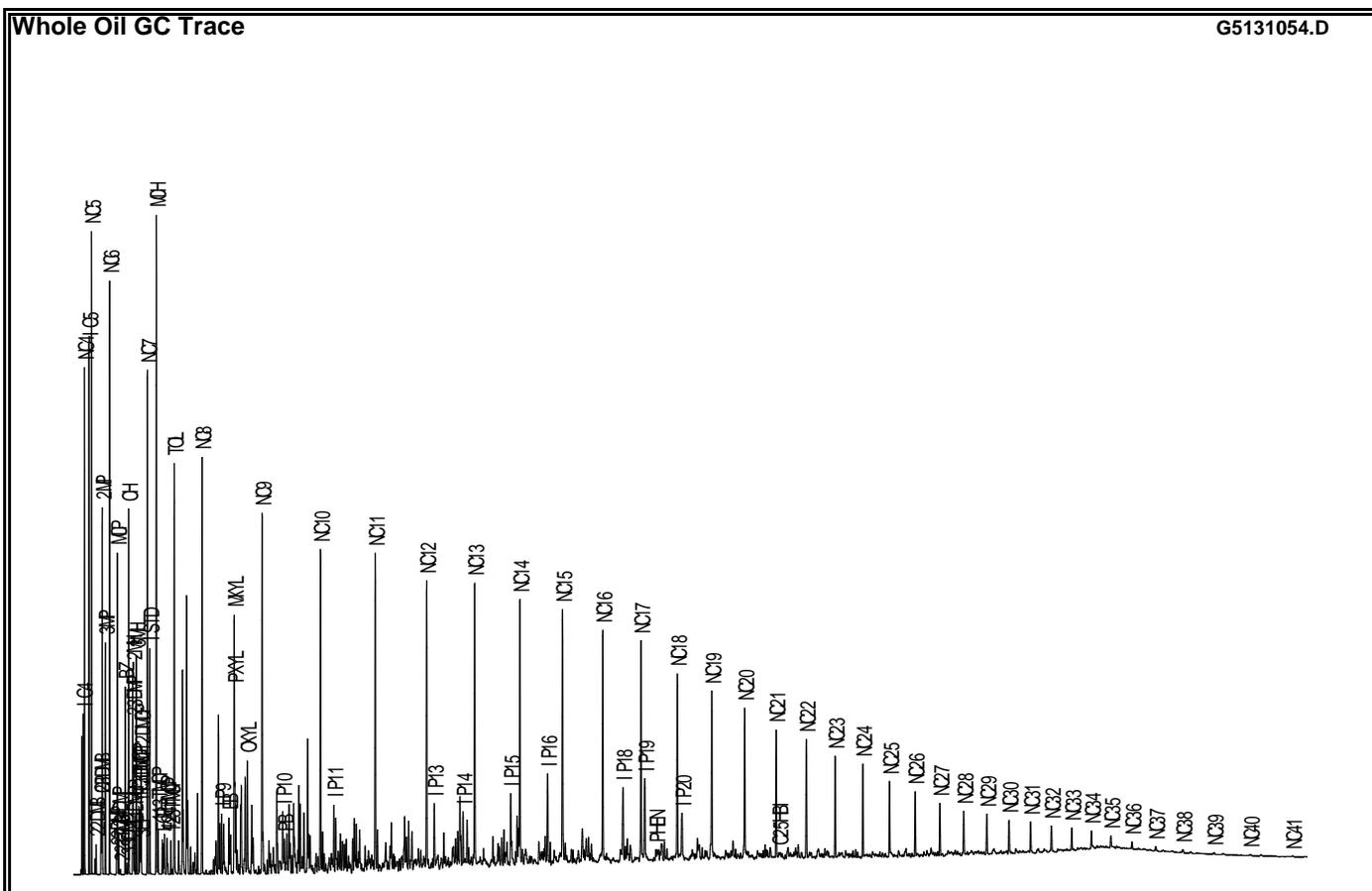


Company:	Department of Health and Human Service	Client ID:	
Country:		Project #:	BH-65414
Basin:		Lab ID:	3403042751
Lease:		Sample Type:	Oil
Block:		Sampling Point:	
Field:		Formation:	
Well Name:	NIOSH Fedan	Geologic Age:	
Latitude:		Top Depth:	
Longitude:		Bottom Depth:	
Preparation:		GC Method:	G02G



