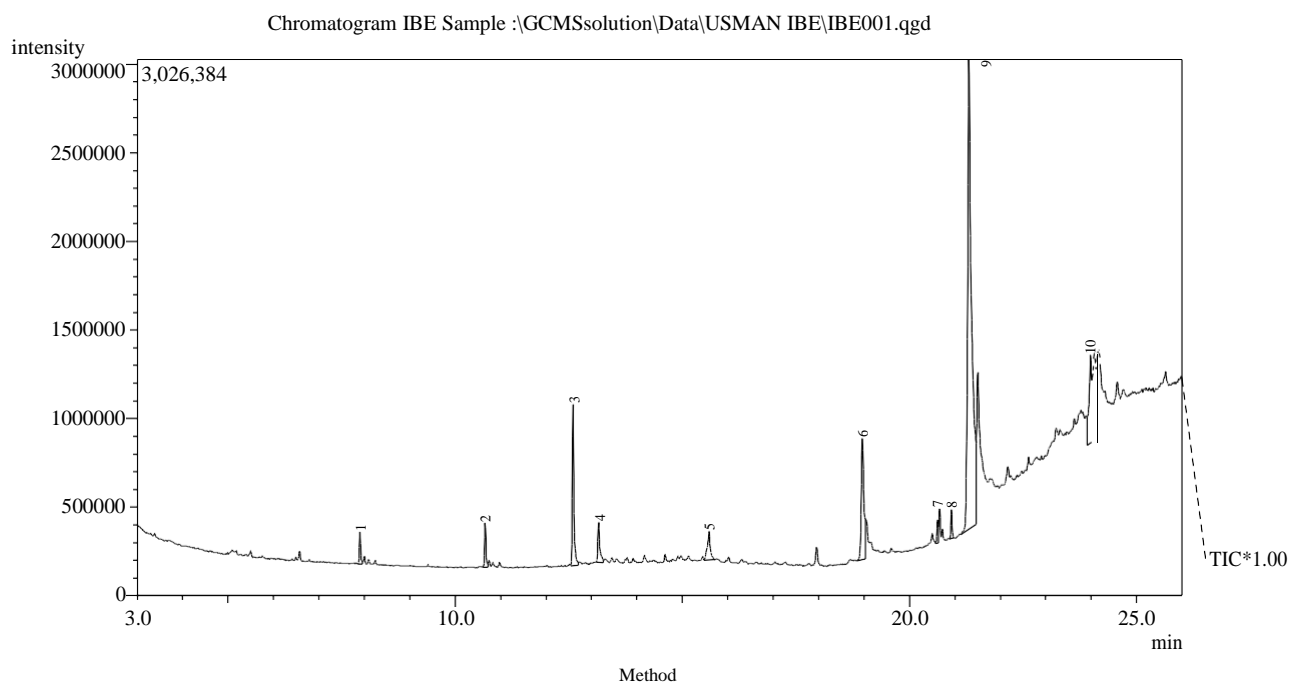


GCMS-QP2010 PLUS  
SHIMADZU,JAPAN

USMAN IBE

SAMPLE: 1E



[Comment]

===== Analytical Line 1 =====

[AOC-20i]

# of Rinses with Presolvent	4
# of Rinses with Solvent(post)	4
# of Rinses with Sample	3
Plunger Speed(Suction)	:High
Viscosity Comp. Time	:0.2 sec
Plunger Speed(Injection)	:High
Syringe Insertion Speed	:High
Injection Mode	:Normal
Pumping Times	5
Inj. Port Dwell Time	:0.3 sec
Terminal Air Gap	:No
Plunger Washing Speed	:High
Washing Volume	:8uL
Syringe Suction Position	:0.0 mm
Syringe Injection Position	:0.0 mm
Use 3 Solvent Vial	:1 vial

[GC-2010]

Column Oven Temp.	:70.0 °C	
Injection Temp.	:250.00 °C	
Injection Mode	:Split	
Flow Control Mode	:Linear Velocity	
Pressure	:104.1 kPa	
Total Flow	:6.2 mL/min	
Column Flow	:1.59 mL/min	
Linear Velocity	:46.3 cm/sec	
Purge Flow	:3.0 mL/min	
Split Ratio	:1.0	
High Pressure Injection	:OFF	
Carrier Gas Saver	:OFF	
Splitter Hold	:OFF	
Oven Temp. Program		
Rate	Temperature(°C)	Hold Time(min)
-	70.0	1.00
10.00	200.0	4.00
15.00	280.0	4.00

< Ready Check Heat Unit >  
 Column Oven : Yes  
 SPL2 : Yes  
 MS : Yes  
 < Ready Check Detector(FTD) >  
 < Ready Check Baseline Drift >  
 < Ready Check Injection Flow >  
 SPL2 Carrier : Yes  
 SPL2 Purge : Yes  
 < Ready Check APC Flow >  
 < Ready Check Detector APC Flow >  
 External Wait :No  
 Equilibrium Time :3.0 min

[GC Program]

[GCMS-QP2010 Plus]  
 IonSourceTemp :200.00 °C  
 Interface Temp. :250.00 °C  
 Solvent Cut Time :2.50 min  
 Detector Gain Mode :Relative  
 Detector Gain :0.00 kV  
 Threshold :2000

[MS Table]

--Group 1 - Event 1--  
 Start Time :3.00min  
 End Time :26.00min  
 ACQ Mode :Scan  
 Event Time :0.50sec  
 Scan Speed :1666  
 Start m/z :40.00  
 End m/z :800.00

Sample Inlet Unit :GC

[MS Program]

Use MS Program :OFF

Peak Report TIC									
Peak#	R.Time	I.Time	F.Time	Area	Area%	Height	Height%	A/H Mark	Name
1	7.905	7.858	7.958	355409	1.36	179787	3.07	1.98	
2	10.661	10.617	10.717	526759	2.02	249532	4.26	2.11	V
3	12.592	12.542	12.708	2212537	8.48	911695	15.57	2.43	V
4	13.158	13.100	13.267	664941	2.55	225057	3.84	2.95	
5	15.590	15.475	15.725	715510	2.74	161138	2.75	4.44	V
6	18.963	18.850	19.033	2880244	11.03	684899	11.70	4.21	V
7	20.623	20.575	20.642	296033	1.13	130525	2.23	2.27	V
8	20.925	20.883	20.975	330220	1.27	162722	2.78	2.03	
9	21.311	21.133	21.467	16087281	61.63	2649340	45.25	6.07	
10	23.985	23.908	24.017	2032274	7.79	499662	8.53	4.07	V
				26101208	100.00	5854357	100.00		

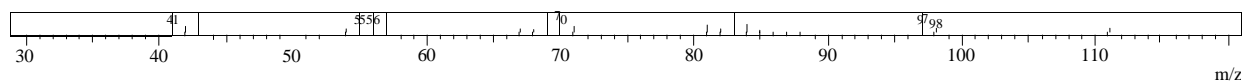
Spectrum

Line#:1 R.Time:7.9(Scan#:590)

MassPeaks:25

RawMode:Single 7.9(590) BasePeak:55(6913)

BG Mode:7.9(591) Group 1 - Event 1



Line#:2 R.Time:10.7(Scan#:920)

MassPeaks:28

RawMode:Single 10.7(920) BasePeak:55(11684)

BG Mode:10.7(922) Group 1 - Event 1

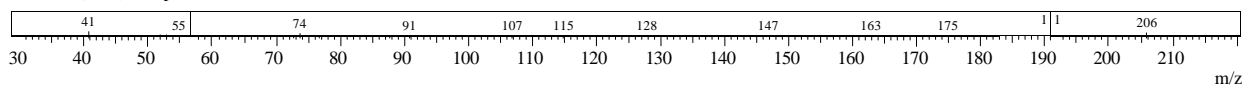


Line#:3 R.Time:12.6(Scan#:1152)

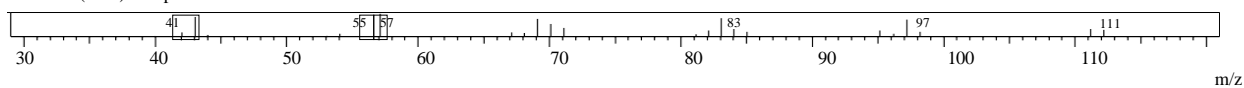
MassPeaks:72

RawMode:Single 12.6(1152) BasePeak:191(156549)

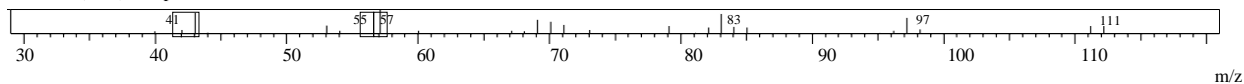
BG Mode:12.6(1156) Group 1 - Event 1



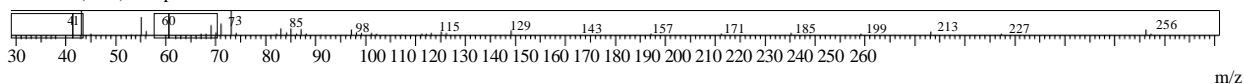
Line#:4 R.Time:13.2(Scan#:1220)  
MassPeaks:27  
RawMode:Single 13.2(1220) BasePeak:55(9576)  
BG Mode:13.2(1222) Group 1 - Event 1



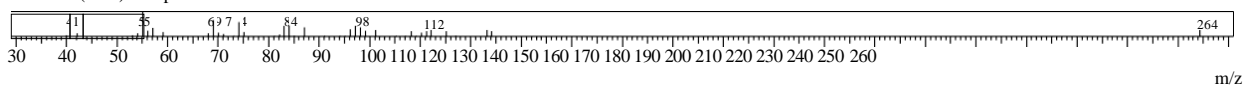
Line#:5 R.Time:15.6(Scan#:1512)  
MassPeaks:30  
RawMode:Single 15.6(1512) BasePeak:57(6583)  
BG Mode:15.6(1514) Group 1 - Event 1



