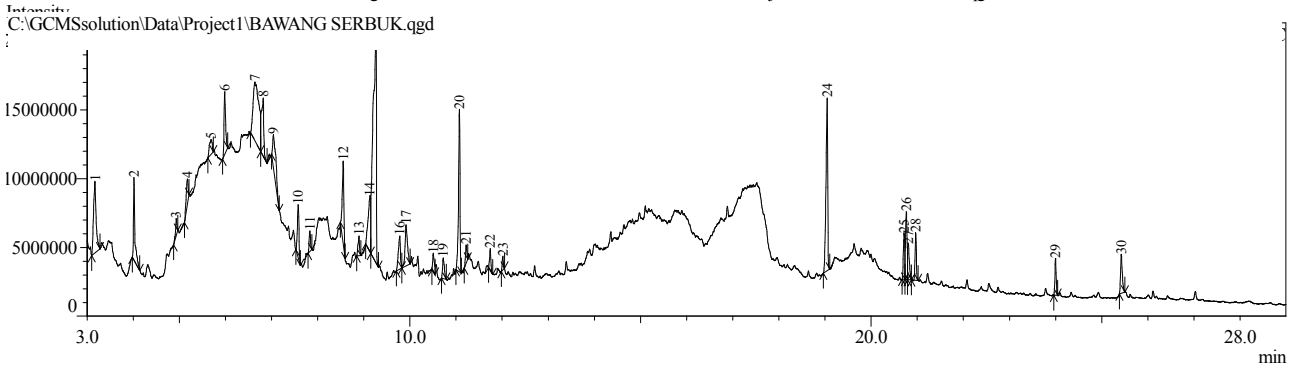
Sample Information

Analyzed by	: Admin
Analyzed	: 10/12/2019 1:19:49 PM
Sample Type	: Unknown
Level #	: 1
Sample Name	: BAWANG SERBUK
Sample ID	: 02268
IS Amount	: [1]=1
Sample Amount	: 1
Dilution Factor	: 1
Vial #	: 16
Injection Volume	: 1.00
Data File	: C:\GCMSsolution\Data\Project1\BAWANG SERBUK.qgd
Org Data File	: C:\GCMSsolution\Data\Project1\BAWANG SERBUK.qgd
Method File	: C:\GCMSsolution\Data\Project1\FITOKIMIA BARU.qgm
Org Method File	: C:\GCMSsolution\Data\Project1\FITOKIMIA BARU.qgm
Report File	:
Tuning File	: C:\GCMSsolution\System\Tune1\10042014.qgt
Modified by	: Admin
Modified	: 10/12/2019 1:48:49 PM

Chromatogram BAWANG SERBUK C:\GCMSsolution\Data\Project1\BAWANG SERBUK.qgd

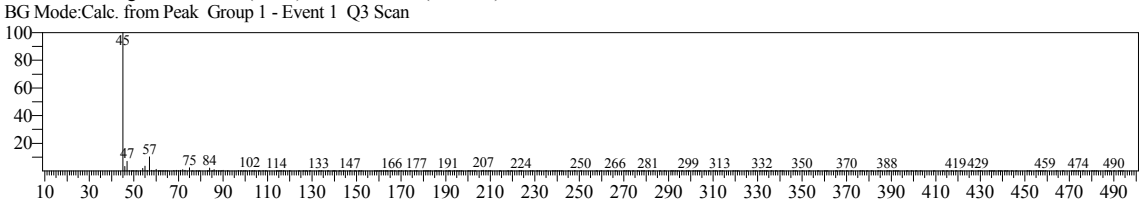


Peak Report TIC

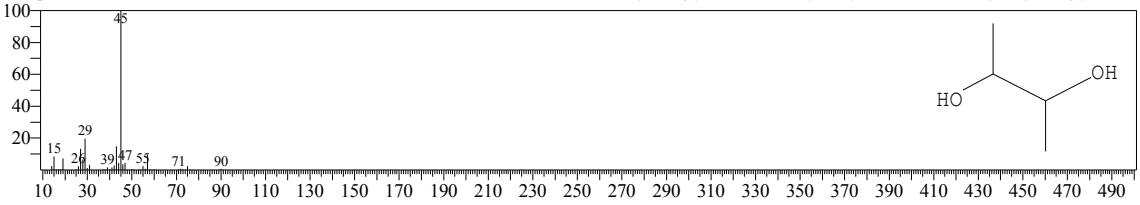
Peak#	R.Time	Area	Area%	Height	Name
1	3.171	20944259	5.40	5202288	2,3-Butanediol (CAS)
2	4.019	13003532	3.35	6039241	2-Furanmethanol (CAS)
3	4.934	2693511	0.69	1271852	.ALPHA.,.BETA.-CROTONOLACTONE
4	5.176	5772556	1.49	1850452	1,2-Cyclopentanedione
5	5.692	4634732	1.19	1065487	5 METHYL FURFURAL
6	5.985	10539454	2.72	4569134	2,3-Dihydro-3,5-dihydroxy-6-methyl-4H-pyran-4-one
7	6.640	38699532	9.97	4263904	Methane, sulfinylbis- (CAS)
8	6.820	14229340	3.67	4234049	Methane, sulfinylbis- (CAS)
9	7.041	12154357	3.13	2398972	2-Octanol, (S)-
10	7.579	8974521	2.31	3854511	Diallyl disulphide
11	7.831	3107912	0.80	1504202	HYDROXY DIMETHYL FURANONE
12	8.553	14529010	3.74	6005200	Hydroperoxide, 1,4-dioxan-2-yl (CAS)
13	8.897	3330373	0.86	1415064	1-Cyclohexene-1-carboxylic acid
14	9.130	16462547	4.24	3957332	Ethanamine, N-ethyl-N-nitroso- (CAS)
15	9.265	82747207	21.33	17427884	2,3-Dihydro-3,5-dihydroxy-6-methyl-4H-pyran-4-one
16	9.780	7587925	1.96	2420154	4H-Pyran-4-one, 3,5-dihydroxy-2-methyl- (CAS)
17	9.915	11141284	2.87	2990369	Benzene, 1,3-bis(tetrahydropyran-2-yloxy)-
18	10.507	3417129	0.88	1324080	2-(1-Methylpropyl)pyrazine
19	10.722	4030532	1.04	1481847	1,2-Benzenediol, 3-methyl-
20	11.074	26160847	6.74	11810673	Trisulfide, di-2-propenyl
21	11.221	2895466	0.75	1327016	Cyclotetrasiloxane, octamethyl- (CAS)
22	11.742	2951441	0.76	1552151	Benzenepropanoic acid, silver(1+) salt (CAS)
23	12.014	1741480	0.45	990682	5-methyl-1,2,3,4-tetrathia-cyclohexane
24	19.047	32133066	8.28	12554889	n-Hexadecanoic acid
25	20.716	7399037	1.91	3337568	9,12-Octadecadienoic acid (Z,Z)-
26	20.762	11271888	2.91	4936290	OCTADEC-9-ENOIC ACID
27	20.810	6288128	1.62	2540047	OCTADEC-9-ENOIC ACID
28	20.967	6594647	1.70	3479589	Octadecanoic acid
29	23.998	5041075	1.30	2678782	Hexadecanoic acid, 2-hydroxy-1-(hydroxymethyl)ethyl ester (CAS)
30	25.424	7495218	1.93	2842080	9-Octadecenal, (Z)- (CAS)
		387972006	100.00	121325789	

Library

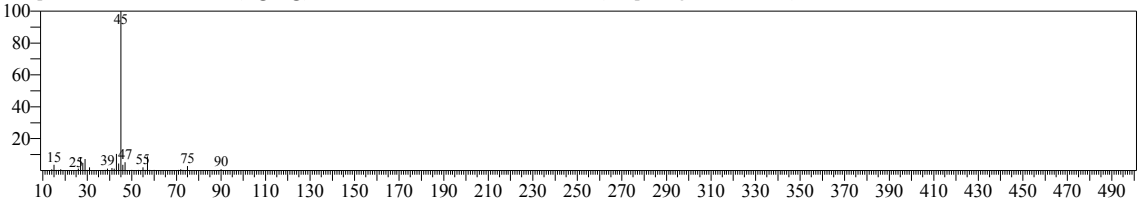
<< Target >>
Line#:1 R.Time:3.170(Scan#:35) MassPeaks:242
RawMode:Averaged 3.165-3.175(34-36) BasePeak:45.00(3748269)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



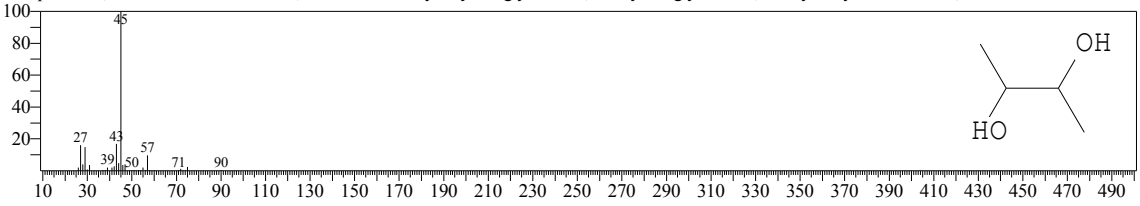
Hit#:1 Entry:5189 Library:Wiley9.lib
SI:96 Formula:C4H10O2 CAS:513-85-9 MolWeight:90 RetIndex:0
CompName:2,3-Butanediol (CAS) $$ Butane-2,3-diol $$ 2,3-BUTANDIOL $$ 2,3-Butylene glycol $$ 2,3-Dihydroxybutane $$ Dimethylethylene glycol $$



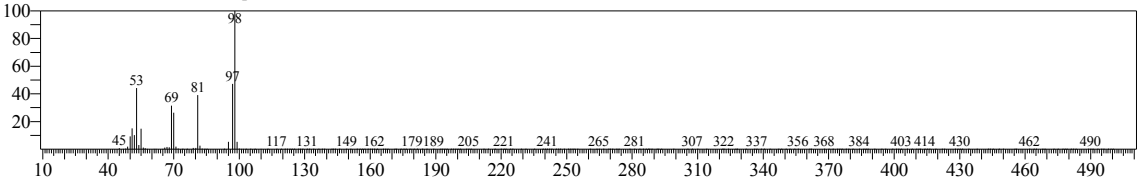
Hit#:2 Entry:5234 Library:Wiley9.lib
SI:96 Formula:C4H10O2 CAS:24347-58-8 MolWeight:90 RetIndex:0
CompName:2,3-Butanediol, [R-(R@,R@)]- $$ 2,3-Butanediol $$ 2,3-Butanediol (computer-generated name)



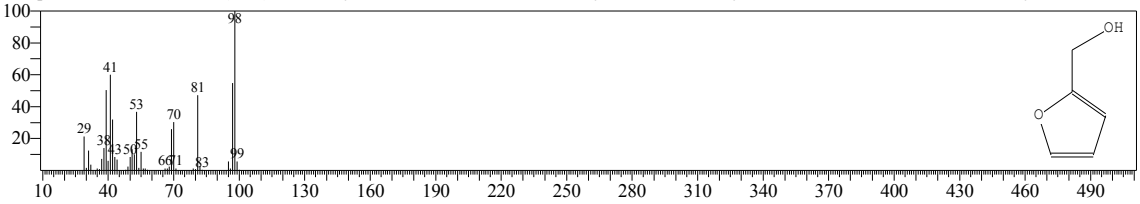
Hit#:3 Entry:1165 Library:NIST08s.LIB
SI:96 Formula:C4H10O2 CAS:513-85-9 MolWeight:90 RetIndex:743
CompName:2,3-Butanediol $$ Butane-2,3-diol $$ Dimethylethylene glycol $$ 2,3-Butylene glycol $$ 2,3-Dihydroxybutane $$ D-2,3-Butane diol $$



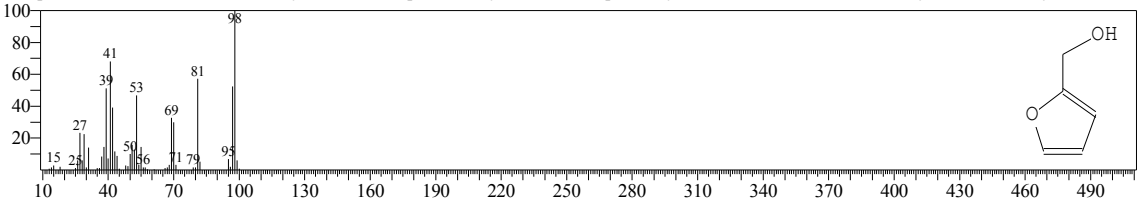
<< Target >>
Line#:2 R.Time:4.020(Scan#:205) MassPeaks:238
RawMode:Averaged 4.015-4.025(204-206) BasePeak:98.00(1517236)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



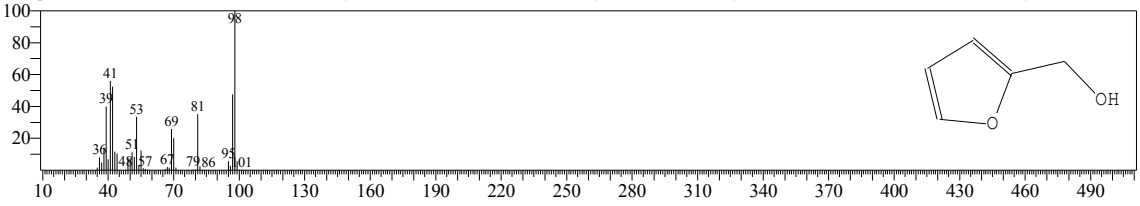
Hit#:1 Entry:6711 Library:Wiley9.lib
SI:97 Formula:C5H6O2 CAS:98-00-0 MolWeight:98 RetIndex:0
CompName:2-Furanmethanol (CAS) $$ Furfuryl alcohol $$ Furfuralcohol $$ Furyl alcohol $$ 2-Furylmethanol $$ 2-Furancarbinol $$ 2-Furylcarbinol $$ 2-



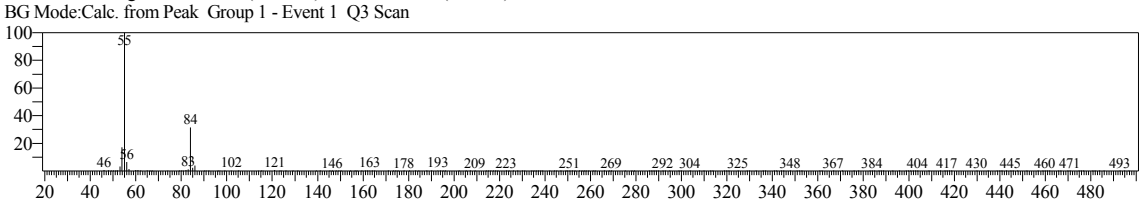
Hit#:2 Entry:1482 Library:NIST08s.LIB
SI:96 Formula:C5H6O2 CAS:98-00-0 MolWeight:98 RetIndex:885
CompName:2-Furanmethanol $$ Furfuryl alcohol $$ .alpha.-Furfuryl alcohol $$ .alpha.-Furylcarbinol $$ Furfuralcohol $$ Furyl alcohol $$ Furylcarbinol $$



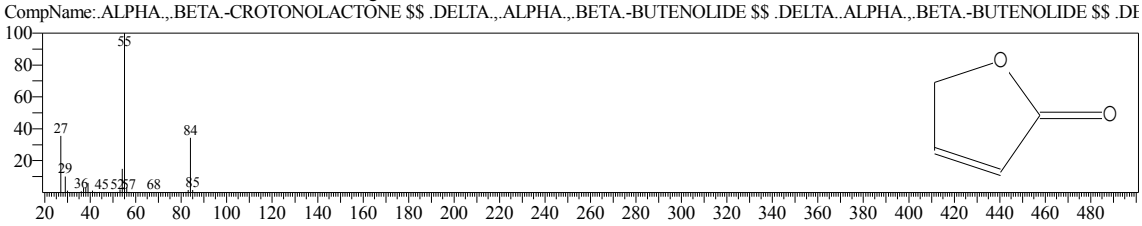
Hit#:3 Entry:6703 Library:Wiley9.lib
SI:96 Formula:C5H6O2 CAS:98-00-0 MolWeight:98 RetIndex:0
CompName:2-Furanmethanol (CAS) $$ Furfuryl alcohol $$ Furfuralcohol $$ Furyl alcohol $$ 2-Furylmethanol $$ 2-Furancarbinol $$ 2-Furylcarbinol $$ 2-



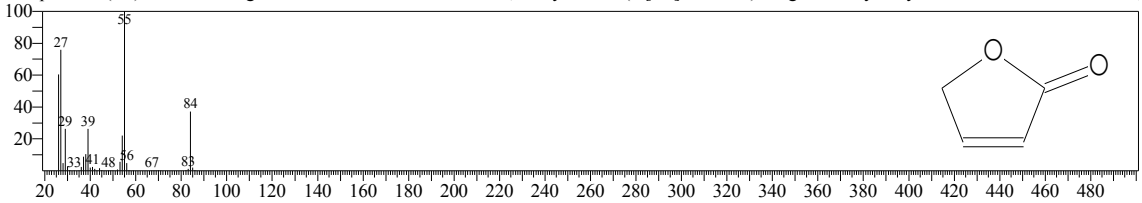
<< Target >>
Line#:3 R.Time:4.935(Scan#:388) MassPeaks:204
RawMode:Averaged 4.930-4.940(387-389) BasePeak:55.00(776391)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



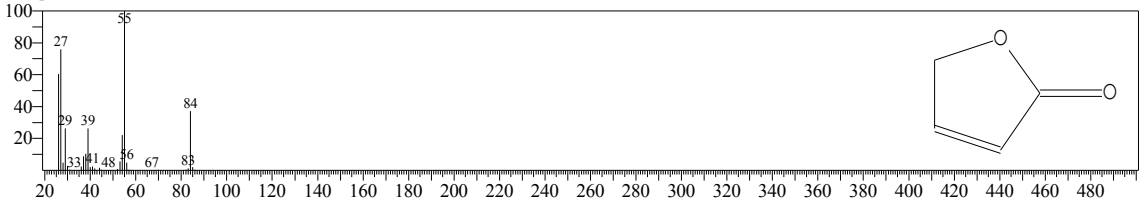
Hit#:1 Entry:3079 Library:Wiley9.lib
SI:96 Formula:C4H4O2 CAS:497-23-4 MolWeight:84 RetIndex:0
CompName:.ALPHA.,.BETA.-CROTONOLACTONE $$ .DELTA.,.ALPHA.,.BETA.-BUTENOLIDE $$ .DELTA..ALPHA.,.BETA.-BUTENOLIDE $$ .DE



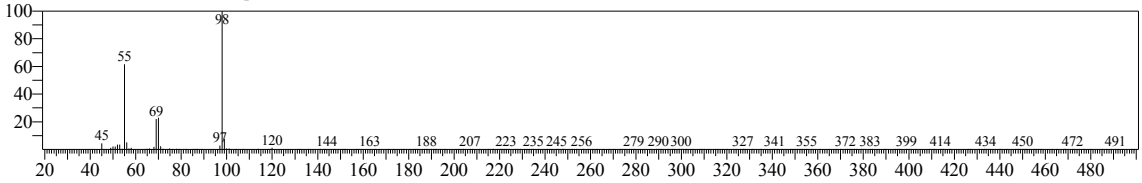
Hit#:2 Entry:687 Library:NIST08s.LIB
SI:95 Formula:C4H4O2 CAS:497-23-4 MolWeight:84 RetIndex:807
CompName:2(5H)-Furanone $$ .gamma.-Crotonolactone $$ 2-Oxo-2,5-dihydrofuran(2-[5H]-furanone) $$ .gamma.-Hydroxycrotonic acid lactone $$ .delta.