WGC parameters	
Pristane/Phytane	1.67
Pristane/nC <sub>17</sub>	0.55
Phytane/nC <sub>18</sub>	0.39
nC <sub>18</sub> /(nC <sub>18</sub> +nC <sub>19</sub> )	0.52
nC <sub>17</sub> /(nC <sub>17</sub> +nC <sub>29</sub> )	0.82
CPI Hunt <sup>4</sup>	1.01
Normal Paraffins	29.6
Isoprenoids	4.4
Cycloparaffins	6.4
Branched (iso-) Paraffins	5.4
BTX aromatics	4.4
Resolved unknowns	49.3

Thompson <sup>1</sup>	
A. BZ/nC <sub>6</sub>	0.38
B. TOL/nC <sub>7</sub>	0.93
C. (nC <sub>6</sub> +nC <sub>7</sub> )/(CH+MCH)	0.92
I. Isoheptane Value	1.69
F. nC <sub>7</sub> /MCH	0.71
U. CH/MCP	1.24
R. nC <sub>7</sub> /2MH	2.59
S. nC <sub>6</sub> /22DMB	21.90
H. Heptane Value	21.86
MCH/nC <sub>7</sub>	1.41
mpXYL/nC <sub>8</sub>	0.81

Mango <sup>2</sup>	
P <sub>1</sub>	19.75
P <sub>2</sub>	15.80
P <sub>3</sub>	5.75
5N <sub>1</sub>	5.69
N <sub>2</sub>	6.66
6N <sub>1</sub>	46.36
K <sub>1</sub>	1.06
K <sub>2</sub>	0.70
5N <sub>1</sub> /6N <sub>1</sub>	0.12
P <sub>3</sub> /N <sub>2</sub>	0.86
ln(24DMP/23DMP)	-0.58

Halpern <sup>3</sup>	
Tr <sub>1</sub>	11.92
Tr <sub>2</sub>	12.77
Tr <sub>3</sub>	5.29
Tr <sub>4</sub>	4.94
Tr <sub>5</sub>	10.22
Tr <sub>7</sub>	1.62
Tr <sub>8</sub>	2.85
C <sub>1</sub>	0.12
C <sub>2</sub>	0.45
C <sub>3</sub>	0.25
C <sub>4</sub>	0.08
C <sub>5</sub>	0.10

<sup>1</sup>Thompson, K.F.M.,1983.GCA:V.47, p.303. <sup>2</sup>Mango, F.D.,1994.GCA: V.58, p.895. <sup>3</sup>Halpern,H.L.,1995.AAPG Bull.: V.79, p.801. <sup>4</sup>Hunt, 1979

<b>Company:</b>	<b>Department of Health and Human Services</b>	<b>Client ID:</b>	
<b>Well Name:</b>	<b>NIOSH Fedan</b>	<b>Project #:</b>	<b>BH-65414</b>
<b>Depth:</b>	<b>-</b>	<b>Lab ID:</b>	<b>3403042751</b>
<b>Sampling Point:</b>		<b>File Name:</b>	<b>G5131054.D</b>

