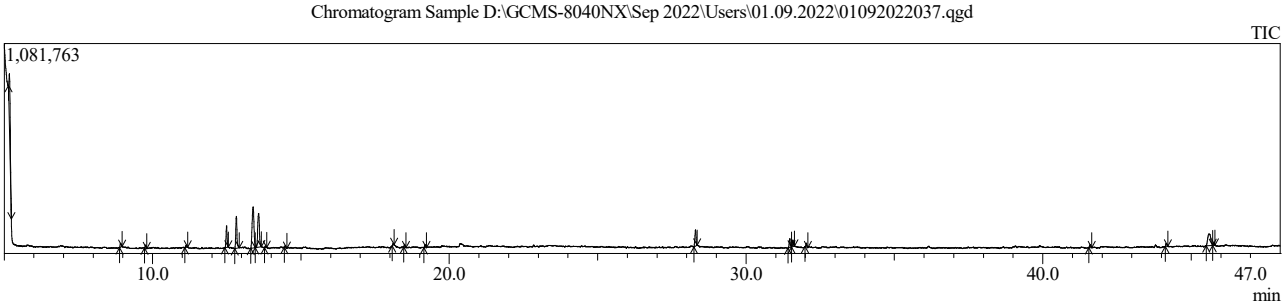
TNAU

Sample Information

Analyzed by	: Admin
Analyzed	: 03-Sep-22 1:14:02 AM
Sample Type	: Unknown
Level #	: 1
Sample Name	: Sample
Sample ID	: 12-3
IS Amount	: [1]=1
Sample Amount	: 1
Dilution Factor	: 1
Vial #	: 13
Injection Volume	: 5.00
Data File	: D:\GCMS-8040NX\Sep 2022\Users\01.09.2022\01092022037.qgd
Org Data File	: D:\GCMS-8040NX\Sep 2022\Users\01.09.2022\01092022037.qgd
Method File	: D:\GCMS-8040NX\Sep 2022\Method\Scan_methodfile.qgm
Org Method File	: D:\GCMS-8040NX\Sep 2022\Method\Scan_methodfile.qgm
Report File	:
Tuning File	: D:\GCMS-8040NX\Aug 2022\Tuning Data\TUNING 25082022.qgt
Modified by	: Admin
Modified	: 05-Sep-22 11:13:16 AM

Chromatogram Sample D:\GCMS-8040NX\Sep 2022\Users\01.09.2022\01092022037.qgd



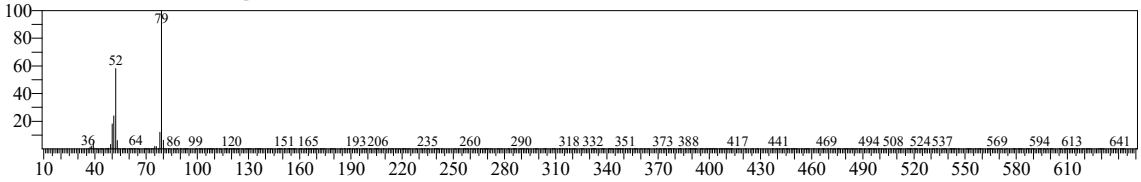
Peak Report TIC

Peak#	R.Time	Area	Area%	Height	Height%	A/H	Similarity	Name
1	5.174	708428	16.34	245702	18.96	2.88	98	Pyridine
2	8.921	54892	1.27	19665	1.52	2.79	85	1-Butanol, 3-methyl-, acetate
3	9.768	39857	0.92	17144	1.32	2.32	94	Pentasiloxane, dodecamethyl-
4	11.134	34914	0.81	12404	0.96	2.81	39	Methyl arachidonate
5	12.492	277971	6.41	110111	8.50	2.52	73	2,5-Cyclohexadiene-1,4-dione, dioxime
6	12.822	456019	10.52	165724	12.79	2.75	73	2,5-Cyclohexadiene-1,4-dione, dioxime
7	13.388	911565	21.02	210680	16.26	4.33	53	Methyl cis-13,16-Docosadienate
8	13.573	690829	15.93	173708	13.40	3.98	53	Methyl cis-13,16-Docosadienate
9	13.800	19696	0.45	8144	0.63	2.42	47	Boric acid-3TMS
10	14.467	27175	0.63	11060	0.85	2.46	72	2,3-Dimethyl-para-anisaldehyde
11	18.112	19024	0.44	9582	0.74	1.99	89	.beta.-D-Glucopyranose, 1,6-anhydro-
12	18.495	33440	0.77	15000	1.16	2.23	91	2,4-Di-tert-butylphenol
13	19.192	38948	0.90	14354	1.11	2.71	86	2,4-Di-tert-butylphenoxytrimethylsilane
14	28.296	218926	5.05	83168	6.42	2.63	95	n-Hexadecanoic acid
15	31.473	123644	2.85	44633	3.44	2.77	94	9,12-Octadecadienoic acid (Z,Z)-
16	31.580	100587	2.32	33681	2.60	2.99	88	cis-9-Hexadecenal
17	32.030	42562	0.98	19485	1.50	2.18	93	Octadecanoic acid
18	41.602	43544	1.00	15435	1.19	2.82	83	Tetracosamethyl-cyclododecasiloxane
19	44.157	40205	0.93	14712	1.14	2.73	88	13-Docosenamide, (Z)-
20	45.601	426512	9.84	60247	4.65	7.08	91	Diosgenin
21	45.767	27012	0.62	11223	0.87	2.41	78	Tetracosamethyl-cyclododecasiloxane
		4335750	100.00	1295862	100.00			