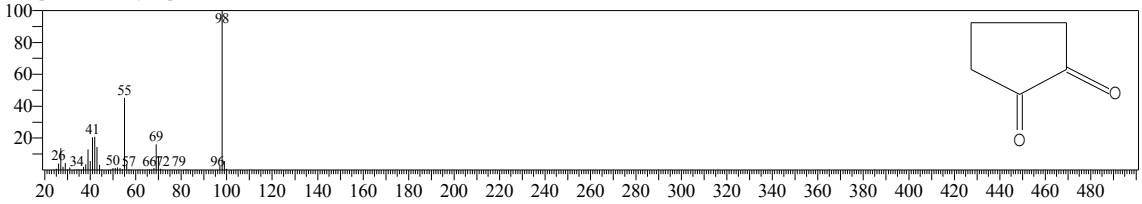
Hit#:3 Entry:3080 Library:Wiley9.lib
SI:95 Formula:C4H4O2 CAS:497-23-4 MolWeight:84 RetIndex:0
CompName:.ALPHA.,.BETA.-CROTONOLACTONE $$ .DELTA.,.ALPHA.,.BETA.-BUTENOLIDE $$ .DELTA..ALPHA.,.BETA.-BUTENOLIDE $$ .DE



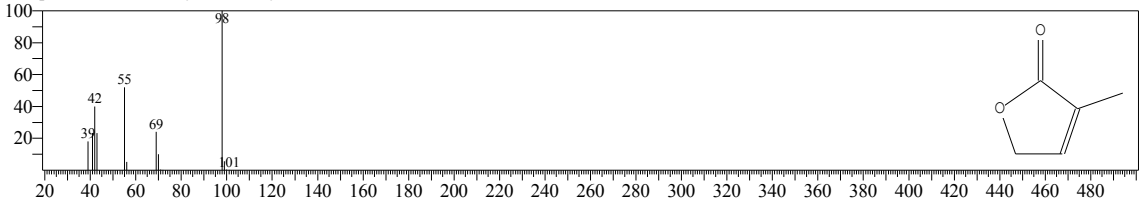
<< Target >>
Line#:4 R.Time:5.175(Scan#:436) MassPeaks:223
RawMode:Averaged 5.170-5.180(435-437) BasePeak:98.00(743158)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



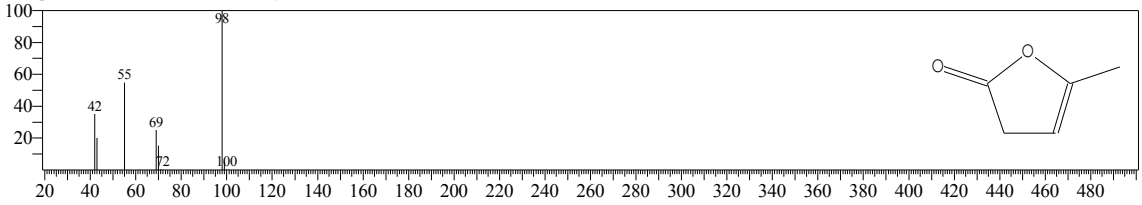
Hit#:1 Entry:6759 Library:Wiley9.lib
SI:92 Formula:C5H6O2 CAS:3008-40-0 MolWeight:98 RetIndex:0
CompName:1,2-Cyclopentanedione $$ 1,2-CYCLOOCTANEDIONE



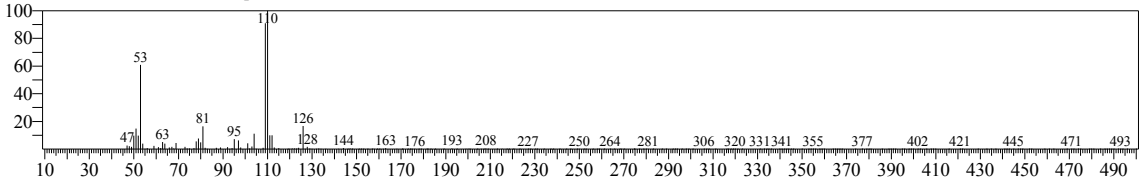
Hit#:2 Entry:6771 Library:Wiley9.lib
SI:90 Formula:C5H6O2 CAS:0-00-0 MolWeight:98 RetIndex:0
CompName:isomeric dihydro - methyl - furanone



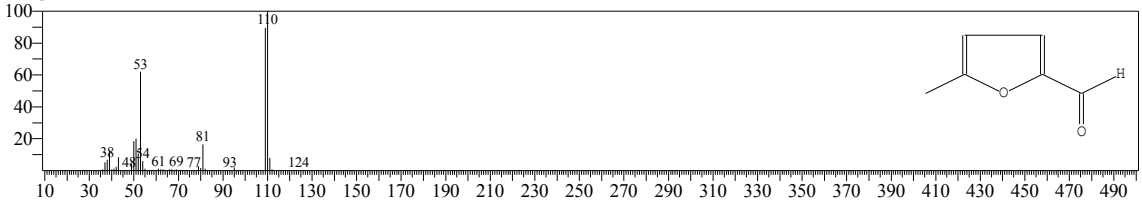
Hit#:3 Entry:6683 Library:Wiley9.lib
SI:89 Formula:C5H6O2 CAS:591-12-8 MolWeight:98 RetIndex:0
CompName:2(3H)-Furanone, 5-methyl- (CAS) $$ 5-METHYL-2-OXO-2,3-DIHYDROFURAN $$ 5-METHYL-3-HYDROFURAN-2-ONE $$ 3-Pentenoic



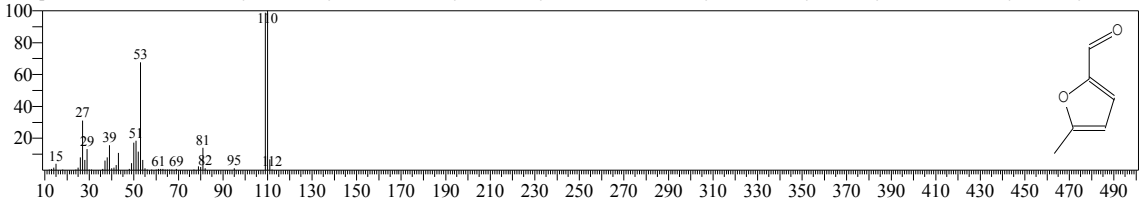
<< Target >>
Line#:5 R.Time:5.690(Scan#:539) MassPeaks:244
RawMode:Averaged 5.685-5.695(538-540) BasePeak:110.00(257062)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



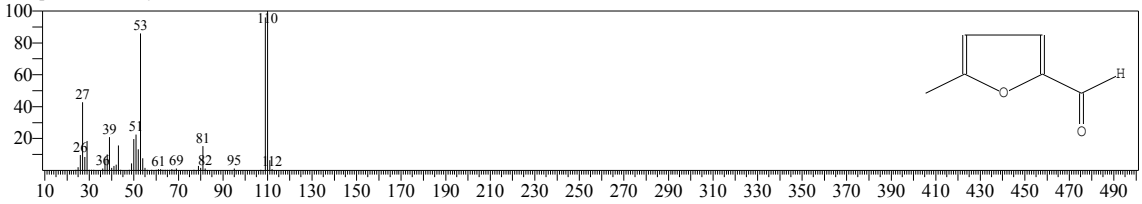
Hit#:1 Entry:12106 Library:Wiley9.lib
SI:89 Formula:C6H6O2 CAS:620-02-0 MolWeight:110 RetIndex:0
CompName:5 METHYL FURFURAL $$ 2-FURANCARBOXALDEHYDE, 5-METHYL- $$ 5-METHYLFURAN-2-CARBALDEHYDE $$ 2-FORMYL-



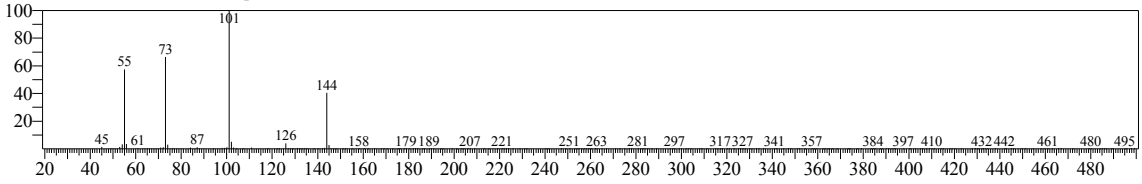
Hit#:2 Entry:2613 Library:NIST08s.LIB
SI:87 Formula:C6H6O2 CAS:620-02-0 MolWeight:110 RetIndex:920
CompName:2-Furancarboxaldehyde, 5-methyl- $$ 2-Furaldehyde, 5-methyl- $$ Furfural, 5-methyl- $$ 2-Formyl-5-methylfuran $$ 2-Methyl-5-formylfuran $



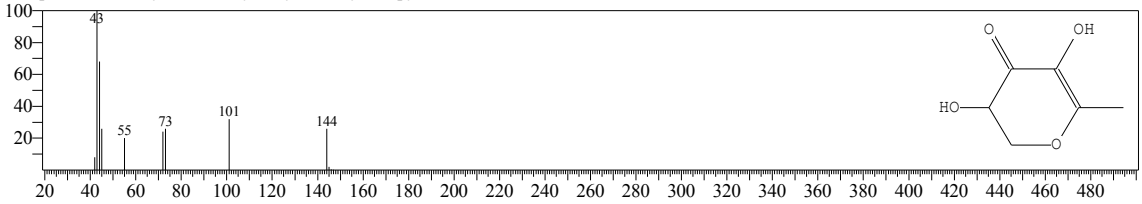
Hit#:3 Entry:12108 Library:Wiley9.lib
SI:86 Formula:C6H6O2 CAS:620-02-0 MolWeight:110 RetIndex:0
CompName:5-methyl furfural $$ 2-FURANCARBOXALDEHYDE, 5-METHYL- $$ 5-METHYLFURAN-2-CARBALDEHYDE $$ 2-FORMYL-5-METH



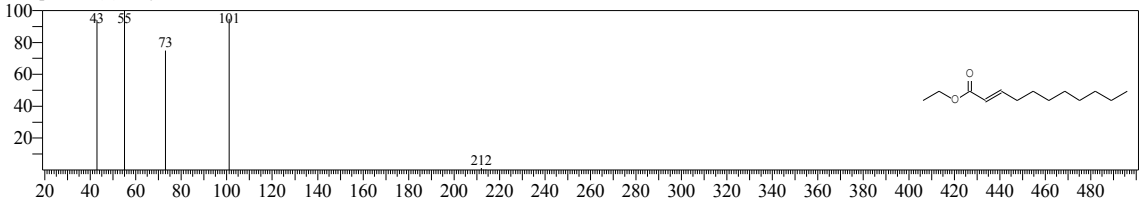
<< Target >>
Line#:6 R.Time:5.985(Scan#:598) MassPeaks:250
RawMode:Averaged 5.980-5.990(597-599) BasePeak:101.00(1468303)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



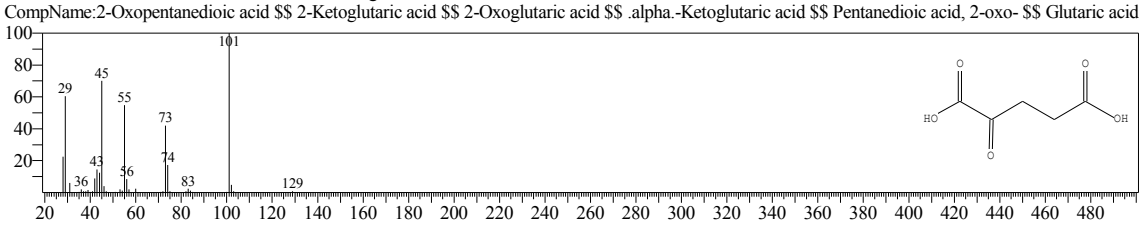
Hit#:1 Entry:42592 Library:Wiley9.lib
SI:82 Formula:C6H8O4 CAS:28564-83-2 MolWeight:144 RetIndex:0
CompName:2,3-Dihydro-3,5-dihydroxy-6-methyl-4H-pyran-4-one



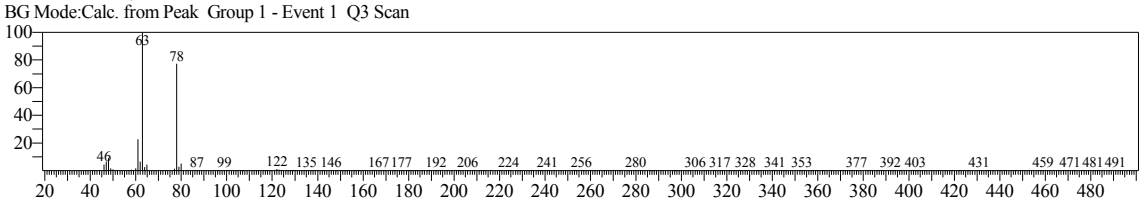
Hit#:2 Entry:170405 Library:Wiley9.lib
SI:81 Formula:C13H24O2 CAS:0-00-0 MolWeight:212 RetIndex:0
CompName:(E) Ethyl Undec-2-enoate



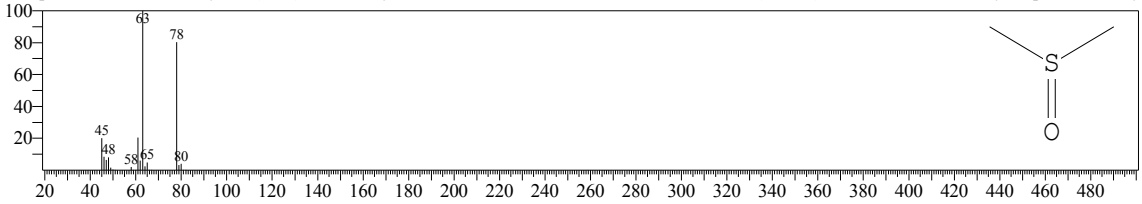
Hit#:3 Entry:44866 Library:Wiley9.lib
SI:80 Formula:C5H6O5 CAS:328-50-7 MolWeight:146 RetIndex:0
CompName:2-Oxopentanedioic acid $$ 2-Ketoglutaric acid $$ 2-Oxoglutaric acid $$ .alpha.-Ketoglutaric acid $$ Pentanedioic acid, 2-oxo- $$ Glutaric acid



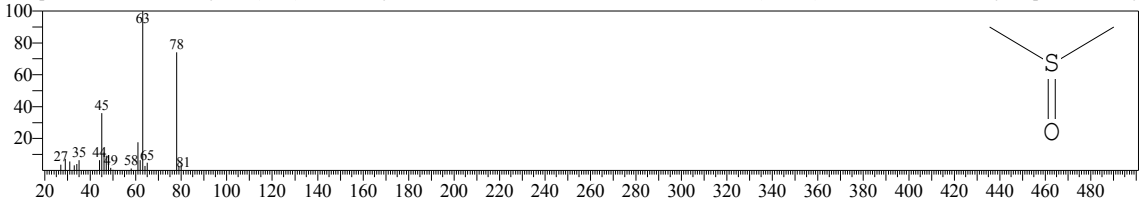
<< Target >>
Line#:7 R.Time:6.640(Scan#:729) MassPeaks:228
RawMode:Averaged 6.635-6.645(728-730) BasePeak:62.95(1315423)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