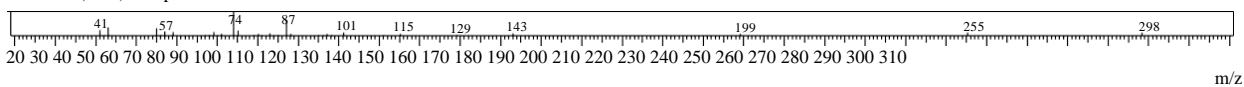
Line#:6 R.Time:19.0(Scan#:1917)  
MassPeaks:62  
RawMode:Single 19.0(1917) BasePeak:73(36097)  
BG Mode:19.0(1922) Group 1 - Event 1



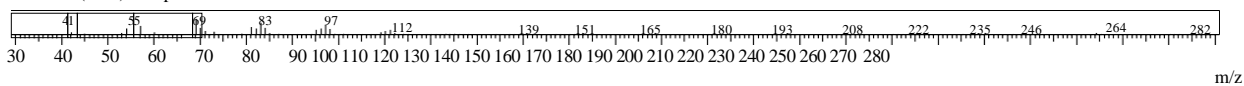
Line#:7 R.Time:20.7(Scan#:2121)  
MassPeaks:48  
RawMode:Single 20.7(2121) BasePeak:55(10434)  
BG Mode:20.6(2118) Group 1 - Event 1



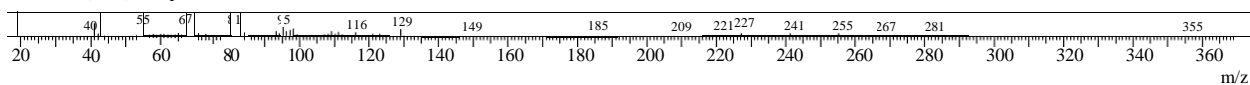
Line#:8 R.Time:20.9(Scan#:2152)  
MassPeaks:44  
RawMode:Single 20.9(2152) BasePeak:74(27083)  
BG Mode:20.9(2155) Group 1 - Event 1



Line#:9 R.Time:21.3(Scan#:2198)  
MassPeaks:105  
RawMode:Single 21.3(2198) BasePeak:55(152013)  
BG Mode:21.4(2206) Group 1 - Event 1



Line#:10 R.Time:24.0(Scan#:2519)  
MassPeaks:89  
RawMode:Single 24.0(2519) BasePeak:55(16852)  
BG Mode:24.0(2524) Group 1 - Event 1

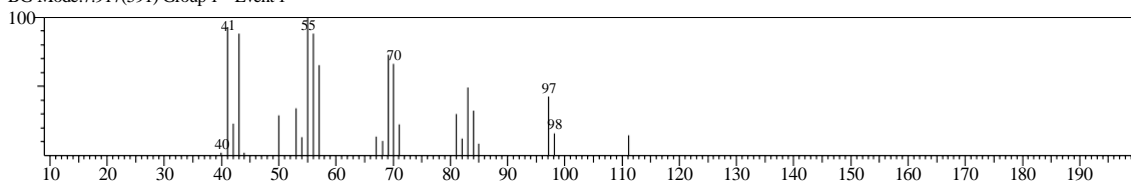


#### Spectrum Comparison

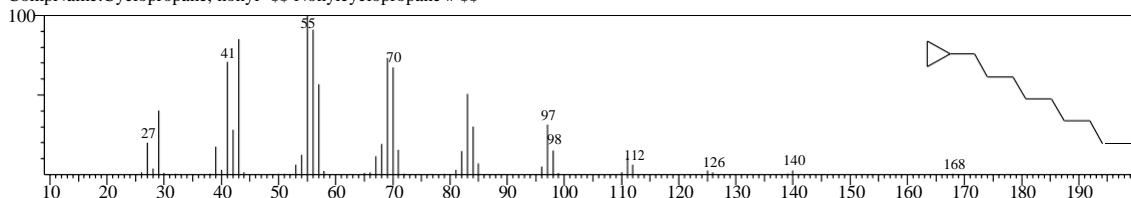
Library

&lt;&lt; Target &gt;&gt;

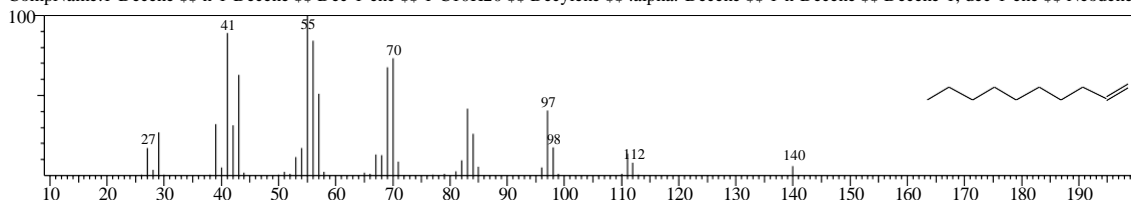
Line#:1 R.Time:7.908(Scan#:590) MassPeaks:25  
RawMode:Single 7.908(590) BasePeak:55.05(6913)  
BG Mode:7.917(591) Group 1 - Event 1



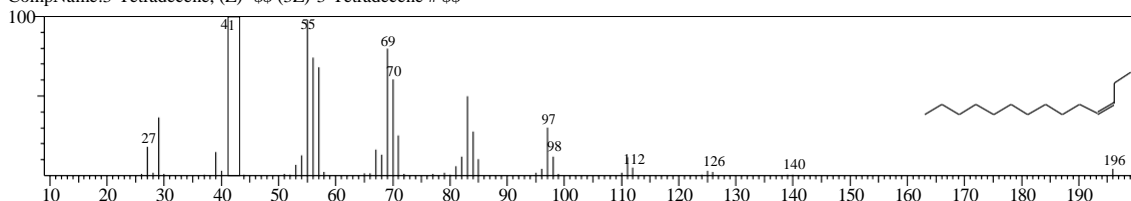
Hit#:1 Entry:23823 Library:NIST05.LIB  
SI:93 Formula:C12H24 CAS:74663-85-7 MolWeight:168 RetIndex:1216  
CompName:Cyclopropane, nonyl- \$\$ Nonylcyclopropane # \$\$



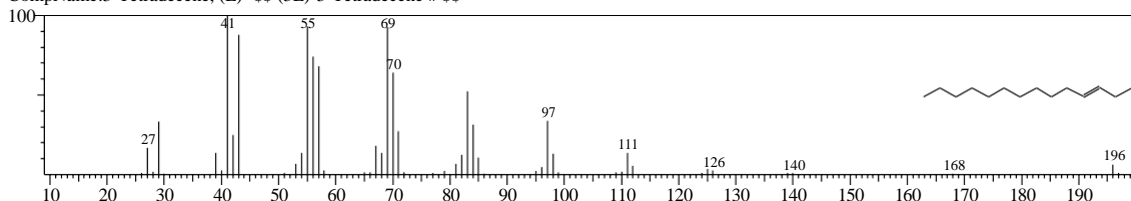
Hit#:2 Entry:6928 Library:NIST05s.LIB  
SI:93 Formula:C10H20 CAS:872-05-9 MolWeight:140 RetIndex:1005  
CompName:1-Decene \$\$ n-1-Decene \$\$ Dec-1-ene \$\$ 1-C10H20 \$\$ Decylene \$\$ .alpha.-Decene \$\$ 1-n-Decene \$\$ Decene-1, dec-1-ene \$\$ Neodene 10 \$\$



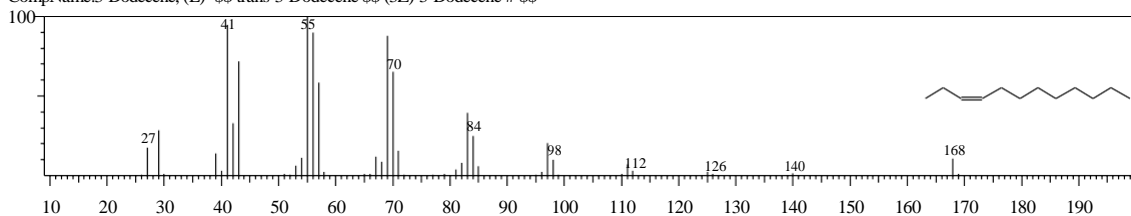
Hit#:3 Entry:39067 Library:NIST05.LIB  
SI:92 Formula:C14H28 CAS:41446-67-7 MolWeight:196 RetIndex:1421  
CompName:3-Tetradecene, (Z)- \$\$ (3Z)-3-Tetradecene # \$\$



Hit#:4 Entry:15745 Library:NIST05s.LIB  
SI:92 Formula:C14H28 CAS:41446-68-8 MolWeight:196 RetIndex:1421  
CompName:3-Tetradecene, (E)- \$\$ (3E)-3-Tetradecene # \$\$

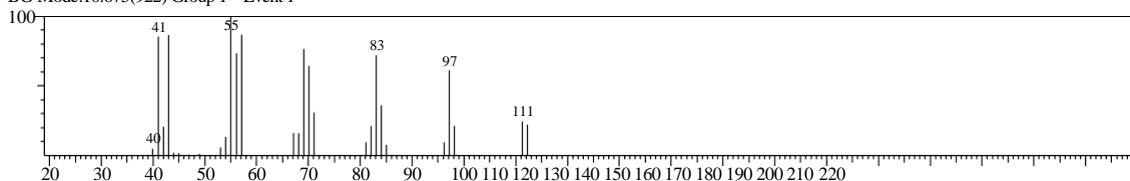


Hit#:5 Entry:23815 Library:NIST05.LIB  
SI:92 Formula:C12H24 CAS:7239-23-8 MolWeight:168 RetIndex:1222  
CompName:3-Dodecene, (E)- \$\$ trans-3-Dodecene \$\$ (3Z)-3-Dodecene # \$\$