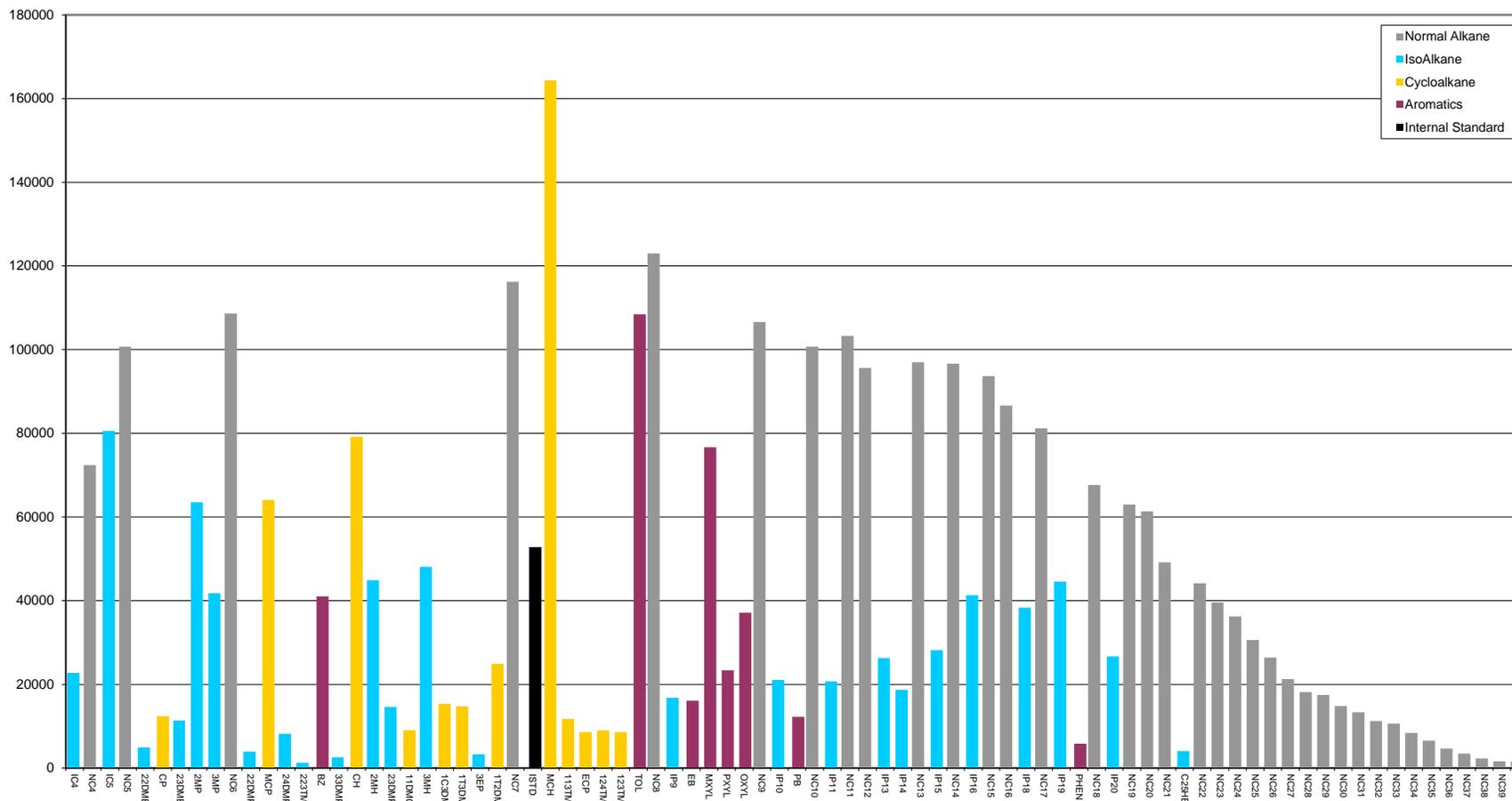
Peak Label	Compound Name	Ret. Time	Area	Height	ppt (Area)	ppt (Hght)
IC4	Iso-alkane C4	3.907	22745	16171	1.48	2.43
NC4	Normal Alkane C4	4.024	72395	51046	4.70	7.68
IC5	Iso-alkane C5	4.459	80606	53871	5.23	8.10
NC5	Normal Alkane C5	4.698	100722	65036	6.54	9.78
22DMB	2,2-Dimethylbutane	5.150	4960	3000	0.32	0.45
CP	Cyclopentane	5.620	12374	7424	0.80	1.12
23DMB	2,3-Dimethylbutane	5.652	11389	6600	0.74	0.99
2MP	2-Methylpentane	5.735	63499	36967	4.12	5.56
3MP	3-Methylpentane	6.033	41786	23350	2.71	3.51
NC6	Normal Alkane C6	6.447	108631	59761	7.05	8.99
22DMP	2,2-Dimethylpentane	7.122	3885	2102	0.25	0.32
MCP	Methylcyclopentane	7.182	64031	32436	4.16	4.88
24DMP	2,4-Dimethylpentane	7.316	8211	4035	0.53	0.61
223TMB	2,2,3-Trimethylbutane	7.473	1247	534	0.08	0.08
BZ	Benzene	7.942	41053	18956	2.66	2.85
33DMP	3,3-Dimethylpentane	8.158	2578	1159	0.17	0.17
CH	Cyclohexane	8.277	79198	36814	5.14	5.54
2MH	2-Methylhexane	8.663	44917	20781	2.92	3.13
23DMP	2,3-Dimethylpentane	8.714	14622	7188	0.95	1.08
11DMCP	1,1-Dimethylcyclopentane	8.803	9099	4025	0.59	0.61
3MH	3-Methylhexane	9.008	48094	22019	3.12	3.31
1C3DMCP	1-cis-3-Dimethylcyclopentane	9.241	15371	6883	1.00	1.04
1T3DMCP	1-trans-3-Dimethylcyclopentane	9.356	14724	6571	0.96	0.99
3EP	3-Ethylpentane	9.419	3290	2211	0.21	0.33
1T2DMCP	1-trans-2-Dimethylcyclopentane	9.472	24904	10874	1.62	1.64