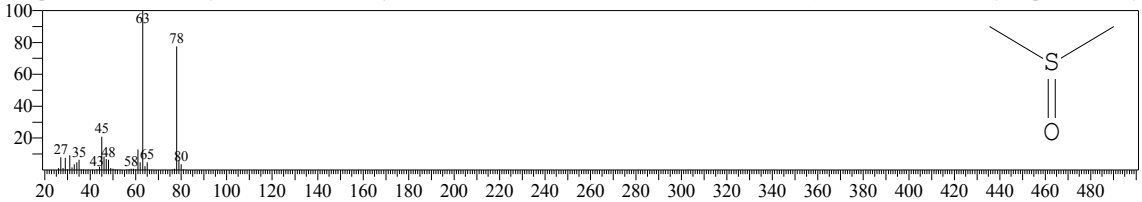
Hit#:1 Entry:2388 Library:Wiley9.lib
SI:97 Formula:C2H6OS CAS:67-68-5 MolWeight:78 RetIndex:0
CompName:Methane, sulfinylbis- (CAS) $$ Dimethyl sulfoxide $$ DMSO $$ DMS 70 $$ DMS 90 $$ (DMSO) $$ SQ 9453 $$ Dimethyl sulphoxide $$ Hy



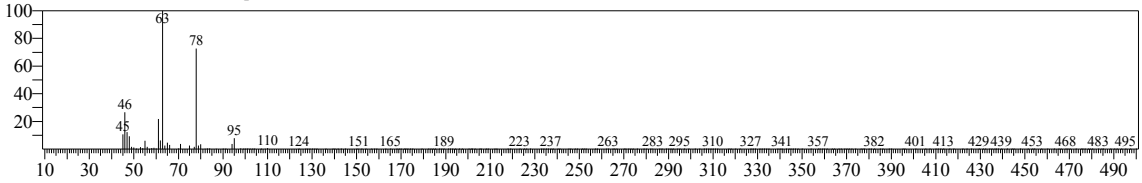
Hit#:2 Entry:2385 Library:Wiley9.lib
SI:96 Formula:C2H6OS CAS:67-68-5 MolWeight:78 RetIndex:0
CompName:Methane, sulfinylbis- (CAS) $$ Dimethyl sulfoxide $$ DMSO $$ DMS 70 $$ DMS 90 $$ (DMSO) $$ SQ 9453 $$ Dimethyl sulphoxide $$ Hy



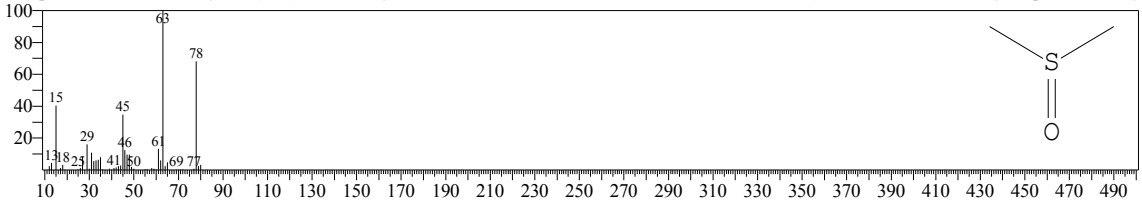
Hit#:3 Entry:2381 Library:Wiley9.lib
SI:96 Formula:C2H6OS CAS:67-68-5 MolWeight:78 RetIndex:0
CompName:Methane, sulfinylbis- (CAS) $$ Dimethyl sulfoxide $$ DMSO $$ DMS 70 $$ DMS 90 $$ (DMSO) $$ SQ 9453 $$ Dimethyl sulphoxide $$ Hy



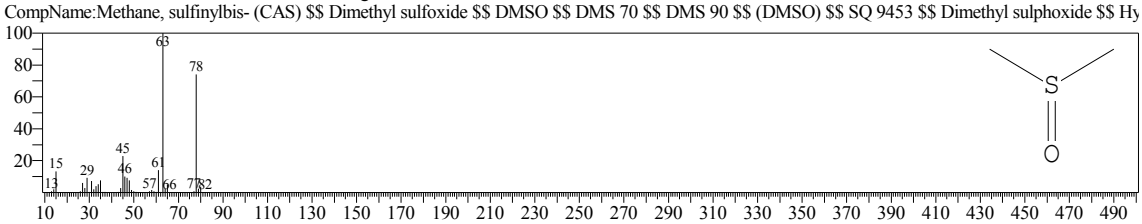
<< Target >>
Line#:8 R.Time:6.820(Scan#:765) MassPeaks:259
RawMode:Averaged 6.815-6.825(764-766) BasePeak:62.95(765483)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



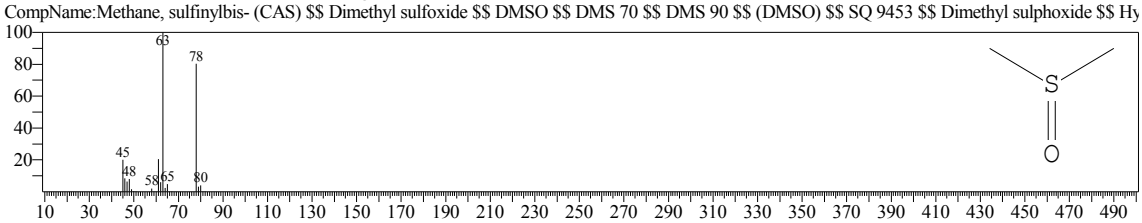
Hit#:1 Entry:2384 Library:Wiley9.lib
SI:90 Formula:C2H6OS CAS:67-68-5 MolWeight:78 RetIndex:0
CompName:Methane, sulfinylbis- (CAS) $$ Dimethyl sulfoxide $$ DMSO $$ DMS 70 $$ DMS 90 $$ (DMSO) $$ SQ 9453 $$ Dimethyl sulphoxide $$ Hy



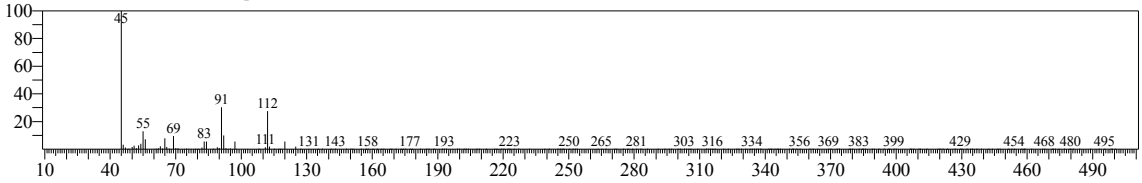
Hit#:2 Entry:2386 Library:Wiley9.lib
SI:90 Formula:C2H6OS CAS:67-68-5 MolWeight:78 RetIndex:0
CompName:Methane, sulfinylbis- (CAS) $$ Dimethyl sulfoxide $$ DMSO $$ DMS 70 $$ DMS 90 $$ (DMSO) $$ SQ 9453 $$ Dimethyl sulphoxide $$ Hy



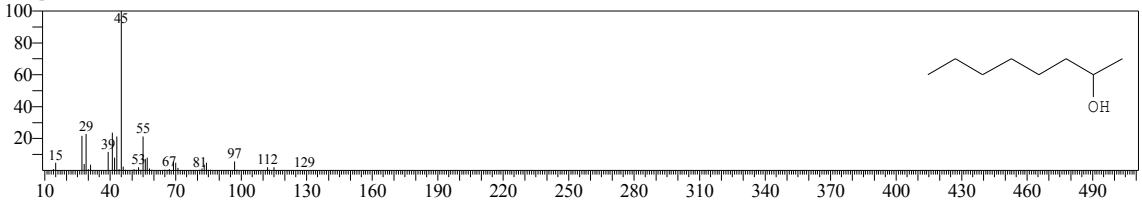
Hit#:3 Entry:2388 Library:Wiley9.lib
SI:90 Formula:C2H6OS CAS:67-68-5 MolWeight:78 RetIndex:0
CompName:Methane, sulfinylbis- (CAS) $$ Dimethyl sulfoxide $$ DMSO $$ DMS 70 $$ DMS 90 $$ (DMSO) $$ SQ 9453 $$ Dimethyl sulphoxide $$ Hy



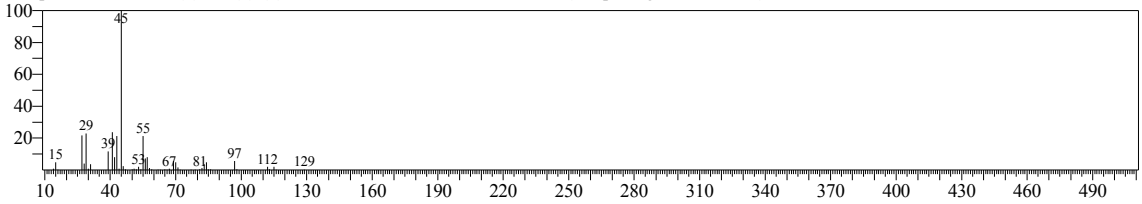
<< Target >>
Line#:9 R.Time:7.040(Scan#:809) MassPeaks:229
RawMode:Averaged 7.035-7.045(808-810) BasePeak:45.00(935522)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



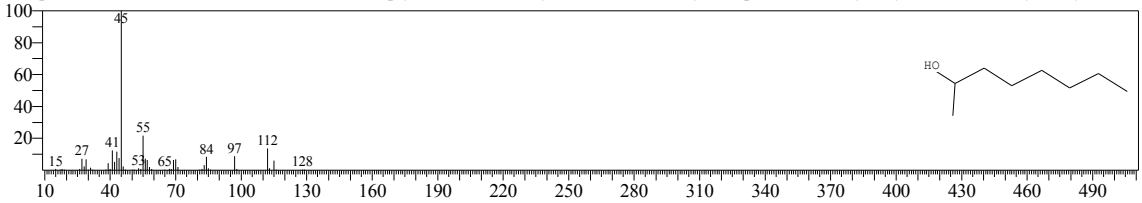
Hit#:1 Entry:5437 Library:NIST08s.LIB
SI:77 Formula:C8H18O CAS:6169-06-8 MolWeight:130 RetIndex:979
CompName:2-Octanol, (S)- $$ (S)-(+)-2-Octanol $$ 2-Octanol # $$



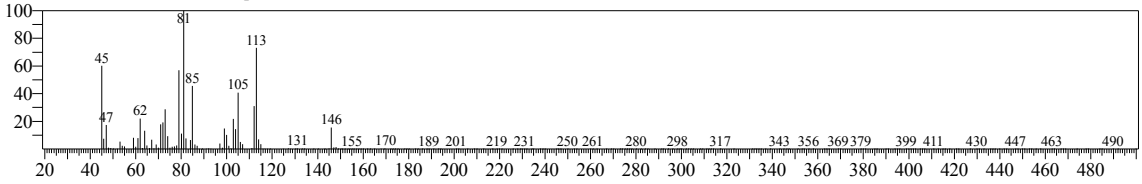
Hit#:2 Entry:28253 Library:Wiley9.lib
SI:77 Formula:C8H18O CAS:6169-06-8 MolWeight:130 RetIndex:0
CompName:2-Octanol, (S)- $$ (S)-(+)-2-Octanol $$ 2-Octanol $$ 2-Octanol (computer-generated name)



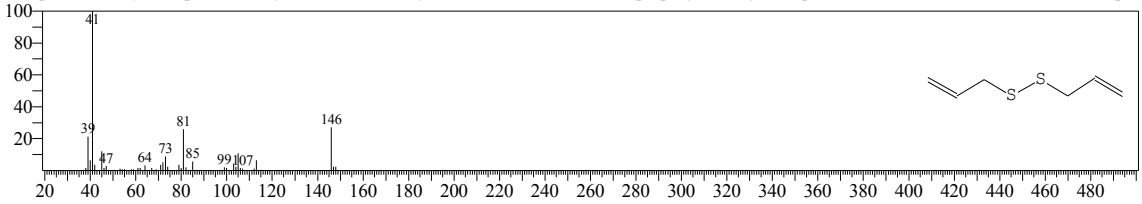
Hit#:3 Entry:28028 Library:Wiley9.lib
SI:77 Formula:C8H18O CAS:123-96-6 MolWeight:130 RetIndex:0
CompName:2-Octanol (CAS) $$ n-Octan-2-ol $$ Capryl alcohol $$ s-Octyl alcohol $$ 1-Methyl-1-heptanol $$ Methylhexylcarbinol $$ Hexylmethylcarbinc



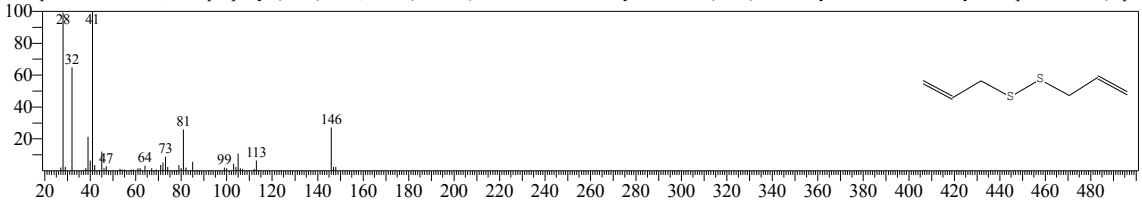
<< Target >>
Line#:10 R.Time:7.580(Scan#:917) MassPeaks:250
RawMode:Averaged 7.575-7.585(916-918) BasePeak:81.05(529805)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



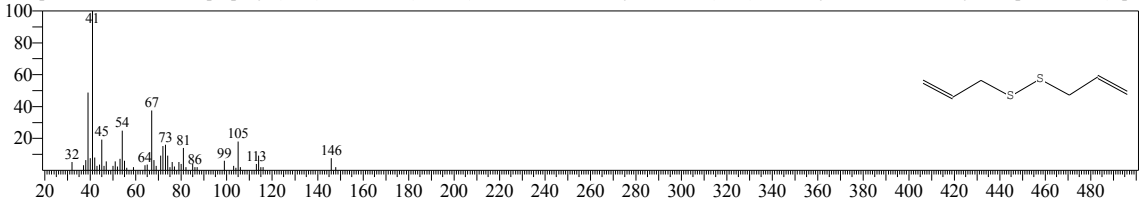
Hit#:1 Entry:7901 Library:NIST08s.LIB
SI:81 Formula:C6H10S2 CAS:2179-57-9 MolWeight:146 RetIndex:1099
CompName:Diallyl disulphide $$ Allyl disulfide $$ Diallyl disulfide $$ Disulfide, di-2-propenyl $$ Allyl disulphide $$ 4,5-Dithia-1,7-octadiene $$ 2-Proper



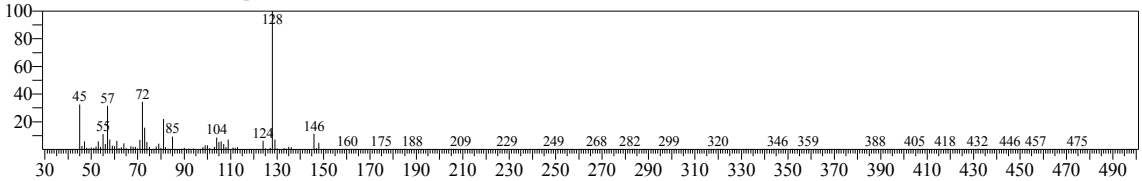
Hit#:2 Entry:45135 Library:Wiley9.lib
SI:81 Formula:C6H10S2 CAS:2179-57-9 MolWeight:146 RetIndex:0
CompName:Disulfide, di-2-propenyl (CAS) $$ 1,2-BIS(ALLYL)DISULFANE $$ Allyl disulfide (CAS) $$ Diallyl disulfide $$ Diallyl disulphide $$ Di(2-pr



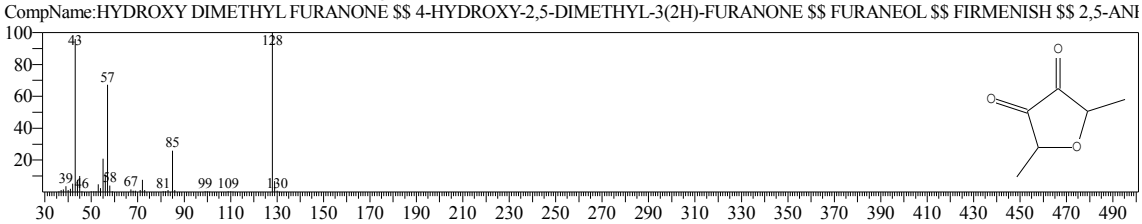
Hit#:3 Entry:45134 Library:Wiley9.lib
SI:71 Formula:C6H10S2 CAS:2179-57-9 MolWeight:146 RetIndex:0
CompName:Disulfide, di-2-propenyl (CAS) $$ 1,2-BIS(ALLYL)DISULFANE $$ Allyl disulfide (CAS) $$ Diallyl disulfide $$ Diallyl disulphide $$ Di(2-pr