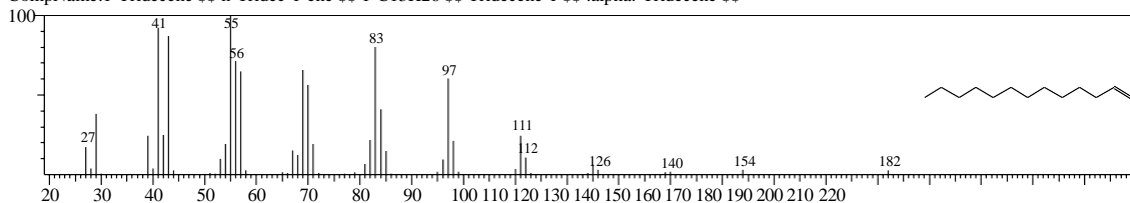


&lt;&lt; Target &gt;&gt;

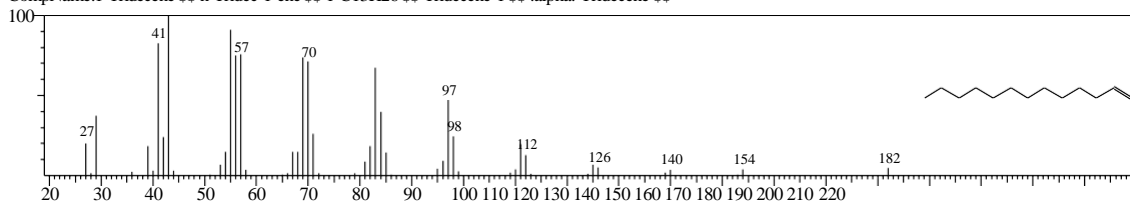
Line#:2 R.Time:10.658(Scan#:920) MassPeaks:28  
RawMode:Single 10.658(920) BasePeak:55.05(11684)  
BG Mode:10.675(922) Group 1 - Event 1



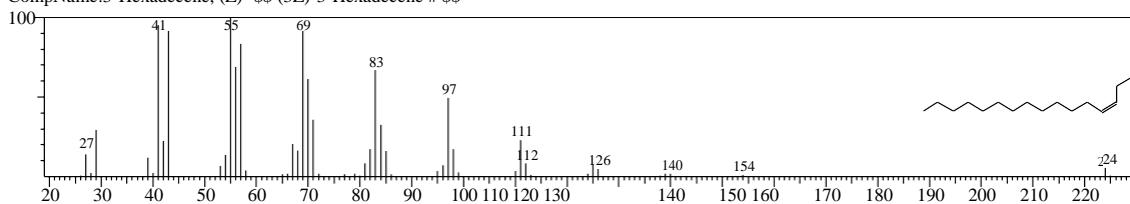
Hit#:1 Entry:13797 Library:NIST05s.LIB  
SI:95 Formula:C<sub>13</sub>H<sub>26</sub> CAS:2437-56-1 MolWeight:182 RetIndex:1304  
CompName:1-Tridecene \$ n-Tridec-1-ene \$ 1-C<sub>13</sub>H<sub>26</sub> \$ Tridecene-1 \$ .alpha.-Tridecene \$



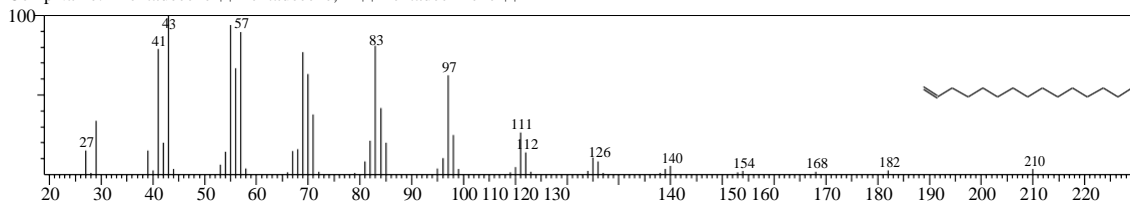
Hit#:2 Entry:31316 Library:NIST05s.LIB  
SI:94 Formula:C<sub>13</sub>H<sub>26</sub> CAS:2437-56-1 MolWeight:182 RetIndex:1304  
CompName:1-Tridecene \$ n-Tridec-1-ene \$ 1-C<sub>13</sub>H<sub>26</sub> \$ Tridecene-1 \$ .alpha.-Tridecene \$



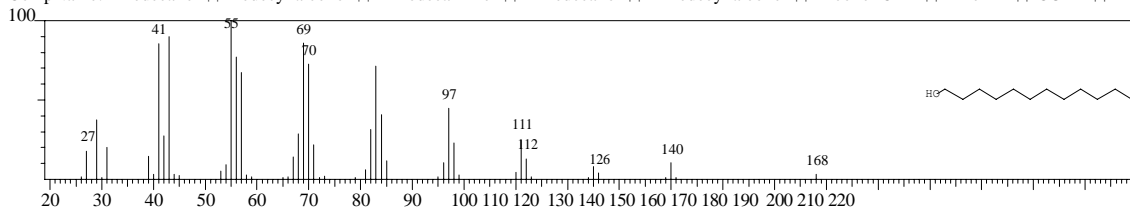
Hit#:3 Entry:55854 Library:NIST05s.LIB  
SI:94 Formula:C<sub>16</sub>H<sub>32</sub> CAS:34303-81-6 MolWeight:224 RetIndex:1620  
CompName:3-Hexadecene, (Z)- \$ (3Z)-3-Hexadecene # \$



Hit#:4 Entry:17440 Library:NIST05s.LIB  
SI:94 Formula:C<sub>15</sub>H<sub>30</sub> CAS:13360-61-7 MolWeight:210 RetIndex:1502  
CompName:1-Pentadecene \$ Pentadecene,1- \$ Pentadec-1-ene \$

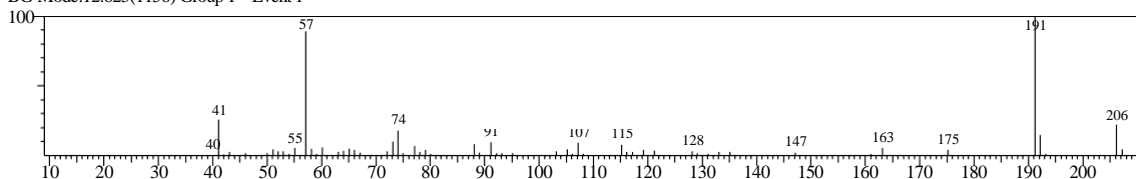


Hit#:5 Entry:14421 Library:NIST05s.LIB  
SI:94 Formula:C<sub>12</sub>H<sub>26</sub>O CAS:112-53-8 MolWeight:186 RetIndex:1457  
CompName:1-Dodecanol \$ Dodecyl alcohol \$ n-Dodecan-1-ol \$ n-Dodecanol \$ n-Dodecyl alcohol \$ Alcohol C-12 \$ Alfol 12 \$ CO 12 \$ Dodecanol-1



&lt;&lt; Target &gt;&gt;

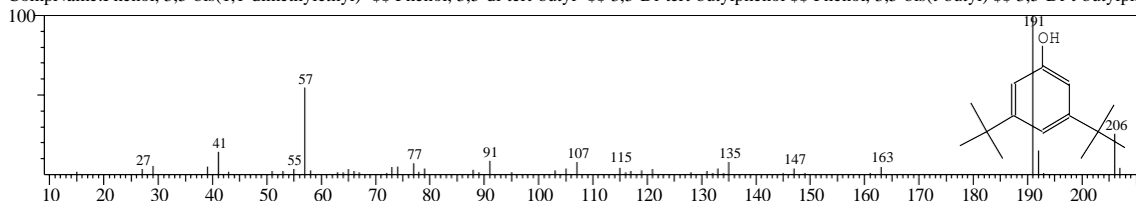
Line#:3 R.Time:12.592(Scan#:1152) MassPeaks:72  
RawMode:Single 12.592(1152) BasePeak:191.20(156549)  
BG Mode:12.625(1156) Group 1 - Event 1



Hit#:1 Entry:16995 Library:NIST05s.LIB

SI:89 Formula:C<sub>14</sub>H<sub>22</sub>O CAS:1138-52-9 MolWeight:206 RetIndex:1555

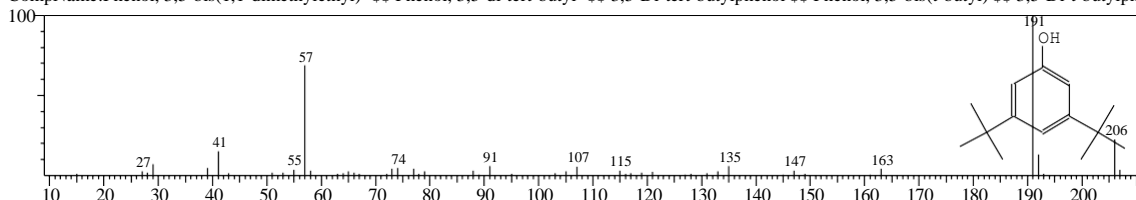
CompName:Phenol, 3,5-bis(1,1-dimethylethyl)- \$ Phenol, 3,5-di-tert-butyl- \$ 3,5-Di-tert-butylphenol \$ Phenol, 3,5-bis(t-butyl) \$ 3,5-Di-t-butylphenol \$



Hit#:2 Entry:16997 Library:NIST05s.LIB

SI:89 Formula:C<sub>14</sub>H<sub>22</sub>O CAS:1138-52-9 MolWeight:206 RetIndex:1555

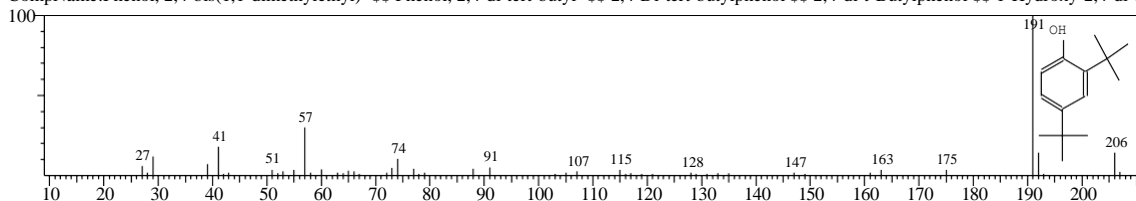
CompName:Phenol, 3,5-bis(1,1-dimethylethyl)- \$ Phenol, 3,5-di-tert-butyl- \$ 3,5-Di-tert-butylphenol \$ Phenol, 3,5-bis(t-butyl) \$ 3,5-Di-t-butylphenol \$