NC7	Normal Alkane C7	10.077	116217	50816	7.54	7.64
ISTD	Internal Standard	10.293	52854	22809	3.43	3.43
MCH	Methylcyclohexane	10.937	164401	66357	10.67	9.98
113TMCP	1,1,3,-Trimethylcyclopentane	11.101	11768	4486	0.76	0.68
ECP	Ethylcyclopentane	11.510	8570	3505	0.56	0.53
124TMCP	1,2,4-Trimethylcyclopentane	11.962	9024	3653	0.59	0.55
123TMCP	1,2,3-Trimethylcyclopentane	12.334	8563	3408	0.56	0.51
TOL	Toluene	12.653	108454	41465	7.04	6.24
NC8	Normal Alkane C8	15.328	122975	41974	7.98	6.31
IP9	Isoprenoid C9	17.203	16786	6062	1.09	0.91
EB	Ethyl-benzene	17.905	16143	5680	1.05	0.85
MXYL	<i>m</i> -xylene	18.415	76688	26116	4.98	3.93
PXYL	<i>p</i> -xylene	18.473	23372	9272	1.52	1.39
OXYL	<i>o</i> -xylene	19.671	37097	11422	2.41	1.72
NC9	Normal Alkane C9	21.091	106624	36416	6.92	5.48
IP10	Isoprenoid C10	23.069	21023	6290	1.36	0.95
PB	Propyl-benzene	23.240	12259	3519	0.80	0.53
NC10	Normal Alkane C10	26.699	100736	32822	6.54	4.94
IP11	Isoprenoid C11	27.992	20735	6898	1.35	1.04
NC11	Normal Alkane C11	31.969	103273	32272	6.70	4.85
NC12	Normal Alkane C12	36.900	95613	29156	6.21	4.38
IP13	Isoprenoid C13	37.634	26296	6650	1.71	1.00
IP14	Isoprenoid C14	40.401	18675	5707	1.21	0.86
NC13	Normal Alkane C13	41.520	96981	28716	6.29	4.32
IP15	Isoprenoid C15	44.969	28151	7363	1.83	1.11
NC14	Normal Alkane C14	45.863	96617	26904	6.27	4.05