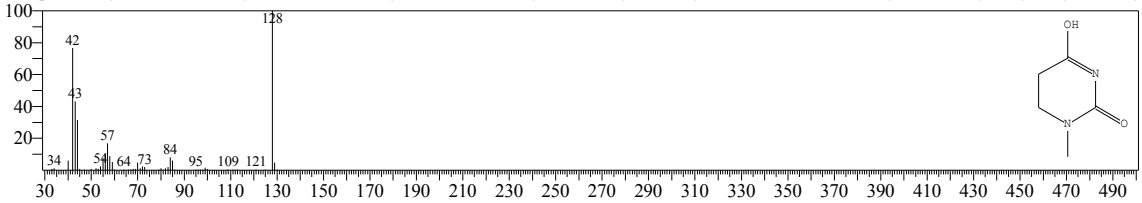
<< Target >>
Line#:11 R.Time:7.830(Scan#:967) MassPeaks:239
RawMode:Averaged 7.825-7.835(966-968) BasePeak:128.00(349521)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



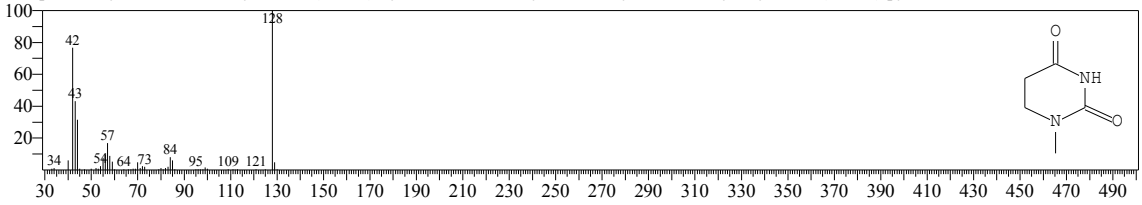
Hit#:1 Entry:25096 Library:Wiley9.lib
SI:72 Formula:C6H8O3 CAS:3658-77-3 MolWeight:128 RetIndex:0
CompName:HYDROXY DIMETHYL FURANONE $$ 4-HYDROXY-2,5-DIMETHYL-3(2H)-FURANONE $$ FURANEOL $$ FIRMENISH $$ 2,5-ANI



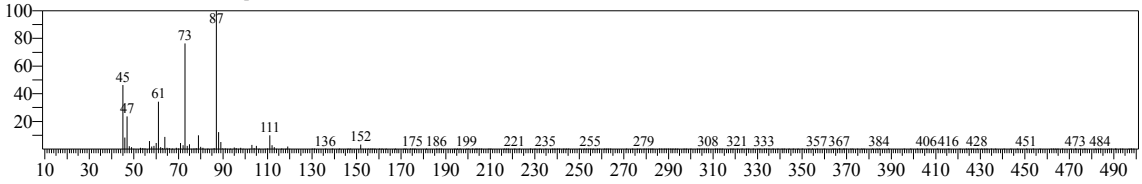
Hit#:2 Entry:24945 Library:Wiley9.lib
SI:70 Formula:C5H8N2O2 CAS:696-11-7 MolWeight:128 RetIndex:0
CompName:Hydrouracil, 1-methyl- $$ 2,4(1H,3H)-Pyrimidinedione, dihydro-1-methyl- $$ 2-Pyrimidinone, 1,2,5,6-tetrahydro-1-methyl-4-hydroxy- $$ 6-Hy



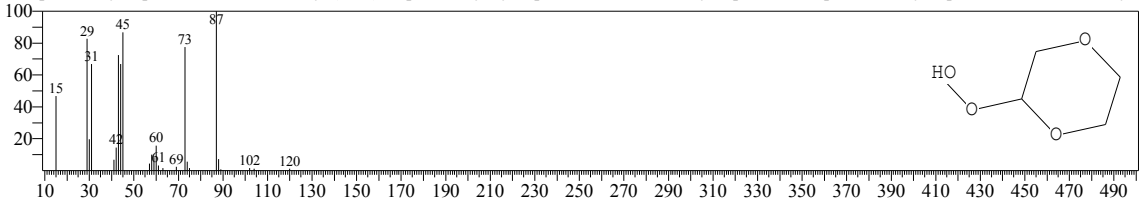
Hit#:3 Entry:4902 Library:NIST08s.LIB
SI:70 Formula:C5H8N2O2 CAS:696-11-7 MolWeight:128 RetIndex:1167
CompName:Hydrouracil, 1-methyl- $$ 2,4(1H,3H)-Pyrimidinedione, dihydro-1-methyl- $$ 1-Methyldihydro-2,4(1H,3H)-pyrimidinedione # $$



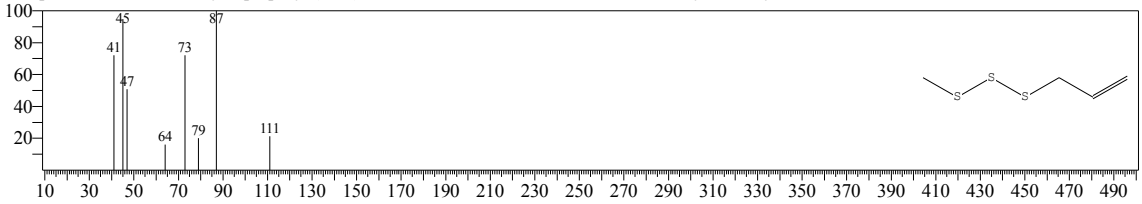
<< Target >>
Line#:12 R.Time:8.555(Scan#:1112) MassPeaks:249
RawMode:Averaged 8.550-8.560(1111-1113) BasePeak:87.00(1476239)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



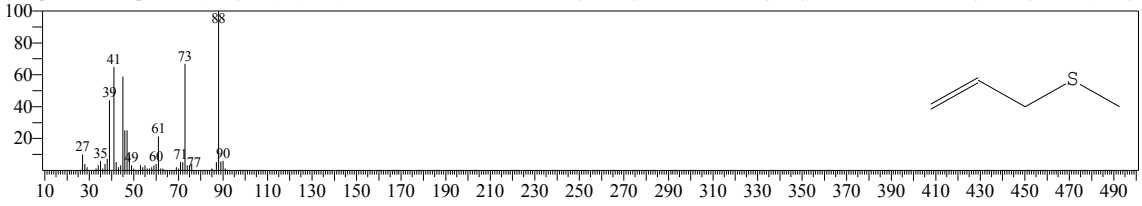
Hit#:1 Entry:18871 Library:Wiley9.lib
SI:78 Formula:C4H8O4 CAS:4722-59-2 MolWeight:120 RetIndex:0
CompName:Hydroperoxide, 1,4-dioxan-2-yl (CAS) $$ p-Dioxanyl hydroperoxide $$ Dioxane hydroperoxide $$ p-Dioxane hydroperoxide $$ 1,4-Dioxanyl



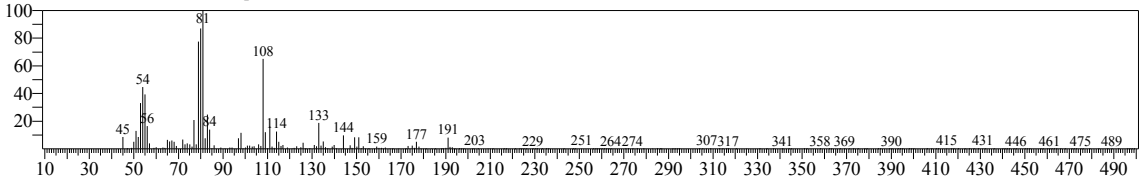
Hit#:2 Entry:52570 Library:Wiley9.lib
SI:77 Formula:C4H8S3 CAS:34135-85-8 MolWeight:152 RetIndex:0
CompName:Trisulfide, methyl 2-propenyl (CAS) $$ METHYL ALLYL TRISULFIDE $$ 1-Allyl-3-methyltrisulfane $$ 1-ALLYL-3-METHYLTRISULFAN



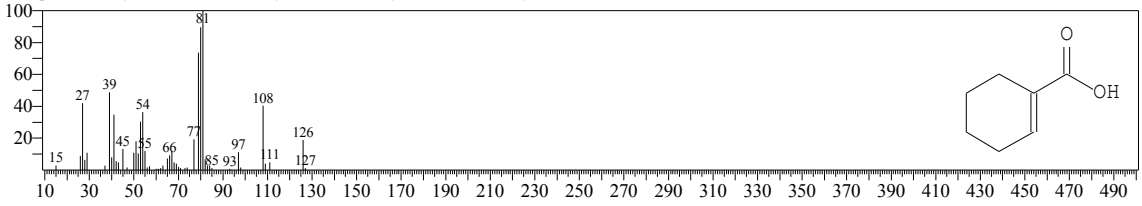
Hit#:3 Entry:4647 Library:Wiley9.lib
SI:75 Formula:C4H8S CAS:10152-76-8 MolWeight:88 RetIndex:0
CompName:1-Propene, 3-(methylthio)- (CAS) $$ 4-THIA-1-PENTENE $$ Allyl methyl sulfide $$ Methyl allyl sulfide $$ Sulfide, allyl methyl $$ 3-(Methy



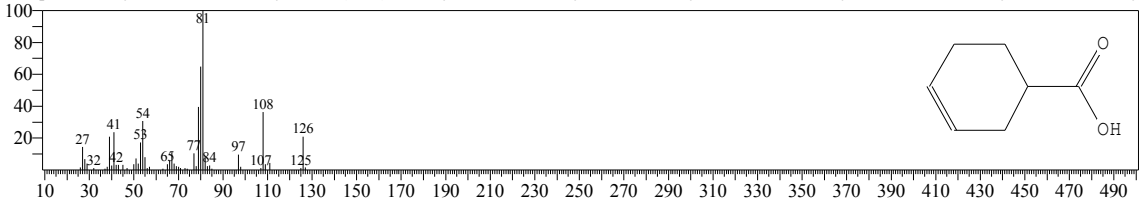
<< Target >>
Line#:13 R.Time:8.895(Scan#:1180) MassPeaks:248
RawMode:Averaged 8.890-8.900(1179-1181) BasePeak:81.05(174348)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



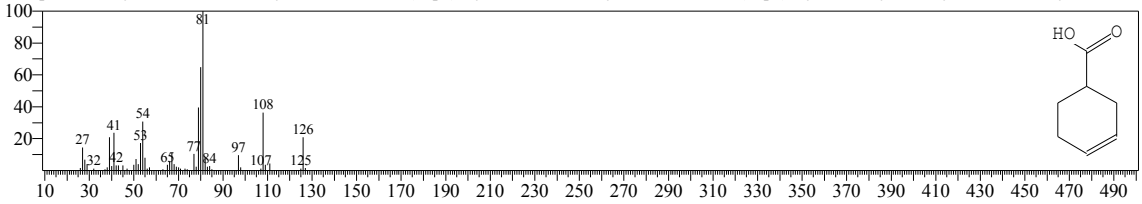
Hit#:1 Entry:4662 Library:NIST08s.LIB
SI:82 Formula:C7H10O2 CAS:636-82-8 MolWeight:126 RetIndex:1148
CompName:1-Cyclohexene-1-carboxylic acid $$ 1-Cyclohexenecarboxylic acid $$



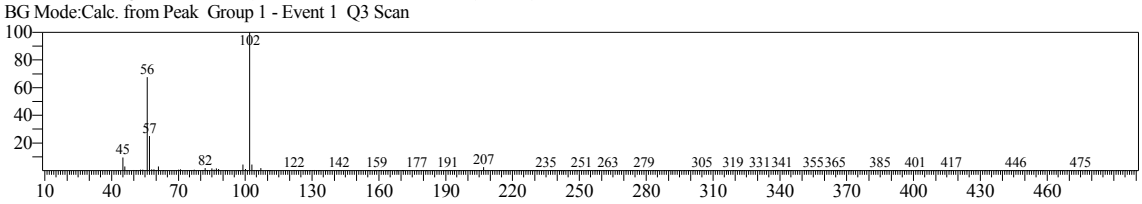
Hit#:2 Entry:23276 Library:Wiley9.lib
SI:77 Formula:C7H10O2 CAS:4771-80-6 MolWeight:126 RetIndex:0
CompName:3-Cyclohexene-1-carboxylic acid (CAS) $$ 3-Cyclohexenecarboxylic acid $$ 1-Cyclohexene-4-carboxylic acid $$ .delta.3-Cyclohexenecarboxy



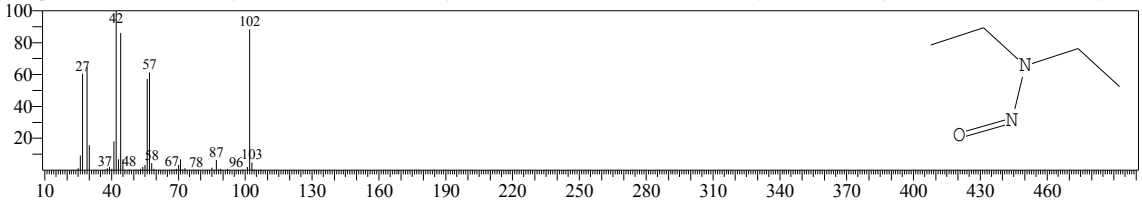
Hit#:3 Entry:4663 Library:NIST08s.LIB
SI:77 Formula:C7H10O2 CAS:4771-80-6 MolWeight:126 RetIndex:1119
CompName:3-Cyclohexene-1-carboxylic acid $$ .delta.(Sup3)-Cyclohexenecarboxylic acid $$ .delta.(Sup3)-Cyclohexenylcarboxylic acid $$ 1-Cyclohexene



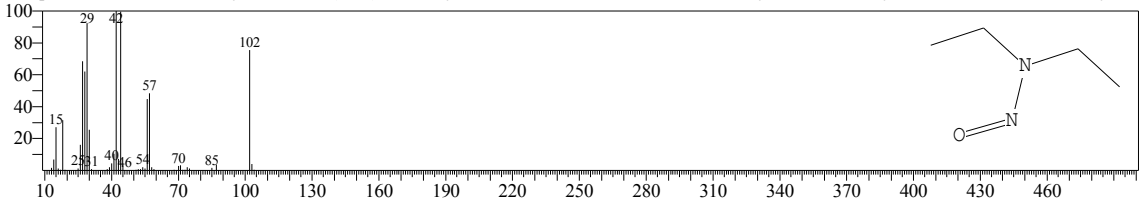
<< Target >>
Line#:14 R.Time:9.130(Scan#:1227) MassPeaks:243
RawMode:Averaged 9.125-9.135(1226-1228) BasePeak:102.00(572027)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



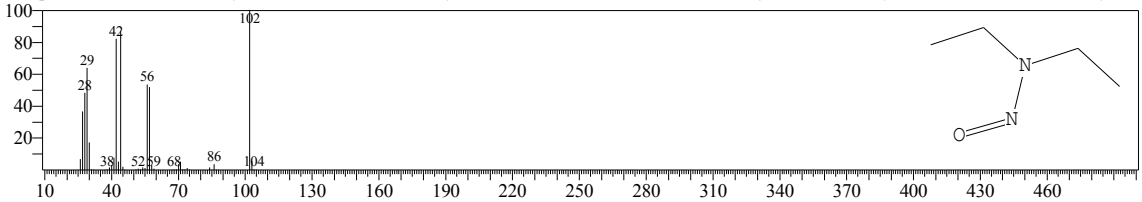
Hit#:1 Entry:9021 Library:Wiley9.lib
SI:87 Formula:C4H10N2O CAS:55-18-5 MolWeight:102 RetIndex:0
CompName:Ethanamine, N-ethyl-N-nitroso- (CAS) $$ Diethylnitrosamine $$ DENA $$ N-Nitrosodiethylamine $$ Diethylnitrosamide $$ Nitrosodiethylami



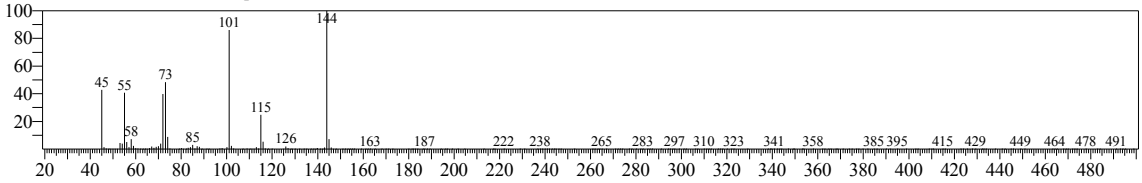
Hit#:2 Entry:9019 Library:Wiley9.lib
SI:87 Formula:C4H10N2O CAS:55-18-5 MolWeight:102 RetIndex:0
CompName:Ethanamine, N-ethyl-N-nitroso- (CAS) $$ Diethylnitrosamine $$ DENA $$ N-Nitrosodiethylamine $$ Diethylnitrosamide $$ Nitrosodiethylami



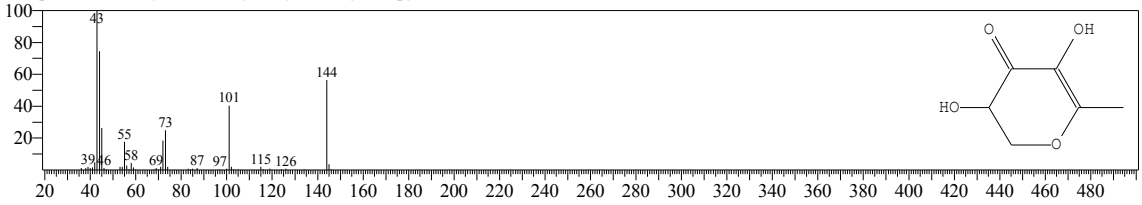
Hit#:3 Entry:9017 Library:Wiley9.lib
SI:87 Formula:C4H10N2O CAS:55-18-5 MolWeight:102 RetIndex:0
CompName:Ethanamine, N-ethyl-N-nitroso- (CAS) $$ Diethylnitrosamine $$ DENA $$ N-Nitrosodiethylamine $$ Diethylnitrosamide $$ Nitrosodiethylami