Hit#:3 Entry:16991 Library:NIST05s.LIB

SI:87 Formula:C<sub>14</sub>H<sub>22</sub>O CAS:96-76-4 MolWeight:206 RetIndex:1555

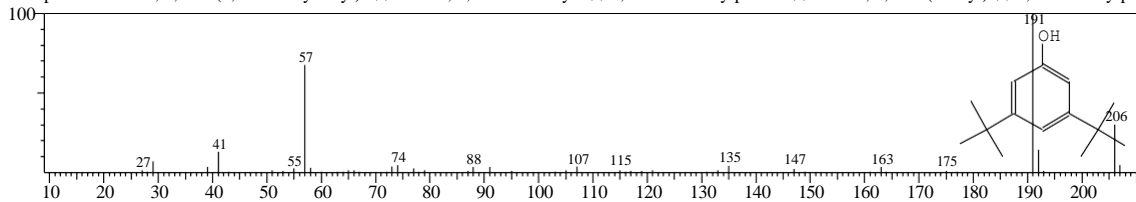
CompName:Phenol, 2,4-bis(1,1-dimethylethyl)- \$ Phenol, 2,4-di-tert-butyl- \$ 2,4-Di-tert-butylphenol \$ 2,4-di-t-Butylphenol \$ 1-Hydroxy-2,4-di-tert-butyl



Hit#:4 Entry:44735 Library:NIST05.LIB

SI:87 Formula:C<sub>14</sub>H<sub>22</sub>O CAS:1138-52-9 MolWeight:206 RetIndex:1555

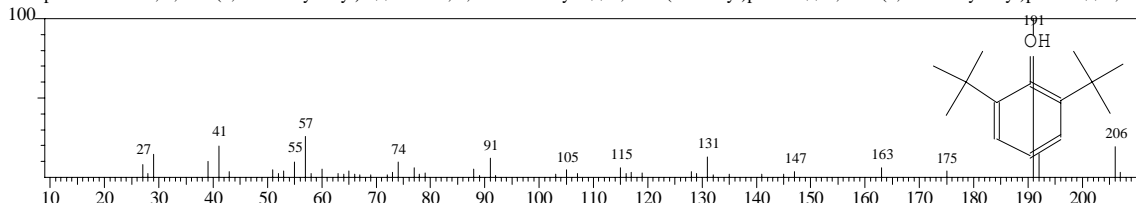
CompName:Phenol, 3,5-bis(1,1-dimethylethyl)- \$ Phenol, 3,5-di-tert-butyl- \$ 3,5-Di-tert-butylphenol \$ Phenol, 3,5-bis(t-butyl) \$ 3,5-Di-t-butylphenol \$



Hit#:5 Entry:16990 Library:NIST05s.LIB

SI:87 Formula:C<sub>14</sub>H<sub>22</sub>O CAS:128-39-2 MolWeight:206 RetIndex:1555

CompName:Phenol, 2,6-bis(1,1-dimethylethyl)- \$ Phenol, 2,6-di-tert-butyl- \$ 2,6-Bis(tert-butyl)phenol \$ 2,6-Bis(1,1-dimethylethyl)phenol \$ 2,6-Di-tert-bu

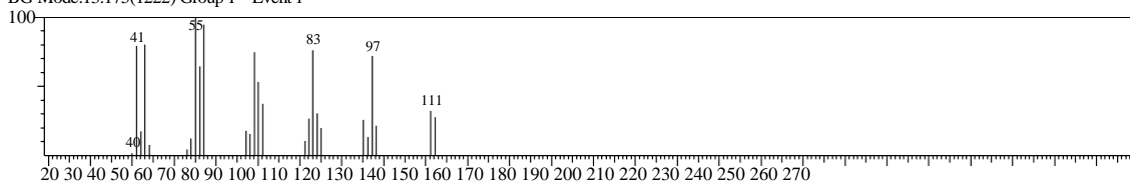


&lt;&lt; Target &gt;&gt;

Line#4 R.Time:13.158(Scan#:1220) MassPeaks:27

RawMode:Single 13.158(1220) BasePeak:55.05(9576)

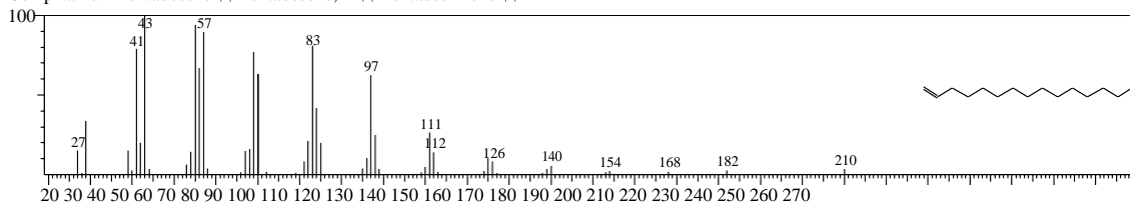
BG Mode:13.175(1222) Group 1 - Event 1



Hit#1 Entry:17440 Library:NIST05s.LIB

SI:94 Formula:C15H30 CAS:13360-61-7 MolWeight:210 RetIndex:1502

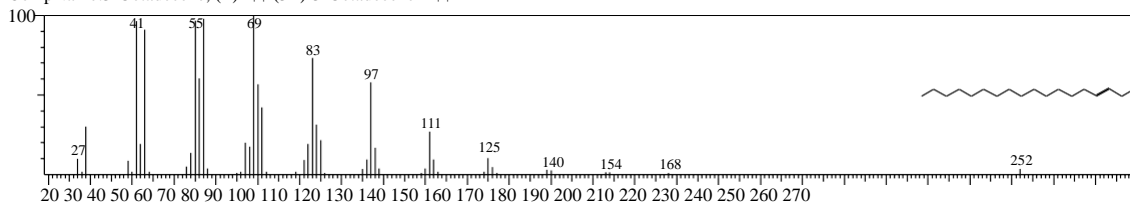
CompName:1-Pentadecene \$\$ Pentadecene,1- \$\$ Pentadec-1-ene \$\$



Hit#2 Entry:72615 Library:NIST05.LIB

SI:93 Formula:C18H36 CAS:7206-19-1 MolWeight:252 RetIndex:1818

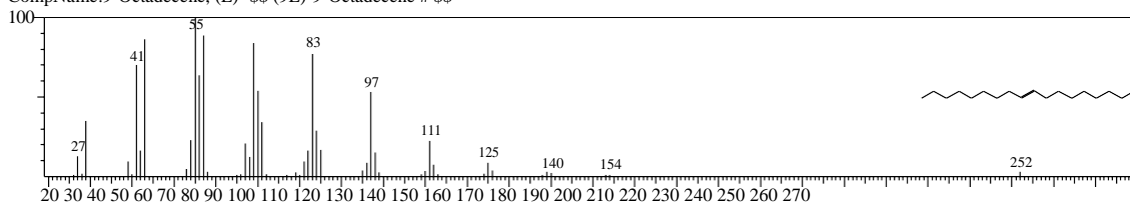
CompName:3-Octadecene, (E)- \$\$ (3E)-3-Octadecene # \$\$



Hit#3 Entry:72607 Library:NIST05.LIB

SI:93 Formula:C18H36 CAS:7206-25-9 MolWeight:252 RetIndex:1818

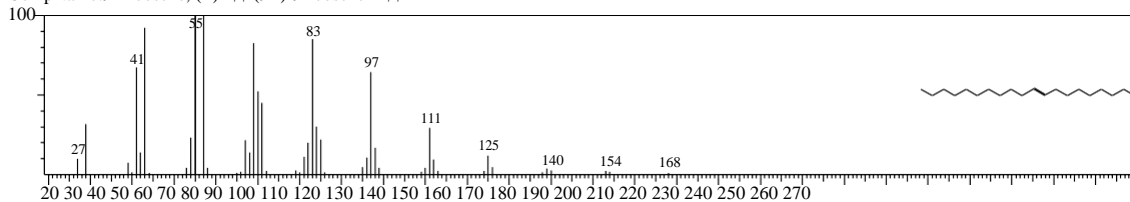
CompName:9-Octadecene, (E)- \$\$ (9E)-9-Octadecene # \$\$



Hit#4 Entry:89386 Library:NIST05.LIB

SI:93 Formula:C20H40 CAS:74685-29-3 MolWeight:280 RetIndex:2017

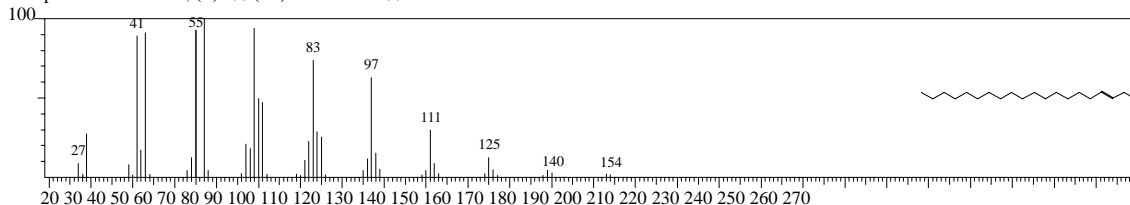
CompName:9-Eicosene, (E)- \$\$ (9E)-9-Eicosene # \$\$



Hit#5 Entry:89387 Library:NIST05.LIB

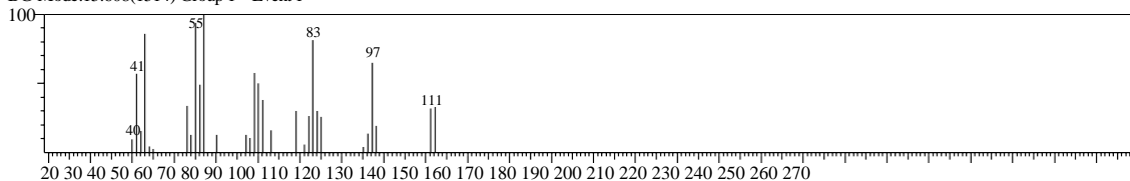
SI:93 Formula:C20H40 CAS:74685-33-9 MolWeight:280 RetIndex:2017

