrylic resin	Information Confidential			0%	0%	
				Business Information				1000	
			Enzyme	Confid	ential		0%	0%	
			Ethoxylated fatty	Information			0%	0%	
			acid	Confidential Business			0 70	07	
			Ethoxylated fatty	Information Confidential			0%	0%	
			acid	Business Information					
			Ethoxylated fatty acid	Confid	ess		0%	09	
			Fatty acids, tall oil	Inform: Confid	ential		0%	09	
				Busine Informa	ation				
			Fatty alcohol polyglycol ether	9043-3	30-5		0%	09	
			surfactant Methyl Isobutyl	108-10	)-1		0%	09	
			Ketone Olefins	Confid			0%	09	
				Busine	ess		0 70	37	
			Olefins	Confid	ential		0%	09	
			Olofina	Informa	ation		000		
			Olefins	Busine	ess		0%	09	
l.			Olefins	Confid	ential		0%	09	
				Busine Informa	ation				
			Quaternary ammonium	68953-	-58-2		0%	09	
			compounds, bis (hydrogenated					1	
			(hydrogenated tallow alkyl) dimethyl,salts with bentonite Reaction product of acetophenone, formaldehyde, thiourea and oleic acid in dimethyl	68527-	49-1		0%	09	
			(hydrogenated tallow alkyl) dimethyl,salts with bentonite Reaction product of acetophenone, formaldehyde, thiourea and oleic	68527- Confid			0%		
			(hydrogenated tallow alkyl) dimethyl,salts with bentonite Reaction product of acetophenone, formaldehyde, thiourea and oleic acid in dimethyl formamide		ential			0%	

NMOCD does not require the reporting of information beyond MSDS data as described in 29 CFR 1910.1200. NMOCD does not require the reporting or disclosure of proprietary, trade secret or confidential business information.