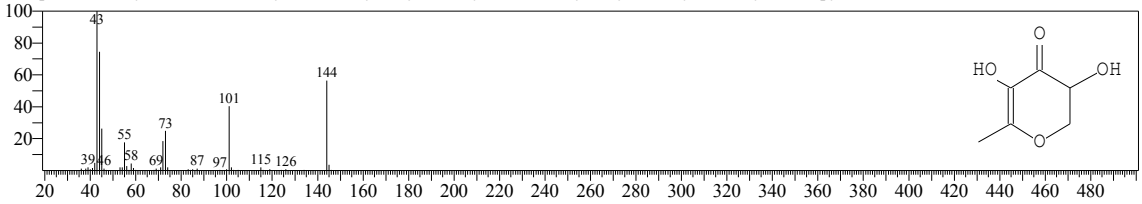
<< Target >>
Line#:15 R.Time:9.265(Scan#:1254) MassPeaks:241
RawMode:Averaged 9.260-9.270(1253-1255) BasePeak:144.00(3352713)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



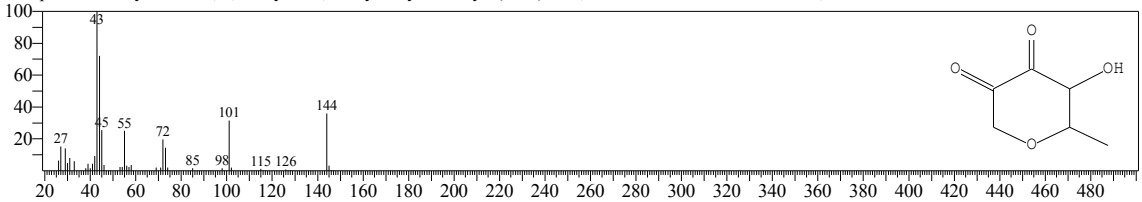
Hit#:1 Entry:42591 Library:Wiley9.lib
SI:93 Formula:C6H8O4 CAS:28564-83-2 MolWeight:144 RetIndex:0
CompName:2,3-Dihydro-3,5-dihydroxy-6-methyl-4H-pyran-4-one



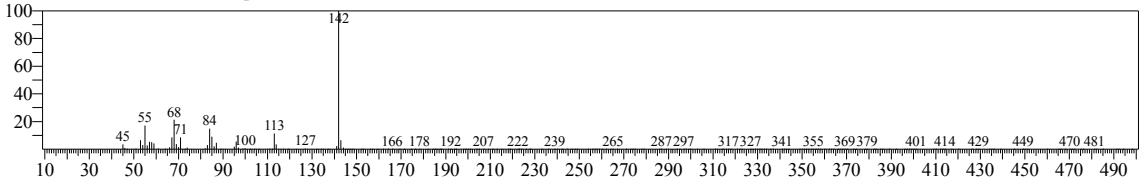
Hit#:2 Entry:7575 Library:NIST08s.LIB
SI:93 Formula:C6H8O4 CAS:28564-83-2 MolWeight:144 RetIndex:1269
CompName:4H-Pyran-4-one, 2,3-dihydro-3,5-dihydroxy-6-methyl- $$ 3,5-Dihydroxy-6-methyl-2,3-dihydro-4H-pyran-4-one # $$



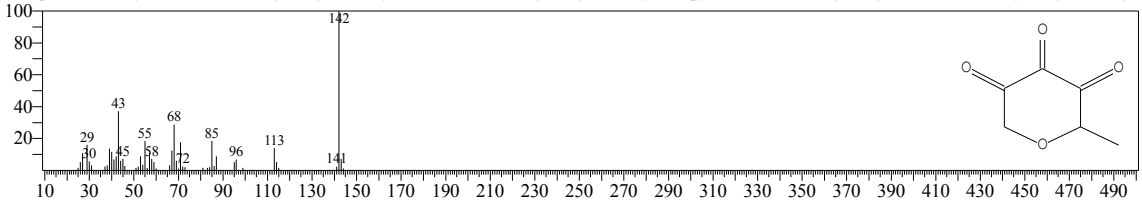
Hit#:3 Entry:42587 Library:Wiley9.lib
SI:90 Formula:C6H8O4 CAS:28564-83-2 MolWeight:144 RetIndex:0
CompName:4H-Pyran-4-one, 2,3-dihydro-3,5-dihydroxy-6-methyl- (CAS) $$ 3,5-DIHYDROXY-2-METHYL-5,6-DIHYDROPYRAN-4-ONE $$ 4H-PYR



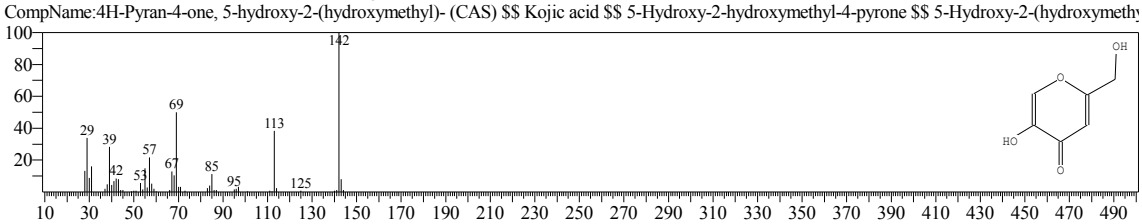
<< Target >>
Line#:16 R.Time:9.780(Scan#:1357) MassPeaks:247
RawMode:Averaged 9.775-9.785(1356-1358) BasePeak:141.95(857963)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



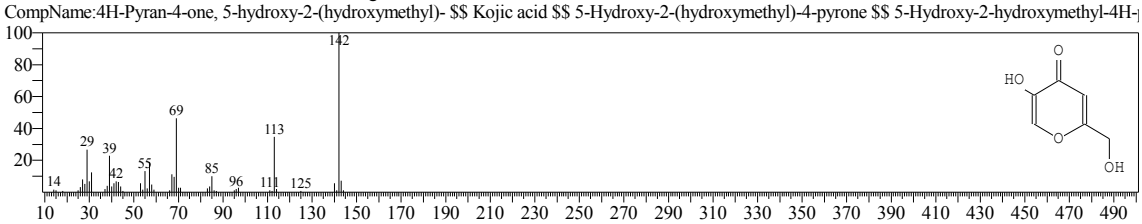
Hit#:1 Entry:40045 Library:Wiley9.lib
SI:90 Formula:C6H6O4 CAS:1073-96-7 MolWeight:142 RetIndex:0
CompName:4H-Pyran-4-one, 3,5-dihydroxy-2-methyl- (CAS) $$ 3,5-Dihydroxy-2-methyl-4H-pyran-4-one $$ 5-Hydroxymaltol $$ 3,5-Dihydroxy-2-methyl-



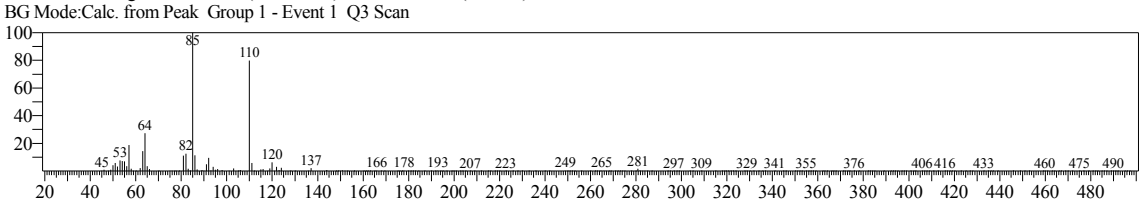
Hit#:2 Entry:40051 Library:Wiley9.lib
SI:85 Formula:C6H6O4 CAS:501-30-4 MolWeight:142 RetIndex:0
CompName:4H-Pyran-4-one, 5-hydroxy-2-(hydroxymethyl)- (CAS) $$ Kojic acid $$ 5-Hydroxy-2-hydroxymethyl-4-pyrone $$ 5-Hydroxy-2-(hydroxymeth



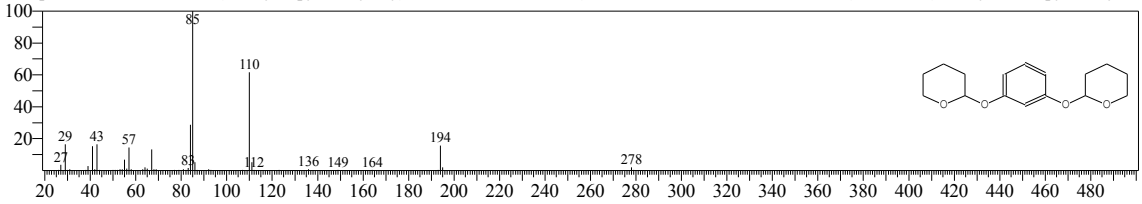
Hit#:3 Entry:7253 Library:NIST08s.LIB
SI:85 Formula:C6H6O4 CAS:501-30-4 MolWeight:142 RetIndex:1306
CompName:4H-Pyran-4-one, 5-hydroxy-2-(hydroxymethyl)- $$ Kojic acid $$ 5-Hydroxy-2-(hydroxymethyl)-4-pyrone $$ 5-Hydroxy-2-hydroxymethyl-4H-



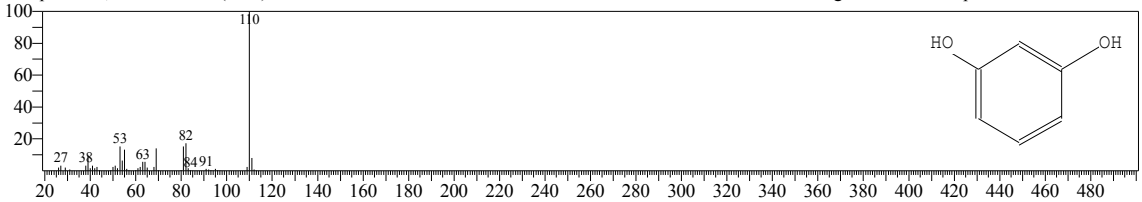
<< Target >>
Line#:17 R.Time:9.915(Scan#:1384) MassPeaks:256
RawMode:Averaged 9.910-9.920(1383-1385) BasePeak:85.00(788462)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



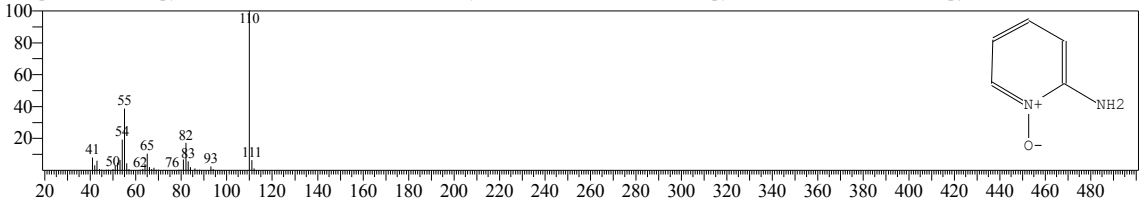
Hit#:1 Entry:324793 Library:Wiley9.lib
SI:76 Formula:C16H22O4 CAS:30778-88-2 MolWeight:278 RetIndex:0
CompName:Benzene, 1,3-bis(tetrahydropyran-2-yloxy)- $$ BENZOL, 1,3-BIS(TETRAHYDROPYRAN-2-YLOXY)- $$ 2-[3-(Tetrahydro-2H-pyran-2-ylox



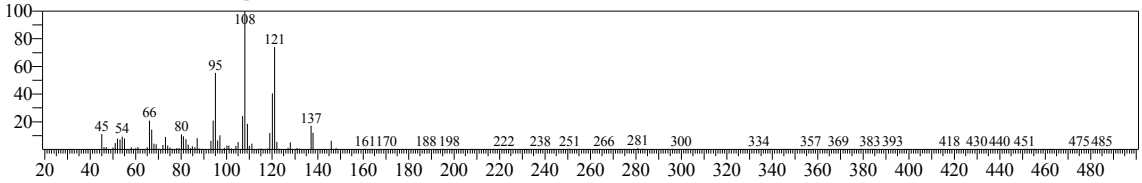
Hit#:2 Entry:12086 Library:Wiley9.lib
SI:71 Formula:C6H6O2 CAS:108-46-3 MolWeight:110 RetIndex:0
CompName:1,3-Benzenediol (CAS) $$ Resorcin $$ CI 76505 $$ C.I. 76505 $$ Resorcinol $$ Nako TGG $$ Pelagol RS $$ Developer R $$ Fourrine 79 $$



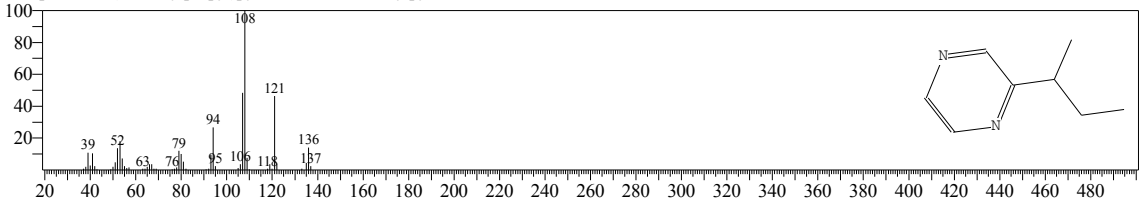
Hit#:3 Entry:12008 Library:Wiley9.lib
SI:71 Formula:C5H6N2O CAS:16867-03-1 MolWeight:110 RetIndex:0
CompName:2-Aminopyridine-N-oxide $$ Nc1ccccn1=O $$ 2-Pyridinamine, 1-oxide $$ 2-Aminopyridine N-oxide $$ 2-Amino-1-pyridine 1-oxide $$ 2-Am



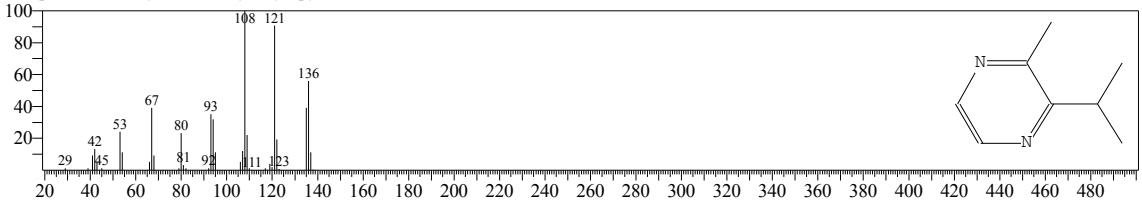
<< Target >>
Line#:18 R.Time:10.505(Scan#:1502) MassPeaks:214
RawMode:Averaged 10.500-10.510(1501-1503) BasePeak:108.00(182327)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



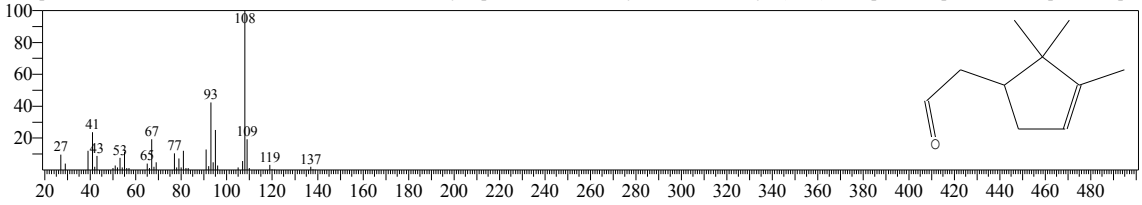
Hit#:1 Entry:32950 Library:Wiley9.lib
SI:71 Formula:C8H12N2 CAS:29460-93-3 MolWeight:136 RetIndex:0
CompName:2-(1-Methylpropyl)pyrazine $$ 2-Sec-butylpyrazine $$ 2-SEC-BUTYLPYRAZINE #



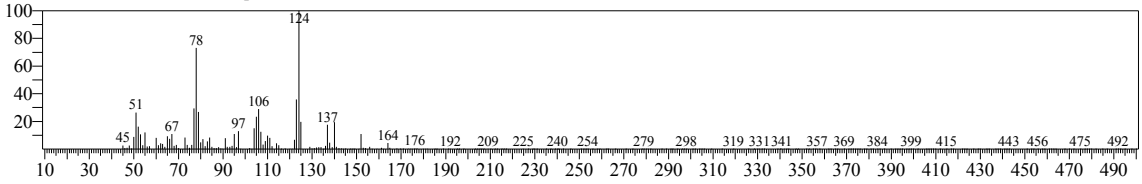
Hit#:2 Entry:32954 Library:Wiley9.lib
SI:69 Formula:C8H12N2 CAS:0-00-0 MolWeight:136 RetIndex:0
CompName:2-methyl-3-(1-methylethyl)-pyrazine