CompName:3-Eicosene, (E)- \$\$ (3E)-3-Eicosene # \$\$

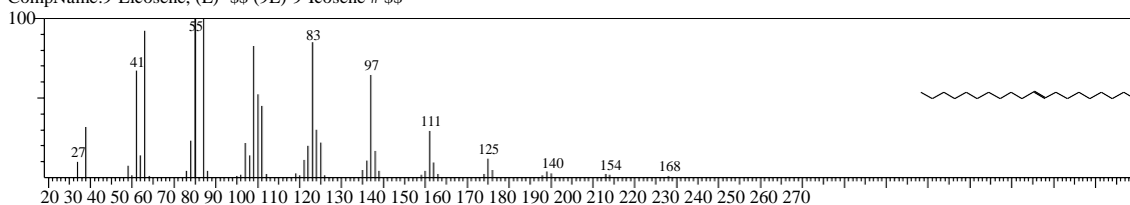


&lt;&lt; Target &gt;&gt;

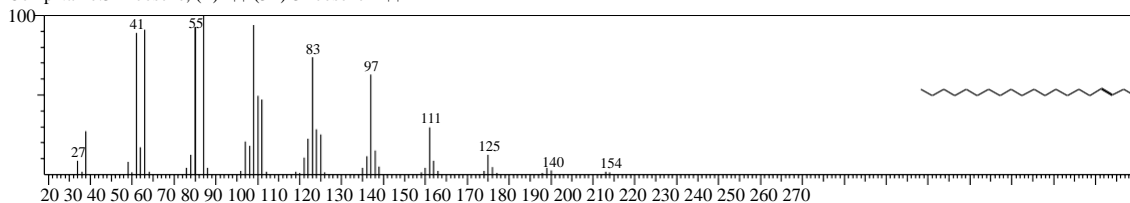
Line#:5 R.Time:15.592(Scan#:1512) MassPeaks:30  
RawMode:Single 15.592(1512) BasePeak:57.10(6583)  
BG Mode:15.608(1514) Group 1 - Event 1



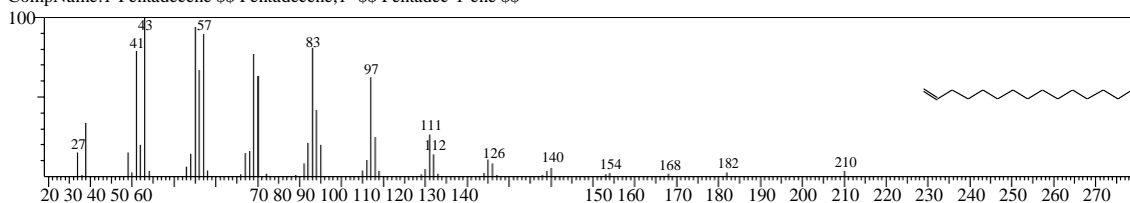
Hit#:1 Entry:89386 Library:NIST05.LIB  
SI:91 Formula:C<sub>20</sub>H<sub>40</sub> CAS:74685-29-3 MolWeight:280 RetIndex:2017  
CompName:9-Eicosene, (E)- \$\$ (9E)-9-Icosene # \$\$



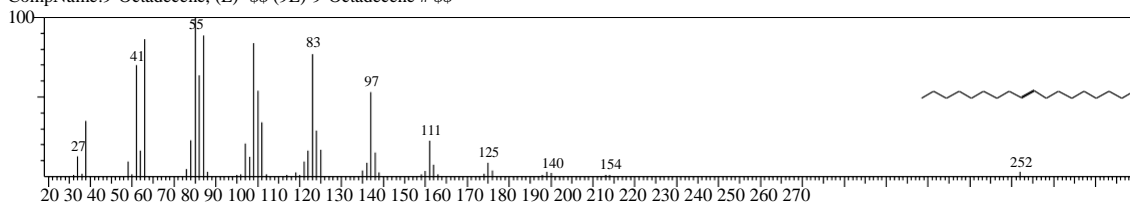
Hit#:2 Entry:89387 Library:NIST05.LIB  
SI:90 Formula:C<sub>20</sub>H<sub>40</sub> CAS:74685-33-9 MolWeight:280 RetIndex:2017  
CompName:3-Eicosene, (E)- \$\$ (3E)-3-Icosene # \$\$



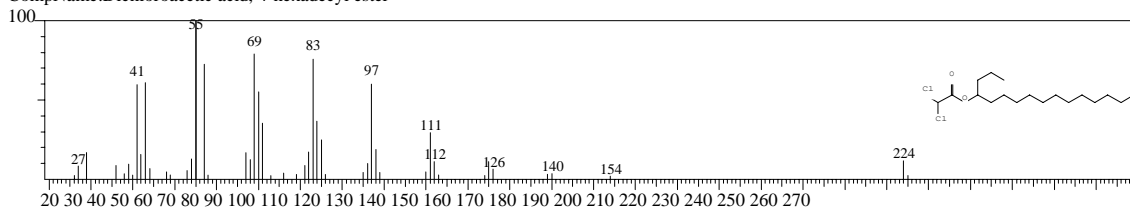
Hit#:3 Entry:17440 Library:NIST05s.LIB  
SI:90 Formula:C<sub>15</sub>H<sub>30</sub> CAS:13360-61-7 MolWeight:210 RetIndex:1502  
CompName:1-Pentadecene, 1- \$\$ Pentadec-1-ene \$\$



Hit#:4 Entry:72607 Library:NIST05.LIB  
SI:90 Formula:C<sub>18</sub>H<sub>36</sub> CAS:7206-25-9 MolWeight:252 RetIndex:1818  
CompName:9-Octadecene, (E)- \$\$ (9E)-9-Octadecene # \$\$



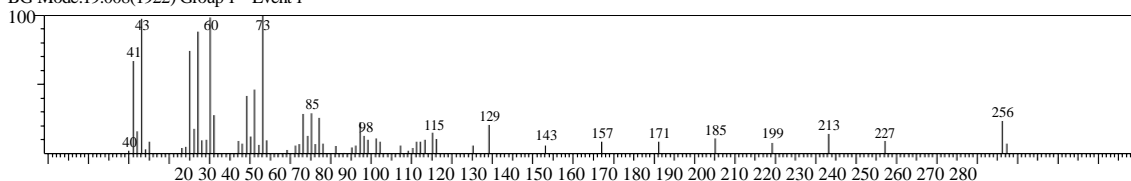
Hit#:5 Entry:127954 Library:NIST05.LIB  
SI:90 Formula:C<sub>18</sub>H<sub>34</sub>Cl<sub>2</sub>O<sub>2</sub> CAS:0-00-0 MolWeight:352 RetIndex:2259  
CompName:Dichloroacetic acid, 4-hexadecyl ester





&lt;&lt; Target &gt;&gt;

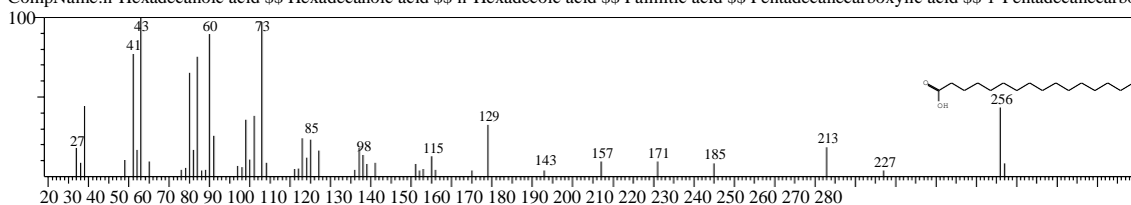
Line#6 R.Time:18.967(Scan#:1917) MassPeaks:62  
RawMode:Single 18.967(1917) BasePeak:73.05(36097)  
BG Mode:19.008(1922) Group 1 - Event 1



Hit#1 Entry:21331 Library:NIST05s.LIB

SI:93 Formula:C16H32O2 CAS:57-10-3 MolWeight:256 RetIndex:1968

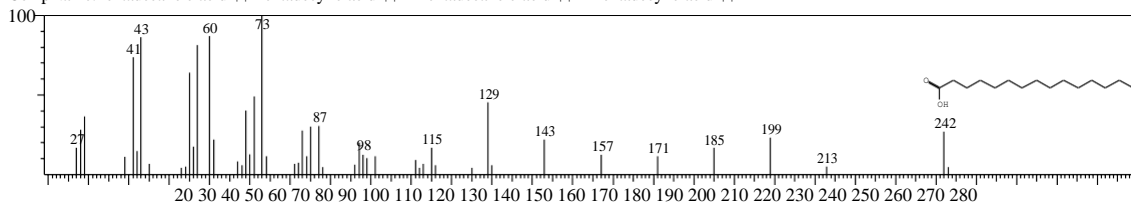
CompName:n-Hexadecanoic acid \$\$ Hexadecanoic acid \$\$ n-Hexadecoic acid \$\$ Palmitic acid \$\$ Pentadecanecarboxylic acid \$\$ 1-Pentadecanecarboxylic acid



Hit#2 Entry:20371 Library:NIST05s.LIB

SI:90 Formula:C15H30O2 CAS:1002-84-2 MolWeight:242 RetIndex:1869

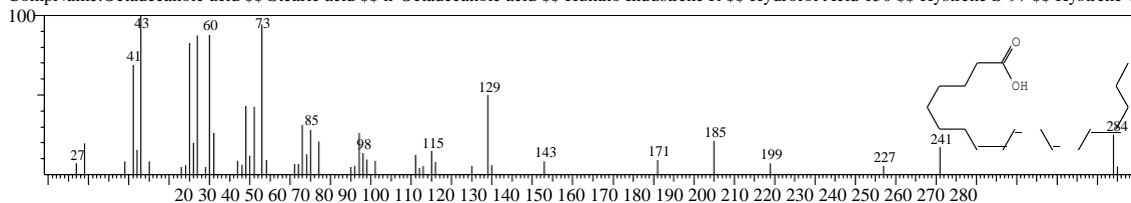
CompName:Pentadecanoic acid \$\$ Pentadecylic acid \$\$ n-Pentadecanoic acid \$\$ n-Pentadecylic acid \$\$