**Company:** Department of Health and Human Services  
**Well Name:** N10SH Fedan  
**Depth:** -  
**Sampling Point:**

**Client ID:**  
**Project #:** BH-65414  
**Lab ID:** 3403042751  
**File Name:** G5131054.D

Histogram Based on Area



Parameter	Formula
<b>WGC Parameters</b>	
Pristane/Phytane	IP19/IP20
Pristane/nC17	IP19/NC17
Phytane/nC18	IP20/NC18
nC <sub>18</sub> /(nC <sub>18</sub> +nC <sub>19</sub> )	NC18/(NC18+NC19)
nC <sub>17</sub> /(nC <sub>17</sub> +nC <sub>29</sub> )	NC17/(NC17+NC29)
CPI Hunt <sup>4</sup>	$((NC23+NC25+NC27+NC29+NC31)+(NC25+NC27+NC29+NC31+NC33))/(2*(NC24+NC26+NC28+NC30+NC32))$
Normal Paraffins	$100*(NC4+NC5+NC6+NC7+NC8+NC9+NC10+NC11+NC12+NC13+NC14+NC15+NC16+NC17+NC18+NC19+NC20+NC21+NC22+NC23+NC24+NC25+NC26+NC27+NC28+NC29+NC30+NC31+NC32+NC33+NC34+NC35+NC36+NC37+NC38+NC39+NC40+NC41)/TOTAL\_RESOLVED$
Isoprenoids	$100*(IP9+IP10+IP11+IP13+IP14+IP15+IP16+IP18+IP19+IP20+C25HBI)/TOTAL\_RESOLVED$
Cycloparaffins	$100*(CP+MCP+CH+11DMCP+1T3DMCP+1T2DMCP+MCH+113TMCP+ECP+124TMCP+123TMCP+1C3DMCP+1C2DMCP)/TOTAL\_RESOLVED$
Branched (iso-) Paraffins	$100*(IC4+IC5+22DMB+23DMB+2MP+3MP+22DMP+24DMP+223TMB+33DMP+2MH+23DMP+3MH+3EP)/TOTAL\_RESOLVED$
BTX aromatics	$100*(BZ+TOL+MXYL+PXYL+OXYL)/TOTAL\_RESOLVED$ (*OXYL added 05/21/2007)
<b>Thompson<sup>1</sup></b>	
BZ/nC6	BZ/NC6
TOL/nC7	TOL/NC7
(nC6+nC7)/(CH+MCH)	(NC6+NC7)/(CH+MCH)
Isoheptane Value	$(2MH+3MH)/(1C3DMCP+1T3DMCP+1T2DMCP)$
nC7/MCH	NC7/MCH
CH/MCP	CH/MCP
nC7/2MH	NC7/2MH
nC6/22DMB	NC6/22DMB
Heptane Value	$100*NC7/(CH+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+1T2DMCP+MCH+NC7)$
MCH/nC7	MCH/NC7
mpXYL/nC8	(MXYL+PXYL)/NC8
<b>Mango<sup>2</sup></b>	
P1	$100*NC7/(22DMP+24DMP+223TMB+33DMP+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+3EP+1T2DMCP+NC7+MCH+ECP+TOL)$
P2	$100*(2MH+3MH)/(22DMP+24DMP+223TMB+33DMP+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+3EP+1T2DMCP+NC7+MCH+ECP+TOL)$
P3	$100*(3EP+33DMP+23DMP+24DMP+22DMP+223TMB)/(22DMP+24DMP+223TMB+33DMP+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+3EP+1T2DMCP+NC7+MCH+ECP+TOL)$
5N1	$100*(ECP+1T2DMCP)/(22DMP+24DMP+223TMB+33DMP+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+3EP+1T2DMCP+NC7+MCH+ECP+TOL)$
N2	$100*(11DMCP+1C3DMCP+1T3DMCP)/(22DMP+24DMP+223TMB+33DMP+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+3EP+1T2DMCP+NC7+MCH+ECP+TOL)$
6N1	$100*(MCH+TOL)/(22DMP+24DMP+223TMB+33DMP+2MH+23DMP+11DMCP+3MH+1C3DMCP+1T3DMCP+3EP+1T2DMCP+NC7+MCH+ECP+TOL)$
K1	$(2MH+23DMP)/(3MH+24DMP)$
K2	$(2MH+3MH)/(2MH+3MH+11DMCP+1C3DMCP+1T3DMCP)$
5N1/6N1	$(ECP+1T2DMCP)/(MCH+TOL)$
P3/N2	$(3EP+33DMP+23DMP+24DMP+22DMP+223TMB)/(11DMCP+1C3DMCP+1T3DMCP)$
ln(24DMP/23DMP)	ln(24DMP/23DMP)
<b>Halpern<sup>3</sup></b>	
Tr1	TOL/11DMCP
Tr2	NC7/11DMCP
Tr3	3MH/11DMCP
Tr4	2MH/11DMCP
Tr5	$(2MH+3MH)/11DMCP$
Tr7	1T3DMCP/11DMCP
Tr8	$(2MH+3MH)/(22DMP+23DMP+24DMP+33DMP+3EP)$
C1	$22DMP/(22DMP+23DMP+24DMP+33DMP+3EP)$
C2	$23DMP/(22DMP+23DMP+24DMP+33DMP+3EP)$
C3	$24DMP/(22DMP+23DMP+24DMP+33DMP+3EP)$
C4	$33DMP/(22DMP+23DMP+24DMP+33DMP+3EP)$
C5	$3EP/(22DMP+23DMP+24DMP+33DMP+3EP)$

1	2	3	4	5	6
INFO	Name	Range Start	Range End	Add Next sample into next	Sample display distance
PAGE	SUMMARY	A1	BE174	workbook	
PAGE	Histogram	A1	I43	workbook	
FILE				GC_WHOIL_[F]_[S]_[D]-[B]_[P].xls	

7	8	9	10	11
<b>Max samples in one page</b>	<b>Tick</b>			