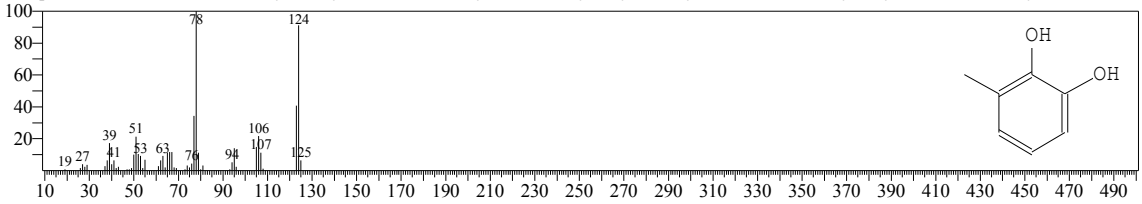
Hit#:3 Entry:54150 Library:Wiley9.lib
SI:69 Formula:C10H16O CAS:4501-58-0 MolWeight:152 RetIndex:0
CompName:.ALPHA.-CAMPHOLENE ALDEHYDE $$ 3-Cyclopentene-1-acetaldehyde, 2,2,3-trimethyl- (CAS) $$ .alpha.-Campholenal $$ .alpha.-Camph



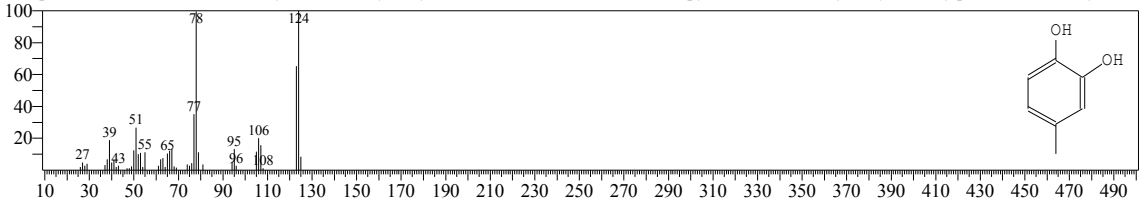
<< Target >>
Line#:19 R.Time:10.720(Scan#:1545) MassPeaks:276
RawMode:Averaged 10.715-10.725(1544-1546) BasePeak:124.05(203506)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



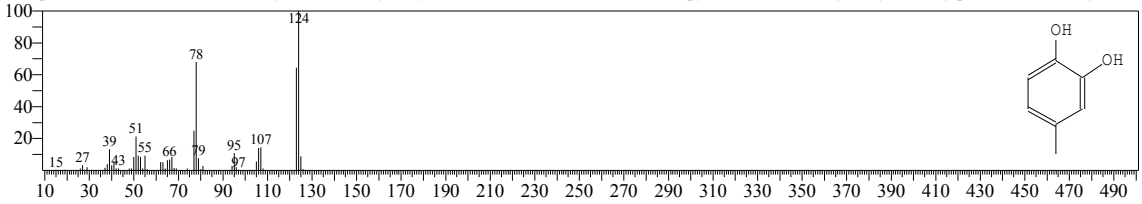
Hit#:1 Entry:4387 Library:NIST08s.LIB
SI:79 Formula:C7H8O2 CAS:488-17-5 MolWeight:124 RetIndex:1235
CompName:1,2-Benzenediol, 3-methyl- $$ Pyrocatechol, 3-methyl- $$ 1,2-Dihydroxy-3-methylbenzene $$ 2,3-Dihydroxytoluene $$ 3-Methyl-1,2-benzenec



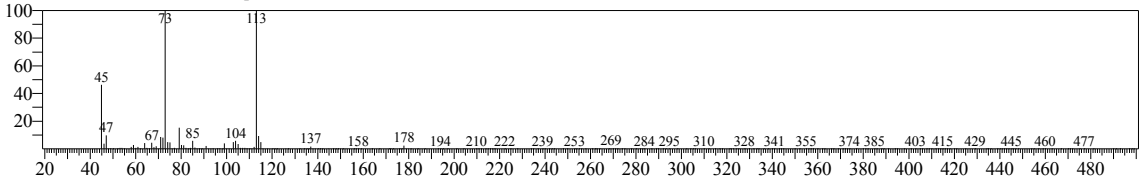
Hit#:2 Entry:4396 Library:NIST08s.LIB
SI:78 Formula:C7H8O2 CAS:452-86-8 MolWeight:124 RetIndex:1235
CompName:1,2-Benzenediol, 4-methyl- $$ 3,4-Dihydroxytoluene $$ Homocatechol $$ Homopyrocatechol $$ 2-Hydroxy-4-methylphenol $$ 4-Methyl-1,2-l



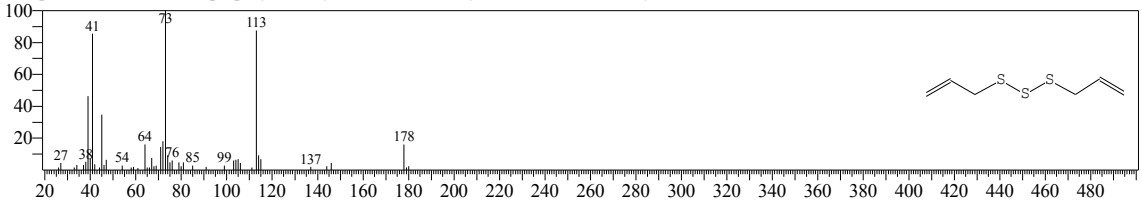
Hit#:3 Entry:4397 Library:NIST08s.LIB
SI:78 Formula:C7H8O2 CAS:452-86-8 MolWeight:124 RetIndex:1235
CompName:1,2-Benzenediol, 4-methyl- $$ 3,4-Dihydroxytoluene $$ Homocatechol $$ Homopyrocatechol $$ 2-Hydroxy-4-methylphenol $$ 4-Methyl-1,2-l



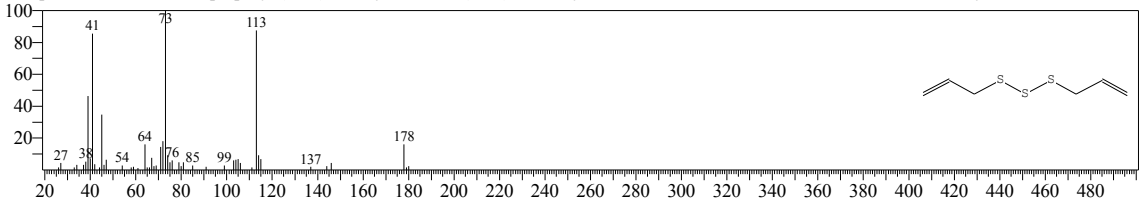
<< Target >>
Line#:20 R.Time:11.075(Scan#:1616) MassPeaks:262
RawMode:Averaged 11.070-11.080(1615-1617) BasePeak:72.95(2928880)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



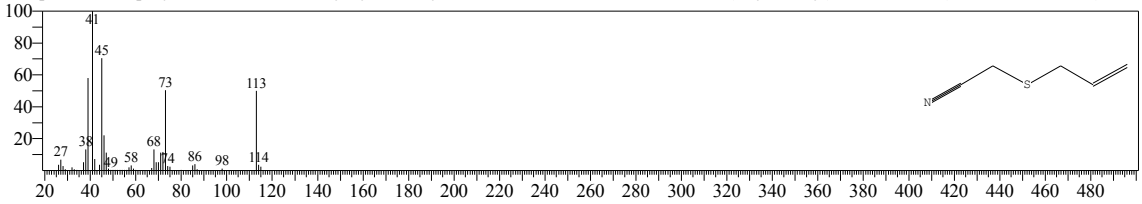
Hit#:1 Entry:13262 Library:NIST08s.LIB
SI:89 Formula:C6H10S3 CAS:2050-87-5 MolWeight:178 RetIndex:1350
CompName:Trisulfide, di-2-propenyl $$ Allyl trisulfide $$ Diallyl trisulfide $$ 1,3-Diallyltrisulfane # $$



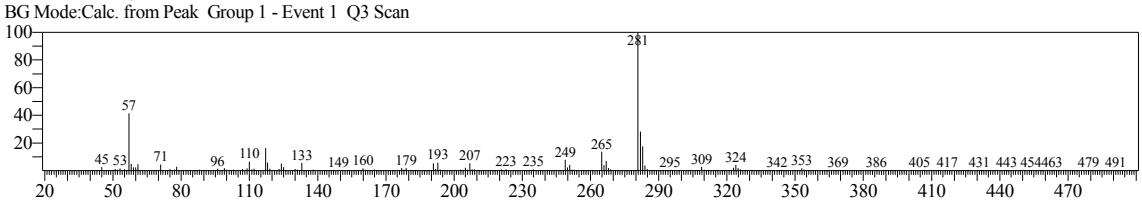
Hit#:2 Entry:96316 Library:Wiley9.lib
SI:89 Formula:C6H10S3 CAS:2050-87-5 MolWeight:178 RetIndex:0
CompName:Trisulfide, di-2-propenyl (CAS) $$ Allyl trisulfide (CAS) $$ Diallyl trisulfide $$ 4,5,6-trithia-nona-1,8-diene $$ 1,3-Diallyltrisulfane $$ 1,3-DI



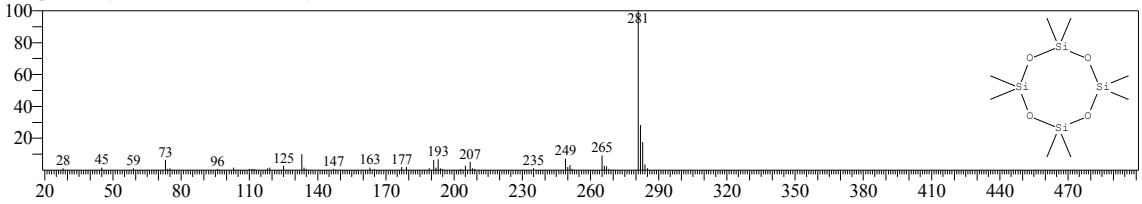
Hit#:3 Entry:14264 Library:Wiley9.lib
SI:78 Formula:C5H7NS CAS:105643-80-9 MolWeight:113 RetIndex:0
CompName:2-Propenylthioacetonitrile $$ Allyl cyanomethyl sulfide $$ CH2=CHCH2SCH2CN $$ (Allylsulfanyl)acetonitrile $$ (ALLYLSULFANYL)ACE



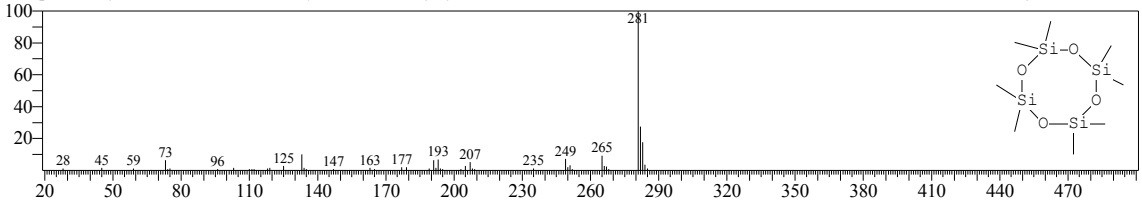
<< Target >>
Line#:21 R.Time:11.220(Scan#:1645) MassPeaks:305
RawMode:Averaged 11.215-11.225(1644-1646) BasePeak:280.90(390397)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



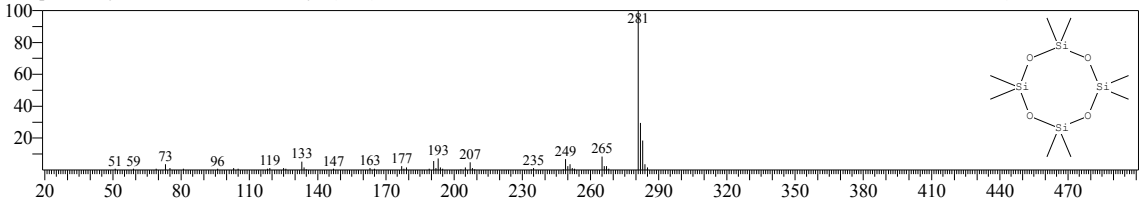
Hit#:1 Entry:363882 Library:Wiley9.lib
SI:76 Formula:C8H24O4Si4 CAS:556-67-2 MolWeight:296 RetIndex:0
CompName:Cyclotetrasiloxane, octamethyl- (CAS) $$ 1,1,3,3,5,5,7,7-OCTAMETHYL-CYCLOOCTASILOXANE $$ OCTAMETHYL-(CYCLIC TETRA



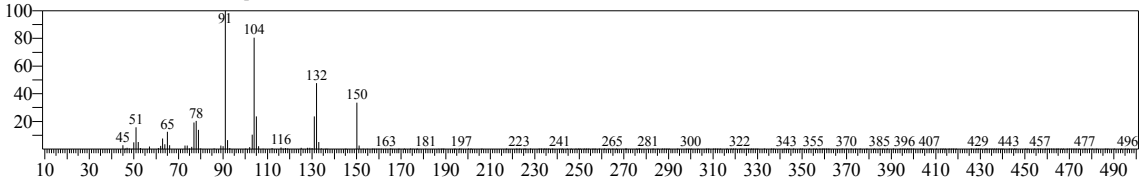
Hit#:2 Entry:24114 Library:NIST08s.LIB
SI:76 Formula:C8H24O4Si4 CAS:556-67-2 MolWeight:296 RetIndex:827
CompName:Cyclotetrasiloxane, octamethyl- $$ Oktamethylcyklotetrasiloxan $$ NUC Silicone VS 7207 $$ CO9810 $$ O9810 $$ Octamethyltetrasiloxane



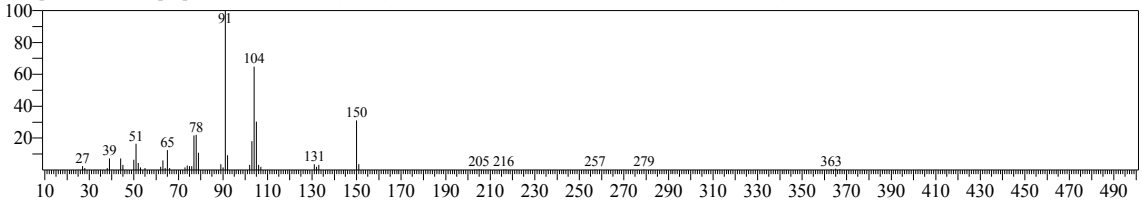
Hit#:3 Entry:363881 Library:Wiley9.lib
SI:74 Formula:C8H24O4Si4 CAS:556-67-2 MolWeight:296 RetIndex:0
CompName:Cyclotetrasiloxane, octamethyl- (CAS) $$ 1,1,3,3,5,5,7,7-OCTAMETHYL-CYCLOOCTASILOXANE $$ OCTAMETHYL-(CYCLIC TETRA



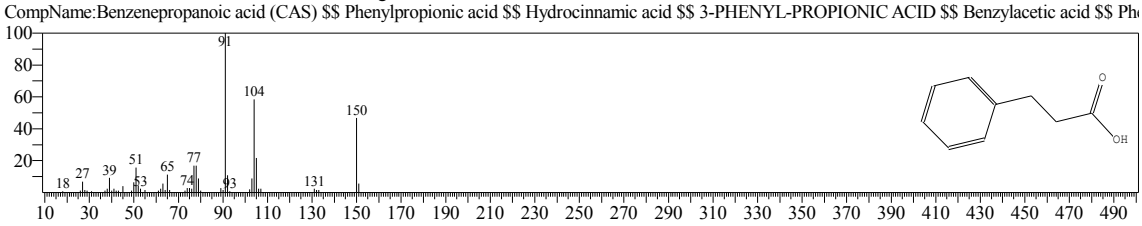
<< Target >>
Line#:22 R.Time:11.740(Scan#:1749) MassPeaks:238
RawMode:Averaged 11.735-11.745(1748-1750) BasePeak:91.00(346773)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



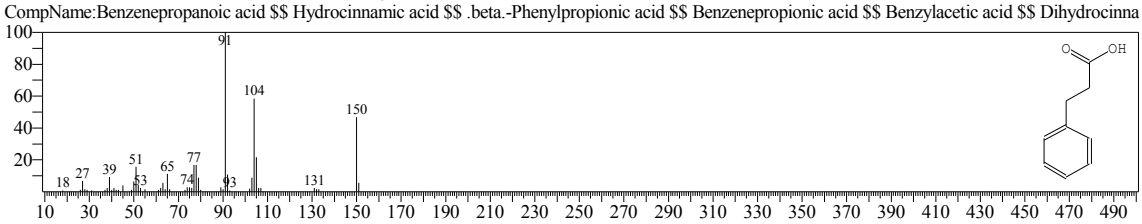
Hit#:1 Entry:275191 Library:Wiley9.lib
SI:91 Formula:C9H10AgO2 CAS:75112-79-7 MolWeight:257 RetIndex:0
CompName:Benzenepropanoic acid, silver(1+) salt (CAS) $$ BENZENEPROPANOIC ACID, SILVER(+1) SALT



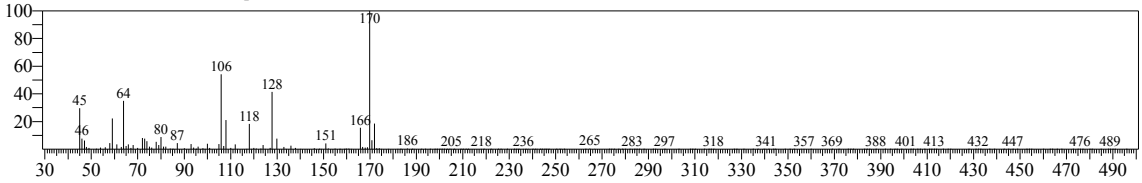
Hit#:2 Entry:50126 Library:Wiley9.lib
SI:90 Formula:C9H10O2 CAS:501-52-0 MolWeight:150 RetIndex:0
CompName:Benzenepropanoic acid (CAS) $$ Phenylpropionic acid $$ Hydrocinnamic acid $$ 3-PHENYL-PROPIONIC ACID $$ Benzylacetic acid $$ Ph