Hit#3 Entry:22979 Library:NIST05s.LIB

SI:90 Formula:C18H36O2 CAS:57-11-4 MolWeight:284 RetIndex:2167

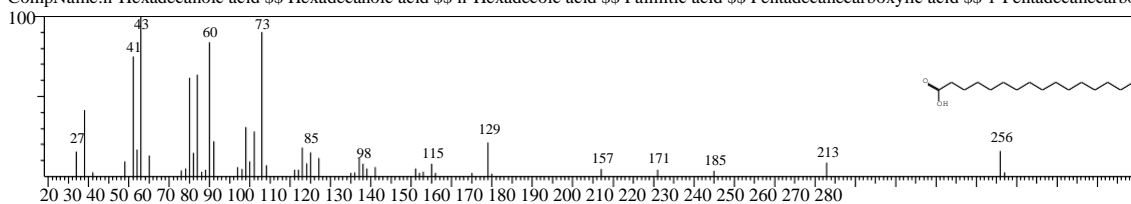
CompName:Octadecanoic acid \$\$ Stearic acid \$\$ n-Octadecanoic acid \$\$ Humko Industrine R \$\$ Hydrofol Acid 150 \$\$ Hystrene S-97 \$\$ Hystrene T-70 \$\$ Hys



Hit#4 Entry:74999 Library:NIST05.LIB

SI:89 Formula:C16H32O2 CAS:57-10-3 MolWeight:256 RetIndex:1968

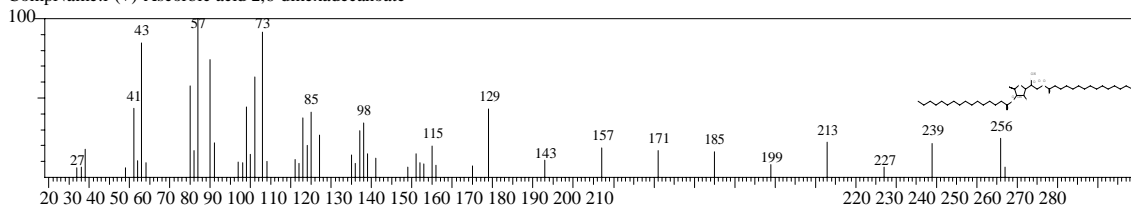
CompName:n-Hexadecanoic acid \$\$ Hexadecanoic acid \$\$ n-Hexadecoic acid \$\$ Palmitic acid \$\$ Pentadecanecarboxylic acid \$\$ 1-Pentadecanecarboxylic acid



Hit#5 Entry:161860 Library:NIST05.LIB

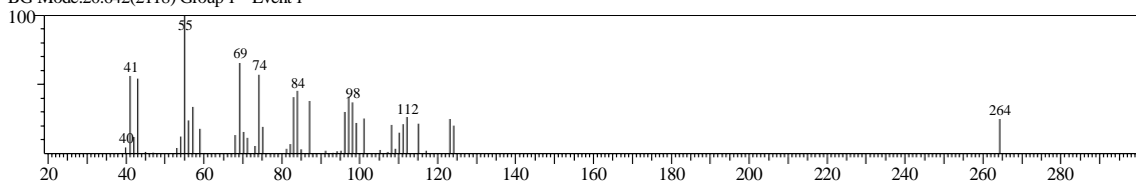
SI:89 Formula:C38H68O8 CAS:28474-90-0 MolWeight:652 RetIndex:4765

CompName:l-(+)-Ascorbic acid 2,6-dihexadecanoate



&lt;&lt; Target &gt;&gt;

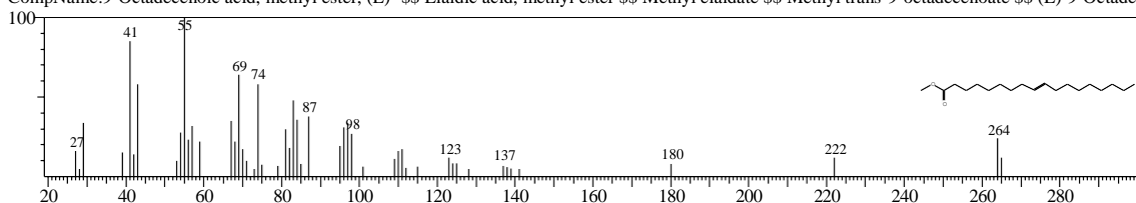
Line#:7 R.Time:20.667(Scan#:2121) MassPeaks:48  
RawMode:Single 20.667(2121) BasePeak:55.05(10434)  
BG Mode:20.642(2118) Group 1 - Event 1



Hit#:1 Entry:23570 Library:NIST05s.LIB

SI:86 Formula:C19H36O2 CAS:1937-62-8 MolWeight:296 RetIndex:2085

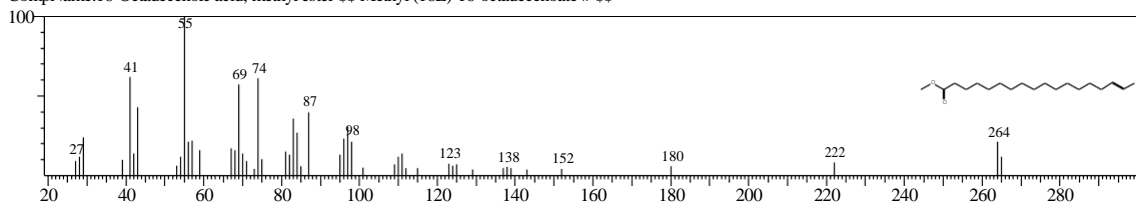
CompName:9-Octadecenoic acid, methyl ester, (E)- \$\$ Elaidic acid, methyl ester \$\$ Methyl elaidate \$\$ Methyl trans-9-octadecenoate \$\$ (E)-9-Octadecenoic aci



Hit#:2 Entry:98790 Library:NIST05.LIB

SI:86 Formula:C19H36O2 CAS:56554-49-5 MolWeight:296 RetIndex:2085

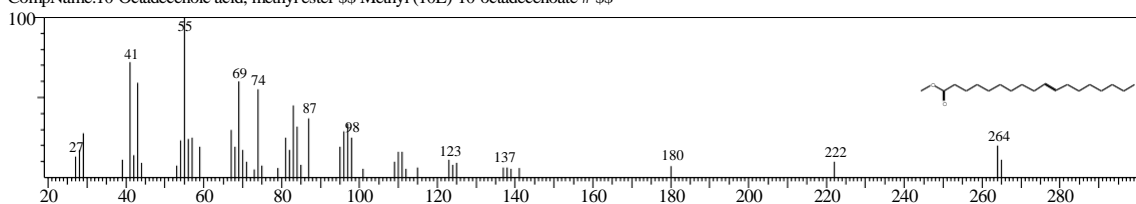
CompName:16-Octadecenoic acid, methyl ester \$\$ Methyl (16E)-16-octadecenoate # \$\$



Hit#:3 Entry:98784 Library:NIST05.LIB

SI:86 Formula:C19H36O2 CAS:13481-95-3 MolWeight:296 RetIndex:2085

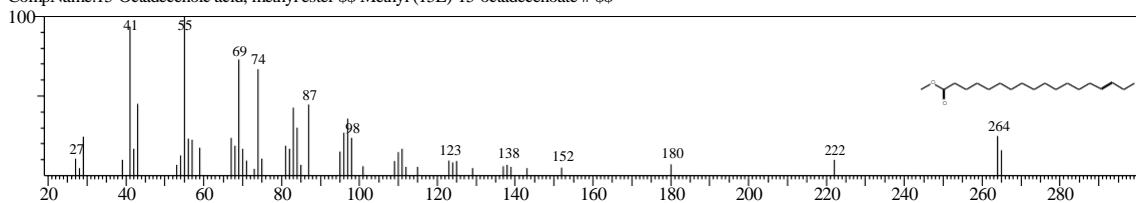
CompName:10-Octadecenoic acid, methyl ester \$\$ Methyl (10E)-10-octadecenoate # \$\$



Hit#:4 Entry:98781 Library:NIST05.LIB

SI:85 Formula:C19H36O2 CAS:4764-72-1 MolWeight:296 RetIndex:2085

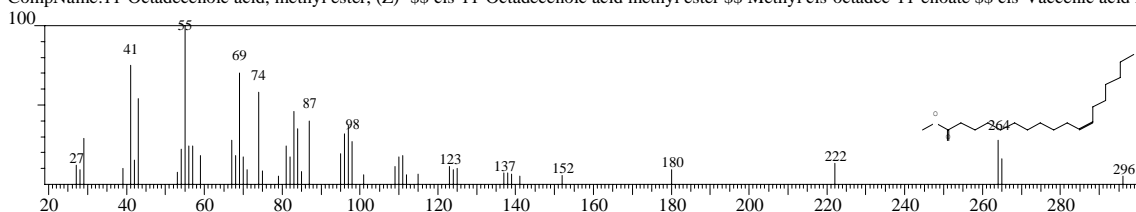
CompName:15-Octadecenoic acid, methyl ester \$\$ Methyl (15E)-15-octadecenoate # \$\$



Hit#:5 Entry:98785 Library:NIST05.LIB

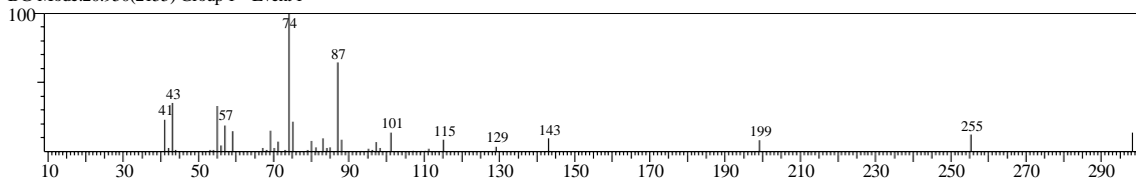
SI:85 Formula:C19H36O2 CAS:1937-63-9 MolWeight:296 RetIndex:2085