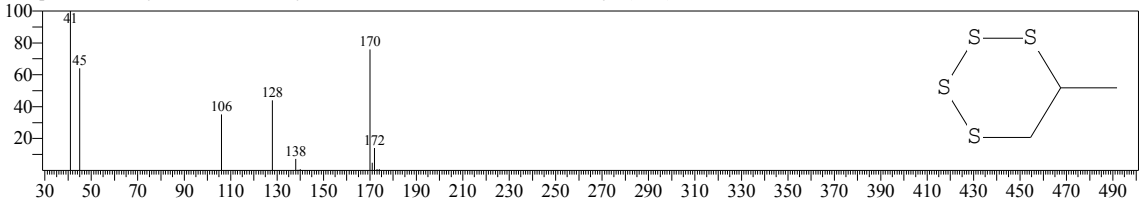
Hit#:3 Entry:8565 Library:NIST08s.LIB
SI:90 Formula:C9H10O2 CAS:501-52-0 MolWeight:150 RetIndex:1349
CompName:Benzenepropanoic acid $$ Hydrocinnamic acid $$ .beta.-Phenylpropionic acid $$ Benzenepropionic acid $$ Benzylacetic acid $$ Dihydrocinna



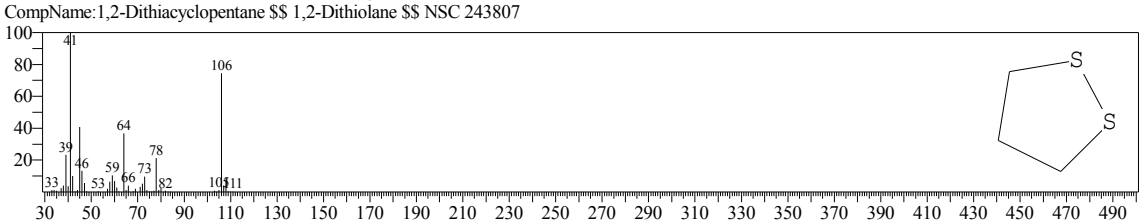
<< Target >>
Line#:23 R.Time:12.015(Scan#:1804) MassPeaks:229
RawMode:Averaged 12.010-12.020(1803-1805) BasePeak:169.90(200134)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



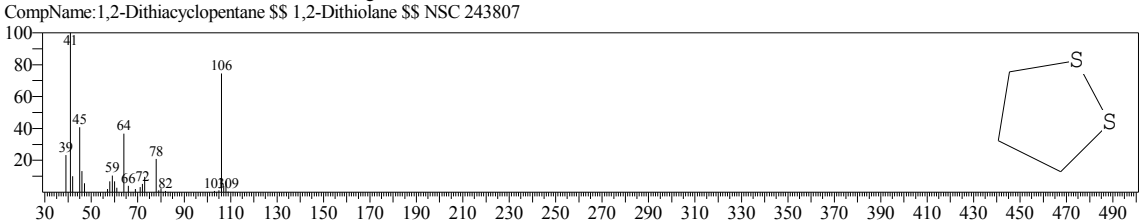
Hit#:1 Entry:82593 Library:Wiley9.lib
SI:65 Formula:C3H6S4 CAS:116664-30-3 MolWeight:170 RetIndex:0
CompName:5-methyl-1,2,3,4-tetrathia-cyclohexane $$ 1,2,3,4-Tetrathiane, 5-methyl- (CAS)



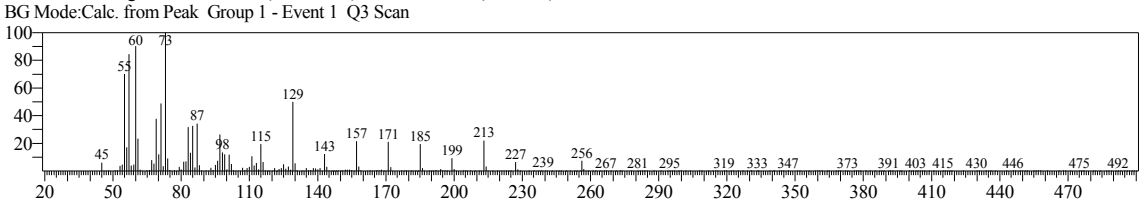
Hit#:2 Entry:10471 Library:Wiley9.lib
SI:58 Formula:C3H6S2 CAS:557-22-2 MolWeight:106 RetIndex:0
CompName:1,2-Dithiacyclopentane $$ 1,2-Dithiolane $$ NSC 243807



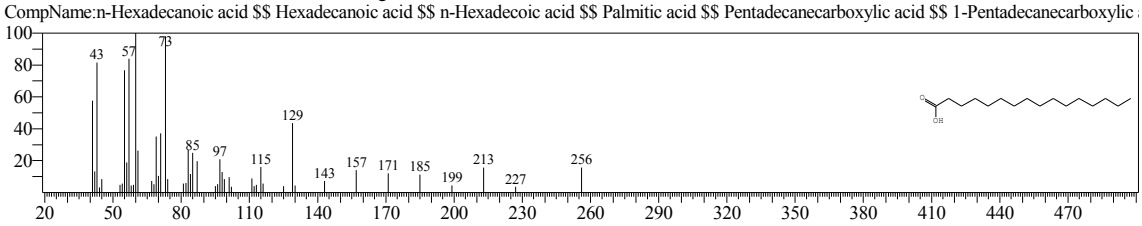
Hit#:3 Entry:10472 Library:Wiley9.lib
SI:57 Formula:C3H6S2 CAS:557-22-2 MolWeight:106 RetIndex:0
CompName:1,2-Dithiacyclopentane $$ 1,2-Dithiolane $$ NSC 243807



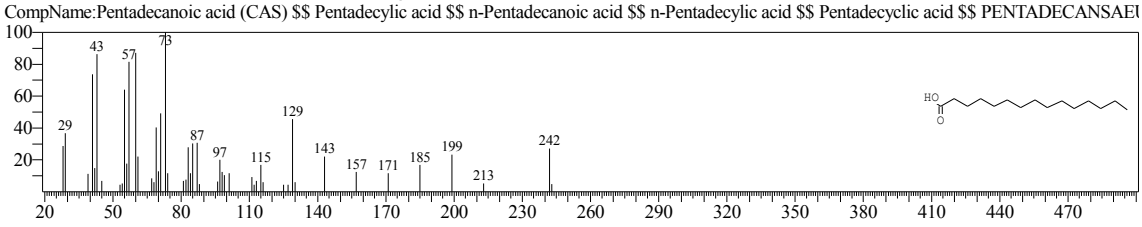
<< Target >>
Line#:24 R.Time:19.045(Scan#:3210) MassPeaks:321
RawMode:Averaged 19.040-19.050(3209-3211) BasePeak:73.00(1148193)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



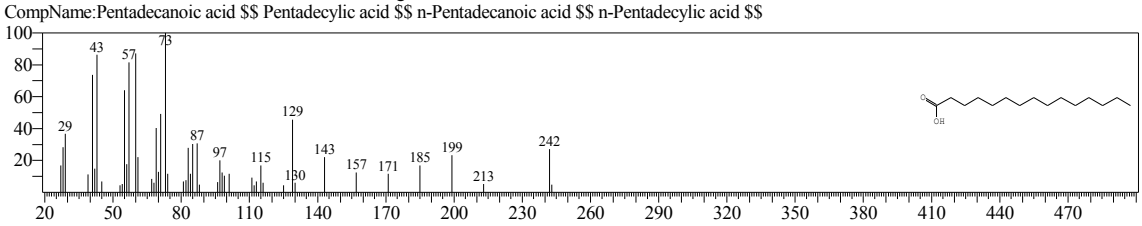
Hit#:1 Entry:21864 Library:NIST08s.LIB
SI:95 Formula:C16H32O2 CAS:57-10-3 MolWeight:256 RetIndex:1968
CompName:n-Hexadecanoic acid $$ Hexadecanoic acid $$ n-Hexadecoic acid $$ Palmitic acid $$ Pentadecanecarboxylic acid $$ 1-Pentadecanecarboxylic



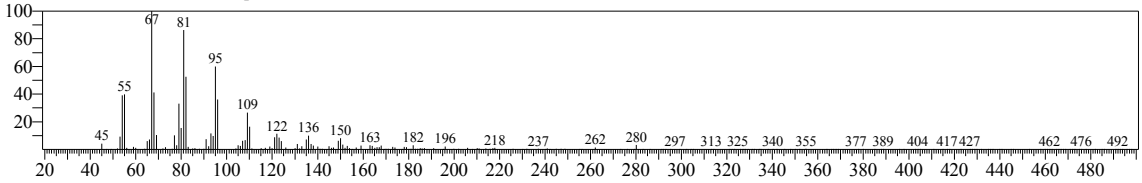
Hit#:2 Entry:241462 Library:Wiley9.lib
SI:93 Formula:C15H30O2 CAS:1002-84-2 MolWeight:242 RetIndex:0
CompName:Pentadecanoic acid (CAS) $$ Pentadecylic acid $$ n-Pentadecanoic acid $$ n-Pentadecylic acid $$ Pentadecyclic acid $$ PENTADECANSAE



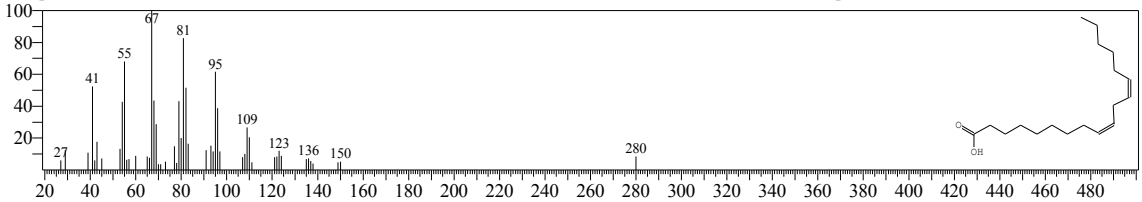
Hit#:3 Entry:20839 Library:NIST08s.LIB
SI:93 Formula:C15H30O2 CAS:1002-84-2 MolWeight:242 RetIndex:1869
CompName:Pentadecanoic acid $$ Pentadecylic acid $$ n-Pentadecanoic acid $$ n-Pentadecylic acid $$



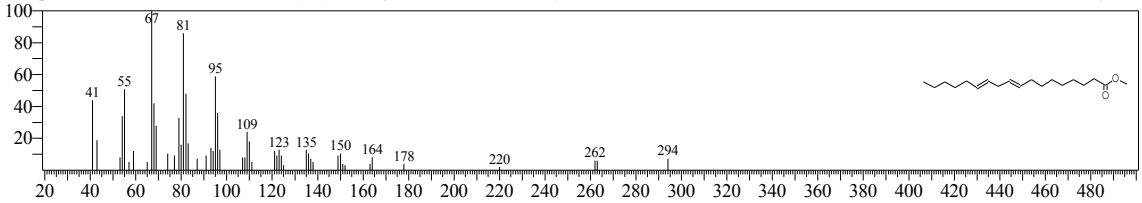
<< Target >>
Line#:25 R.Time:20.715(Scan#:3544) MassPeaks:256
RawMode:Averaged 20.710-20.720(3543-3545) BasePeak:67.00(228138)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



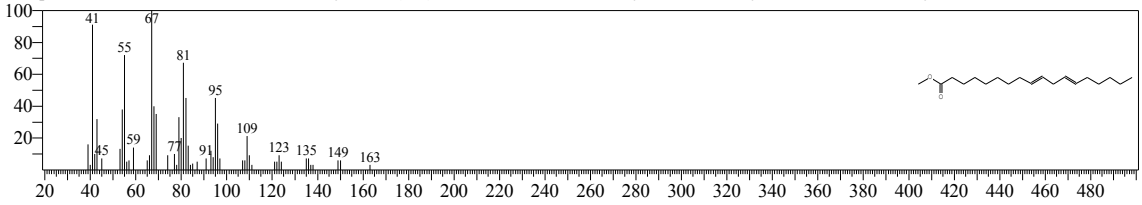
Hit#:1 Entry:23296 Library:NIST08s.LIB
SI:89 Formula:C18H32O2 CAS:60-33-3 MolWeight:280 RetIndex:2183
CompName:9,12-Octadecadienoic acid (Z,Z)- $$ cis-9,cis-12-Octadecadienoic acid $$ cis,cis-Linoleic acid $$ Grape seed oil $$ Linoleic $$ Linoleic acid $



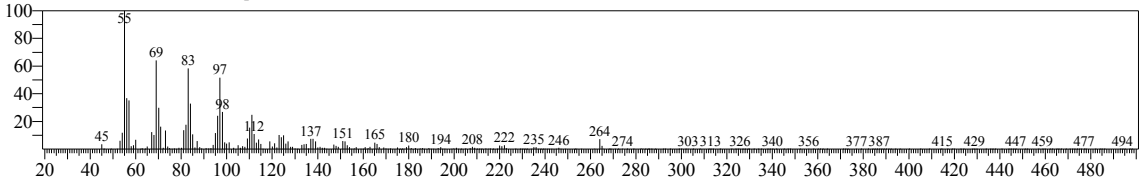
Hit#:2 Entry:361819 Library:Wiley9.lib
SI:88 Formula:C19H34O2 CAS:112-63-0 MolWeight:294 RetIndex:0
CompName:9,12-Octadecadienoic acid (Z,Z)-, methyl ester (CAS) $$ Methyl linoleate $$ METHYL CIS-9,CIS-12-OCTADECADIENOATE $$ Methyl oct



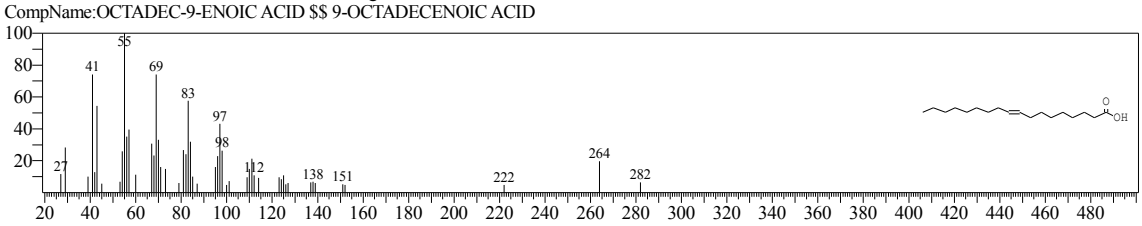
Hit#:3 Entry:24058 Library:NIST08s.LIB
SI:88 Formula:C19H34O2 CAS:2566-97-4 MolWeight:294 RetIndex:2093
CompName:9,12-Octadecadienoic acid, methyl ester, (E,E)- $$ Linolelaidic acid, methyl ester $$ Methyl linolelaidate $$ Methyl trans,trans-9,12-octadecadi



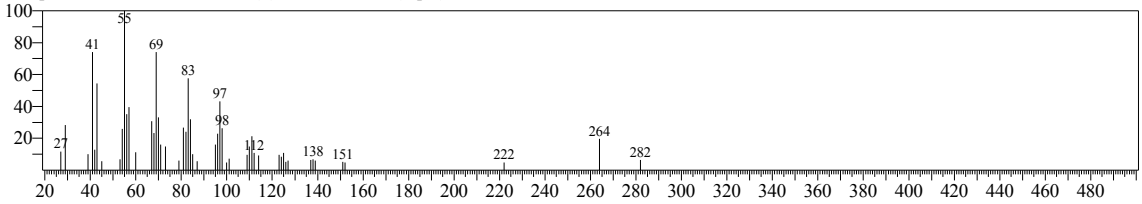
<< Target >>
Line#:26 R.Time:20.760(Scan#:3553) MassPeaks:294
RawMode:Averaged 20.755-20.765(3552-3554) BasePeak:55.00(288781)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



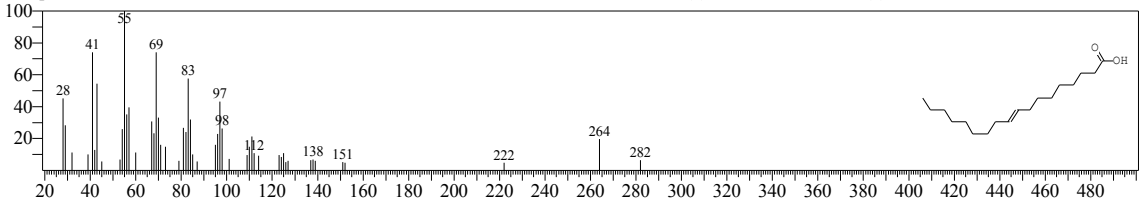
Hit#:1 Entry:335034 Library:Wiley9.lib
SI:92 Formula:C18H34O2 CAS:0-00-0 MolWeight:282 RetIndex:0
CompName:OCTADEC-9-ENOIC ACID $$ 9-OCTADECENOIC ACID



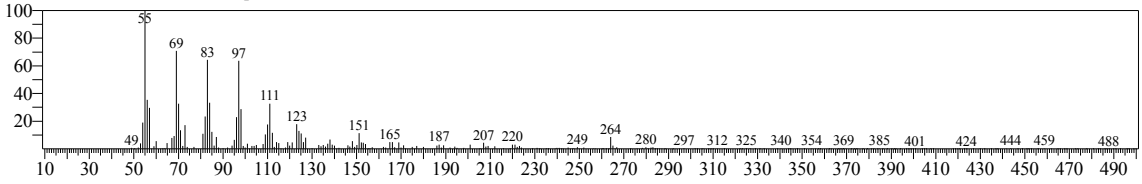
Hit#:2 Entry:335087 Library:Wiley9.lib
SI:92 Formula:C18H34O2 CAS:112-79-8 MolWeight:282 RetIndex:0
CompName:9-Octadecenoic acid, (E)- $$ trans-.delta.(sup 9)-Octadecenoic acid $$ trans-.delta.9-Octadecenoic acid $$ trans-Octadec-9-enoic acid $$ trans-



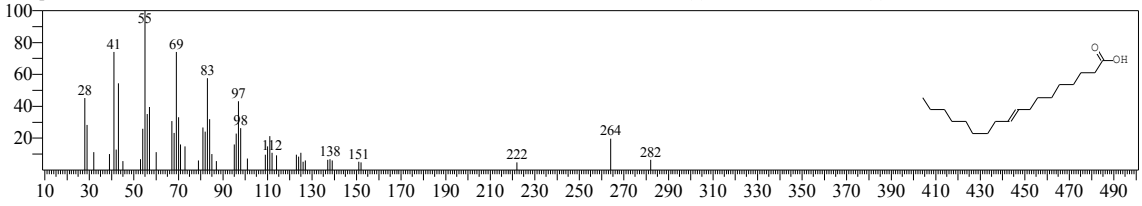
Hit#:3 Entry:335033 Library:Wiley9.lib
SI:92 Formula:C18H34O2 CAS:0-00-0 MolWeight:282 RetIndex:0
CompName:OCTADEC-9-ENOIC ACID $$ 9-OCTADECENOIC ACID $$ HEPTADECEN-(8)-CARBONSAEURE-(1)



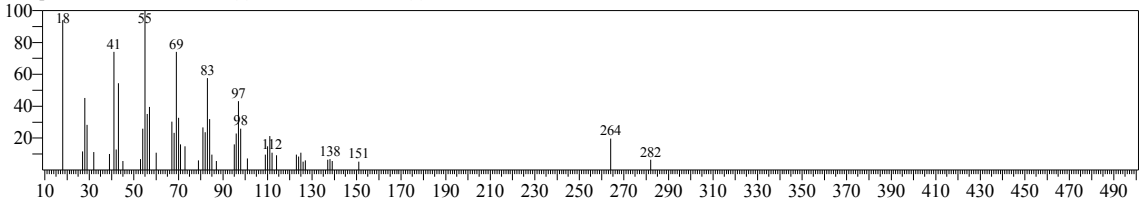
<< Target >>
Line#:27 R.Time:20.810(Scan#:3563) MassPeaks:266
RawMode:Averaged 20.805-20.815(3562-3564) BasePeak:55.00(84528)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



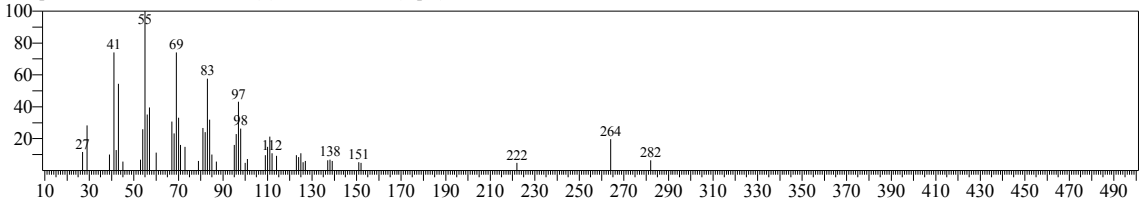
Hit#:1 Entry:335033 Library:Wiley9.lib
SI:90 Formula:C18H34O2 CAS:0-00-0 MolWeight:282 RetIndex:0
CompName:OCTADEC-9-ENOIC ACID $$ 9-OCTADECENOIC ACID $$ HEPTADECEN-(8)-CARBONSAEURE-(1)



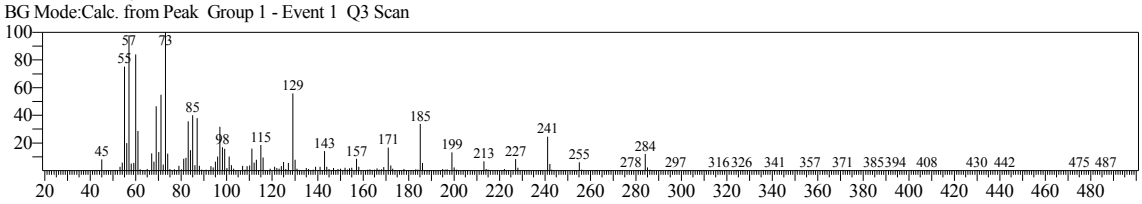
Hit#:2 Entry:335027 Library:Wiley9.lib
SI:90 Formula:C18H34O2 CAS:0-00-0 MolWeight:282 RetIndex:0
CompName:HEPTADECENE-(8)-CARBONIC ACID-(1)



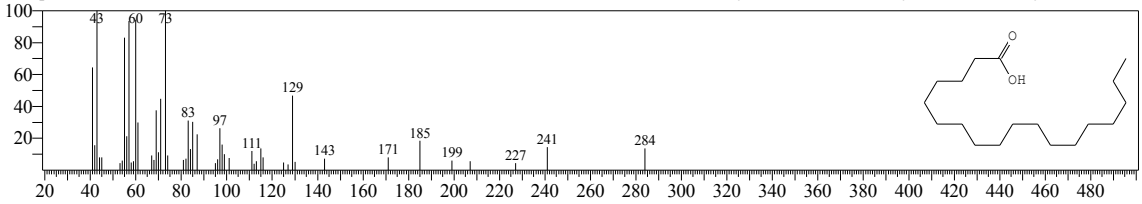
Hit#:3 Entry:335087 Library:Wiley9.lib
SI:90 Formula:C18H34O2 CAS:112-79-8 MolWeight:282 RetIndex:0
CompName:9-Octadecenoic acid, (E)- $$ trans-.delta.(sup 9)-Octadecenoic acid $$ trans-.delta.9-Octadecenoic acid $$ trans-Octadec-9-enoic acid $$ trans-



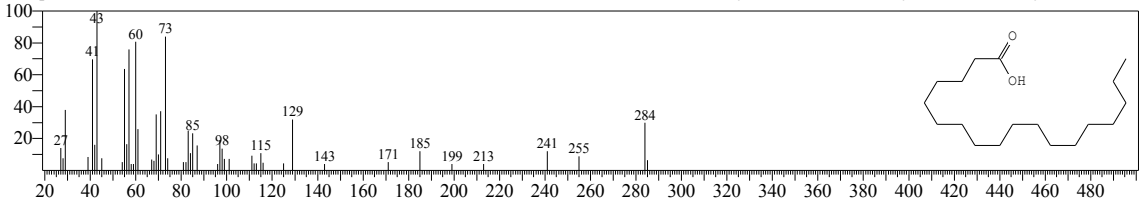
<< Target >>
Line#:28 R.Time:20.965(Scan#:3594) MassPeaks:271
RawMode:Averaged 20.960-20.970(3593-3595) BasePeak:73.00(264782)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



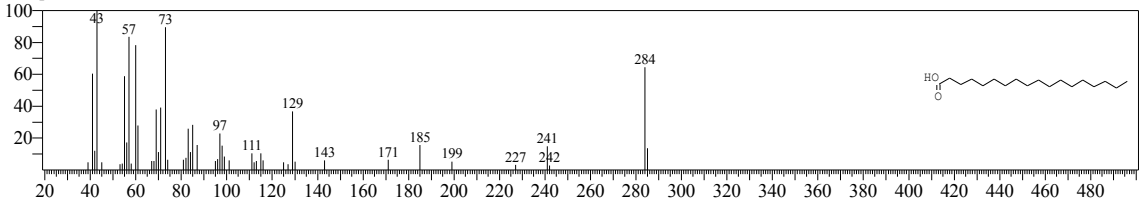
Hit#:1 Entry:23514 Library:NIST08s.LIB
SI:93 Formula:C18H36O2 CAS:57-11-4 MolWeight:284 RetIndex:2167
CompName:Octadecanoic acid $$ Stearic acid $$ n-Octadecanoic acid $$ Humko Industrene R $$ Hydrofol Acid 150 $$ Hystrene S-97 $$ Hystrene T-70 $



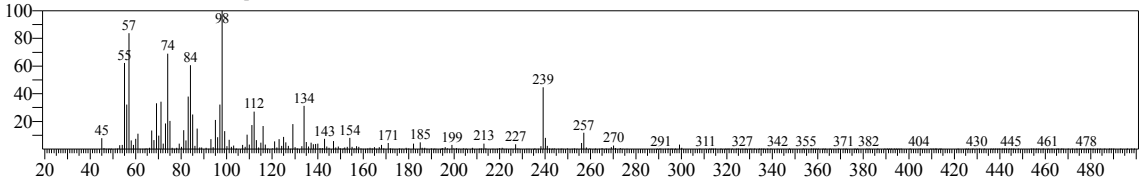
Hit#:2 Entry:23513 Library:NIST08s.LIB
SI:91 Formula:C18H36O2 CAS:57-11-4 MolWeight:284 RetIndex:2167
CompName:Octadecanoic acid $$ Stearic acid $$ n-Octadecanoic acid $$ Humko Industrene R $$ Hydrofol Acid 150 $$ Hystrene S-97 $$ Hystrene T-70 $



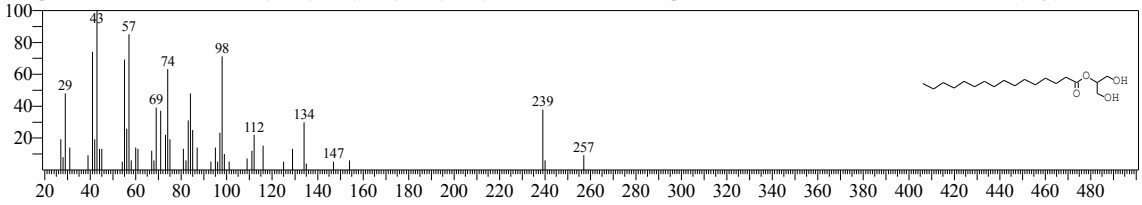
Hit#:3 Entry:339650 Library:Wiley9.lib
SI:91 Formula:C18H36O2 CAS:57-11-4 MolWeight:284 RetIndex:0
CompName:OCTADECANOIC ACID $$ STEARATE $$ STEARIC ACID $$ 1-HEPTADECANECARBOXYLIC ACID $$ 1-HEPTADECANECARBOX



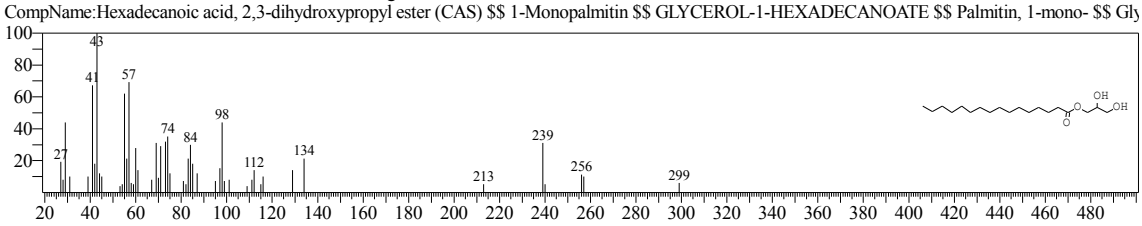
<< Target >>
Line#:29 R.Time:24.000(Scan#:4201) MassPeaks:311
RawMode:Averaged 23.995-24.005(4200-4202) BasePeak:98.00(218370)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



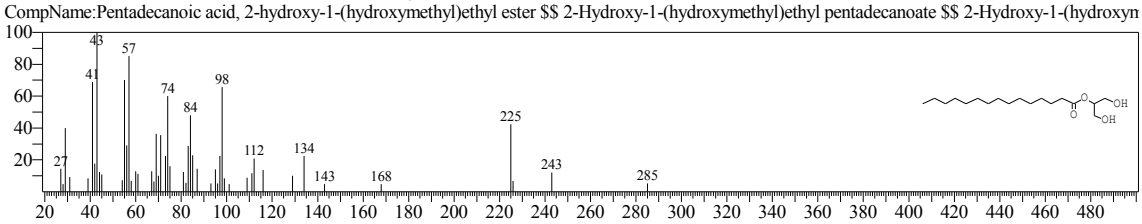
Hit#:1 Entry:436193 Library:Wiley9.lib
SI:93 Formula:C19H38O4 CAS:23470-00-0 MolWeight:330 RetIndex:0
CompName:Hexadecanoic acid, 2-hydroxy-1-(hydroxymethyl)ethyl ester (CAS) $$ 2-Monopalmitin $$ Palmitin, 2-mono- $$ 2-Hexadecanoyl glycerol $$ 2-



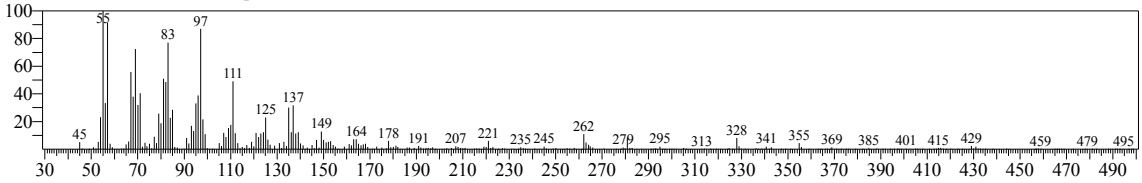
Hit#:2 Entry:436190 Library:Wiley9.lib
SI:86 Formula:C19H38O4 CAS:542-44-9 MolWeight:330 RetIndex:0
CompName:Hexadecanoic acid, 2,3-dihydroxypropyl ester (CAS) $$ 1-Monopalmitin $$ GLYCEROL-1-HEXADECANOATE $$ Palmitin, 1-mono- $$ Gly



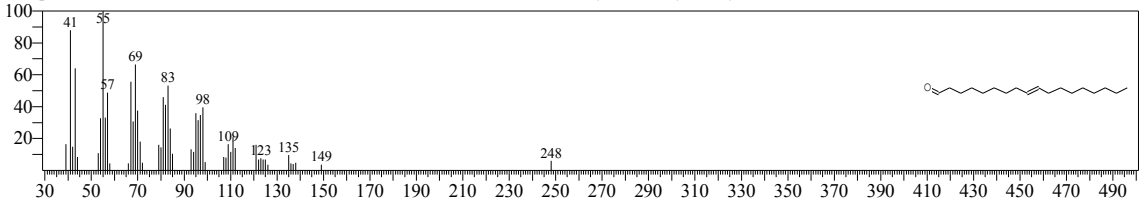
Hit#:3 Entry:408553 Library:Wiley9.lib
SI:85 Formula:C18H36O4 CAS:98863-01-5 MolWeight:316 RetIndex:0
CompName:Pentadecanoic acid, 2-hydroxy-1-(hydroxymethyl)ethyl ester $$ 2-Hydroxy-1-(hydroxymethyl)ethyl pentadecanoate $$ 2-Hydroxy-1-(hydroxym



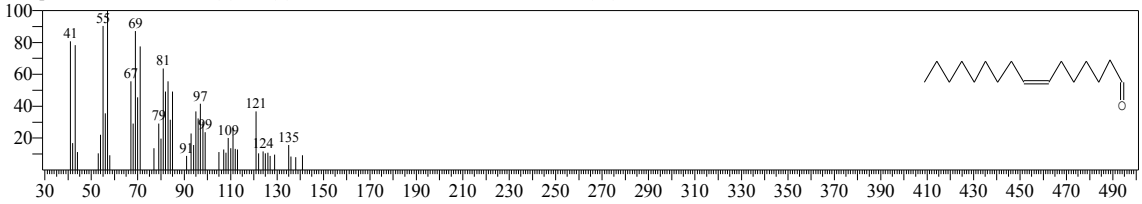
<< Target >>
Line#:30 R.Time:25.425(Scan#:4486) MassPeaks:315
RawMode:Averaged 25.420-25.430(4485-4487) BasePeak:55.00(171911)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



Hit#:1 Entry:298349 Library:Wiley9.lib
SI:87 Formula:C18H34O CAS:2423-10-1 MolWeight:266 RetIndex:0
CompName:9-Octadecenal, (Z)- (CAS) $$ CIS-OCTADEC-9-ENAL $$ Olealdehyde $$ Oleylaldehyde $$ Z-9-Octadecenal $$ cis-9-Octadecenal $$ (9Z)-9-



Hit#:2 Entry:20502 Library:NIST08s.LIB
SI:87 Formula:C16H30O CAS:56797-40-1 MolWeight:238 RetIndex:1808
CompName:7-Hexadecenal, (Z)- $$ (Z)-7-Hexadecenal $$ Z-7-Hexadecenal $$ (7Z)-7-Hexadecenal # $$



Hit#:3 Entry:232676 Library:Wiley9.lib
SI:87 Formula:C16H30O CAS:56797-40-1 MolWeight:238 RetIndex:0
CompName:Z-7-Hexadecenal $$ 7-Hexadecenal, (Z)- $$ (Z)-7-Hexadecenal $$ (7Z)-7-Hexadecenal