CompName:11-Octadecenoic acid, methyl ester, (Z)- \$\$ cis-11-Octadecenoic acid methyl ester \$\$ Methyl cis-octadec-11-enoate \$\$ cis-Vaccenic acid methyl est



&lt;&lt; Target &gt;&gt;

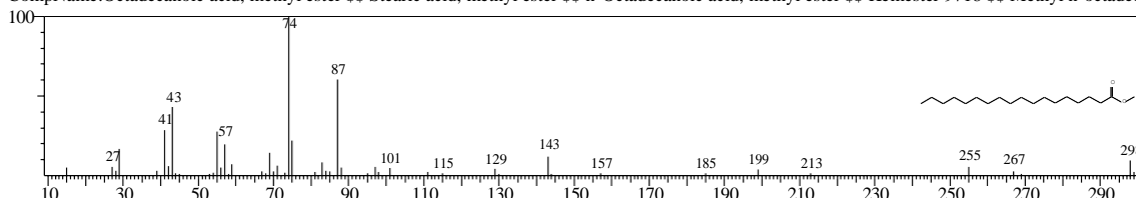
Line# 8 R.Time: 20.925 (Scan#: 2152) MassPeaks: 44  
RawMode: Single 20.925 (2152) BasePeak: 74.10 (27083)  
BG Mode: 20.950 (2155) Group 1 - Event 1



Hit#: 1 Entry: 100069 Library: NIST05.LIB

SI: 92 Formula: C<sub>19</sub>H<sub>38</sub>O<sub>2</sub> CAS: 112-61-8 MolWeight: 298 RetIndex: 2077

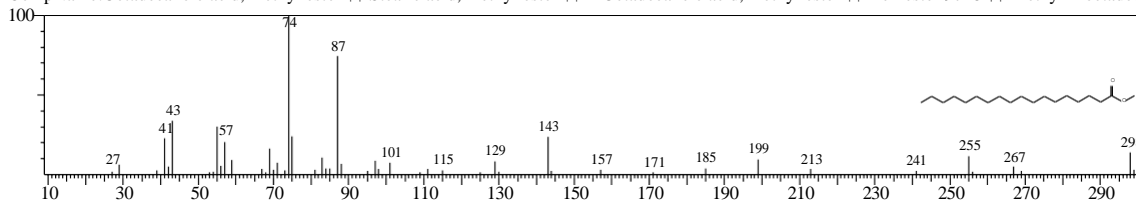
CompName: Octadecanoic acid, methyl ester \$\$\$\$ Stearic acid, methyl ester \$\$\$\$ n-Octadecanoic acid, methyl ester \$\$\$\$ Kemester 9718 \$\$\$\$ Methyl n-octadecanoate \$\$\$\$



Hit#: 2 Entry: 23659 Library: NIST05s.LIB

SI: 91 Formula: C<sub>19</sub>H<sub>38</sub>O<sub>2</sub> CAS: 112-61-8 MolWeight: 298 RetIndex: 2077

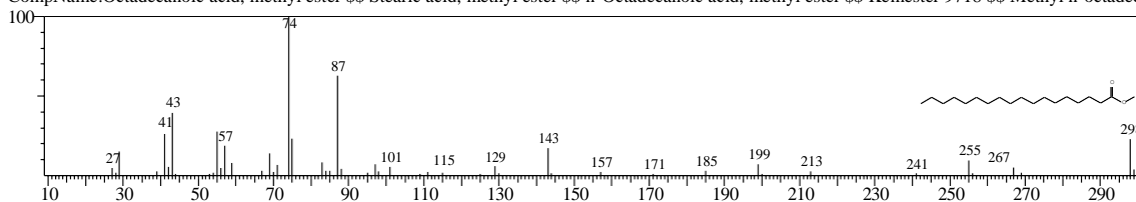
CompName: Octadecanoic acid, methyl ester \$\$\$\$ Stearic acid, methyl ester \$\$\$\$ n-Octadecanoic acid, methyl ester \$\$\$\$ Kemester 9718 \$\$\$\$ Methyl n-octadecanoate \$\$\$\$



Hit#: 3 Entry: 23658 Library: NIST05s.LIB

SI: 91 Formula: C<sub>19</sub>H<sub>38</sub>O<sub>2</sub> CAS: 112-61-8 MolWeight: 298 RetIndex: 2077

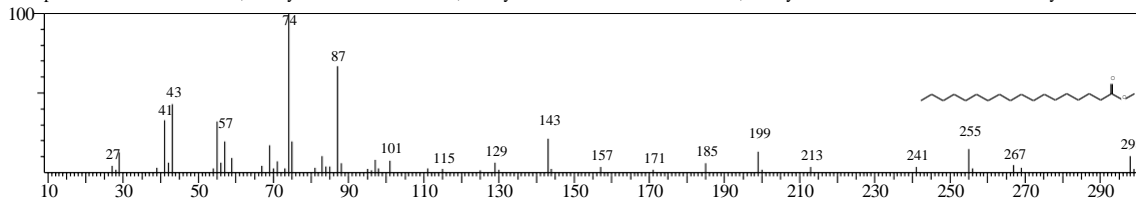
CompName: Octadecanoic acid, methyl ester \$\$\$\$ Stearic acid, methyl ester \$\$\$\$ n-Octadecanoic acid, methyl ester \$\$\$\$ Kemester 9718 \$\$\$\$ Methyl n-octadecanoate \$\$\$\$



Hit#: 4 Entry: 23657 Library: NIST05s.LIB

SI: 90 Formula: C<sub>19</sub>H<sub>38</sub>O<sub>2</sub> CAS: 112-61-8 MolWeight: 298 RetIndex: 2077

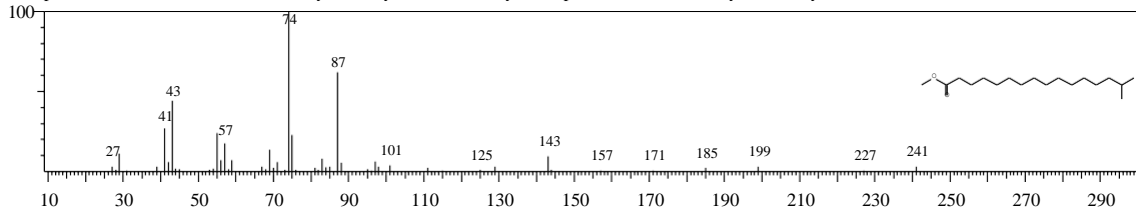
CompName: Octadecanoic acid, methyl ester \$\$\$\$ Stearic acid, methyl ester \$\$\$\$ n-Octadecanoic acid, methyl ester \$\$\$\$ Kemester 9718 \$\$\$\$ Methyl n-octadecanoate \$\$\$\$



Hit#: 5 Entry: 22987 Library: NIST05s.LIB

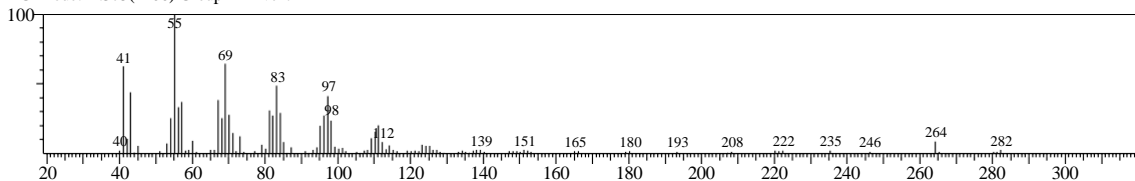
SI: 90 Formula: C<sub>18</sub>H<sub>36</sub>O<sub>2</sub> CAS: 6929-04-0 MolWeight: 284 RetIndex: 1914

CompName: Hexadecanoic acid, 15-methyl-, methyl ester \$\$\$\$ Methyl isoheptadecanoate \$\$\$\$ Methyl 15-methylhexadecanoate \$\$\$\$



&lt;&lt; Target &gt;&gt;

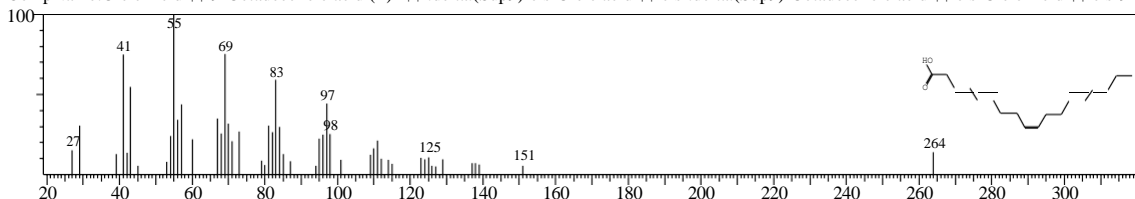
Line#9 R.Time:21.308(Scan#:2198) MassPeaks:105  
RawMode:Single 21.308(2198) BasePeak:55.05(152013)  
BG Mode:21.375(2206) Group 1 - Event 1



Hit#1 Entry:90577 Library:NIST05.LIB

SI:93 Formula:C18H34O2 CAS:112-80-1 MolWeight:282 RetIndex:2175

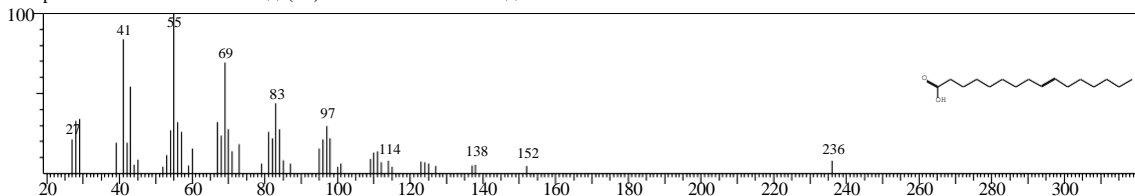
CompName:Oleic Acid \$ 9-Octadecenoic acid (Z)- \$\$.delta.(Sup9)-cis-Oleic acid \$ cis-.delta.(Sup9)-Octadecenoic acid \$ cis-Oleic Acid \$ cis-9-Octadecen



Hit#2 Entry:73685 Library:NIST05.LIB

SI:92 Formula:C16H30O2 CAS:2091-29-4 MolWeight:254 RetIndex:1976

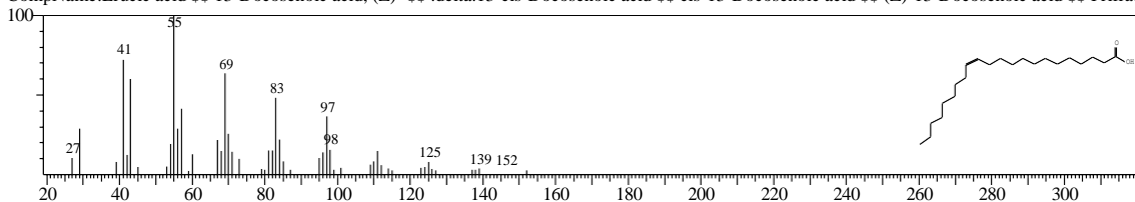
CompName:9-Hexadecenoic acid \$ (9E)-9-Hexadecenoic acid # \$



Hit#3 Entry:121691 Library:NIST05.LIB

SI:91 Formula:C22H42O2 CAS:112-86-7 MolWeight:338 RetIndex:2572

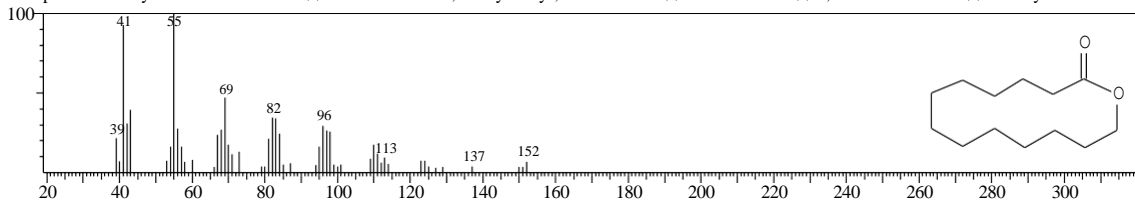
CompName:Erucic acid \$ 13-Docosenoic acid, (Z)- \$\$.delta.13-cis-Docosenoic acid \$ cis-13-Docosenoic acid \$ Prifrac 2990 \$



Hit#4 Entry:17628 Library:NIST05s.LIB

SI:90 Formula:C13H24O2 CAS:1725-04-8 MolWeight:212 RetIndex:1904

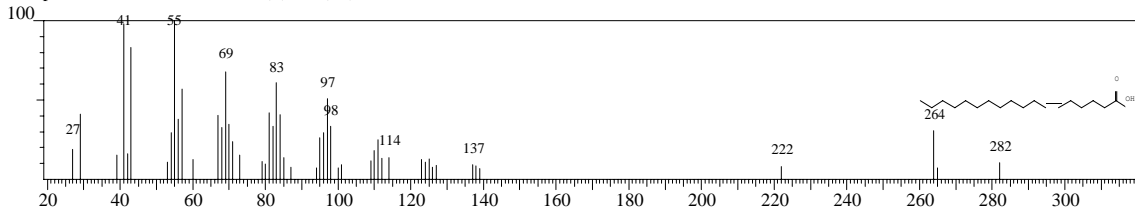
CompName:Oxacyclotetradecan-2-one \$ Tridecanoic acid, 13-hydroxy-, .mu.-lactone \$ Tridecanolide \$ 1,13-Tridecanolide \$ -Oxacyclotetradecan-2-one \$



Hit#5 Entry:90568 Library:NIST05.LIB

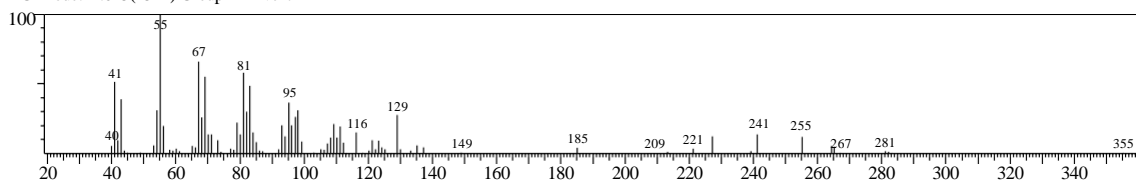
SI:90 Formula:C18H34O2 CAS:593-39-5 MolWeight:282 RetIndex:2175

CompName:6-Octadecenoic acid, (Z)- \$ (6Z)-6-Octadecenoic acid # \$



&lt;&lt; Target &gt;&gt;

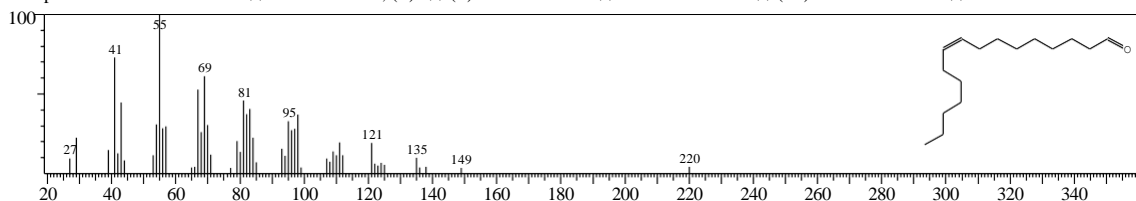
Line#:10 R.Time:23.983(Scan#:2519) MassPeaks:89  
RawMode:Single 23.983(2519) BasePeak:55.05(16852)  
BG Mode:24.025(2524) Group 1 - Event 1



Hit#:1 Entry:20037 Library:NIST05s.LIB

SI:86 Formula:C16H30O CAS:56219-04-6 MolWeight:238 RetIndex:1808

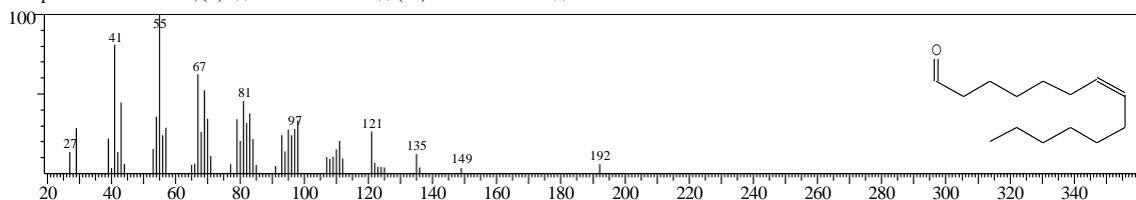
CompName:cis-9-Hexadecenal \$\$ 9-Hexadecenal, (Z)- \$\$ (Z)-9-Hexadecenal \$\$ Z-9-Hexadecenal \$\$ (9Z)-9-Hexadecenal # \$\$



Hit#:2 Entry:47275 Library:NIST05s.LIB

SI:86 Formula:C14H26O CAS:65128-96-3 MolWeight:210 RetIndex:1609

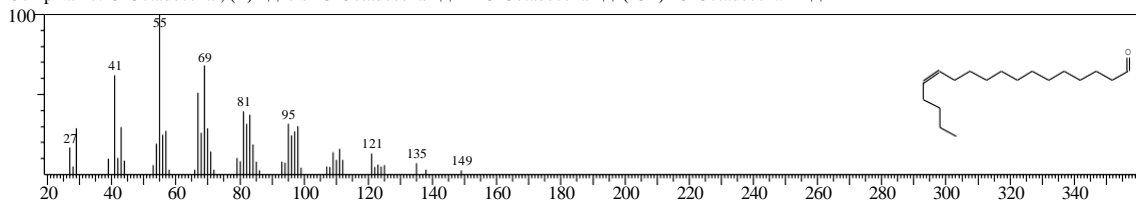
CompName:7-Tetradecenal, (Z)- \$\$ Z-7-Tetradecenal \$\$ (7Z)-7-Tetradecenal # \$\$



Hit#:3 Entry:21955 Library:NIST05s.LIB

SI:85 Formula:C18H34O CAS:58594-45-9 MolWeight:266 RetIndex:2007

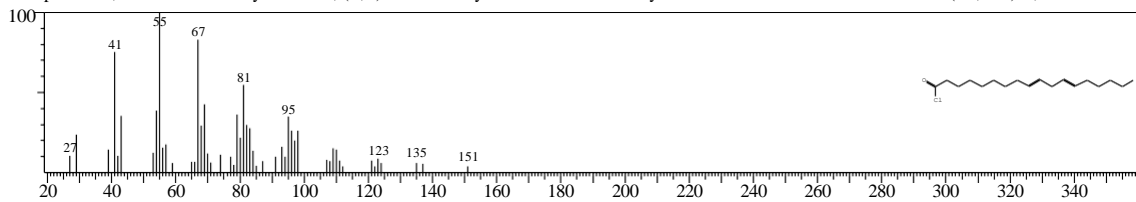
CompName:13-Octadecenal, (Z)- \$\$ cis-13-Octadecenal \$\$ Z-13-Octadecenal \$\$ (13Z)-13-Octadecenal # \$\$



Hit#:4 Entry:23649 Library:NIST05s.LIB

SI:85 Formula:C18H31ClO CAS:7459-33-8 MolWeight:298 RetIndex:2139

CompName:9,12-Octadecadienoyl chloride, (Z,Z)- \$\$ Linoleoyl chloride \$\$ Lineoleoyl chloride \$\$ Linoleic acid chloride \$\$ (9E,12E)-9,12-Octadecadienoyl ch



Hit#:5 Entry:47277 Library:NIST05s.LIB

SI:85 Formula:C14H26O CAS:85896-31-7 MolWeight:210 RetIndex:1591

CompName:13-Tetradecenal