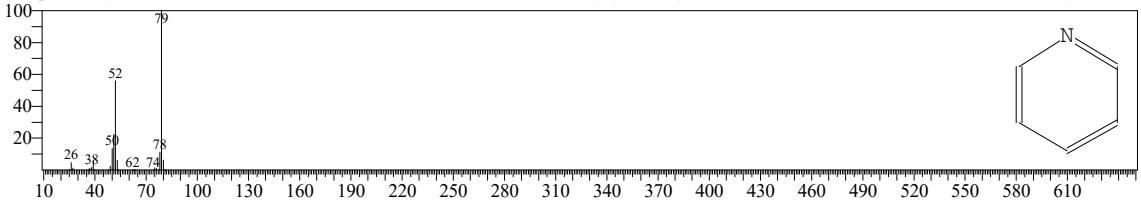
TNAU

Library

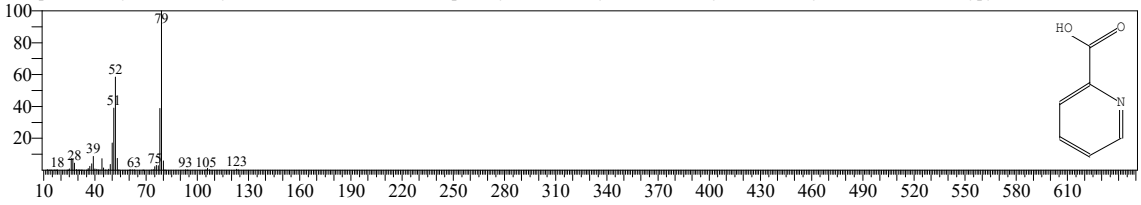
<< Target >>
Line#:1 R.Time:5.175(Scan#:36) MassPeaks:335
RawMode:Averaged 5.170-5.180(35-37) BasePeak:79.05(101115)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



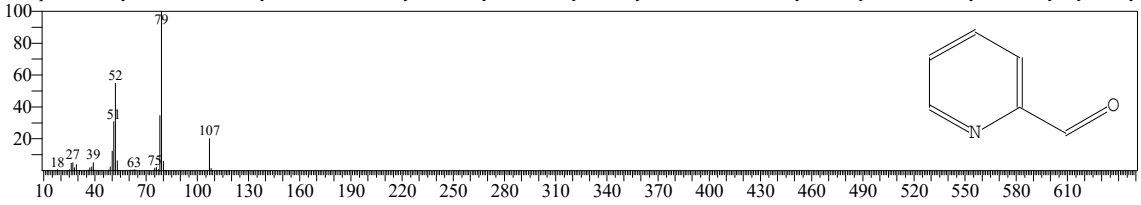
Hit#:1 Entry:498 Library:NIST20M1.lib
SI:98 Formula:C5H5N CAS:110-86-1 MolWeight:79 RetIndex:674
CompName:Pyridine $$ Azabenzene $$ Azine $$ NCI-C55301 $$ Piridina $$ Pirydyna $$ Pyridin $$ Rcra waste number U196 $$ UN 1282 $$ Pyr $$ CP



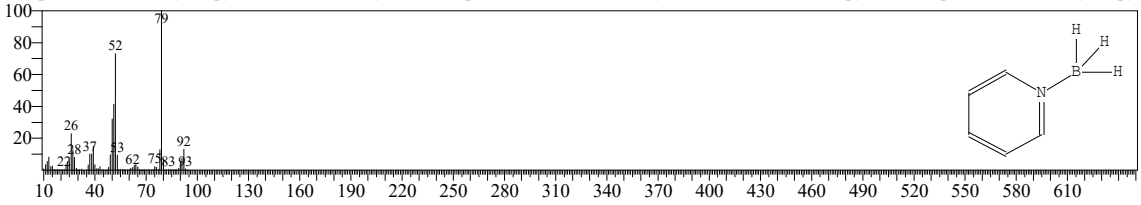
Hit#:2 Entry:6149 Library:NIST20M1.lib
SI:92 Formula:C6H5NO2 CAS:98-98-6 MolWeight:123 RetIndex:1144
CompName:2-Pyridinecarboxylic acid $$ Picolinic acid $$ .alpha.-Pyridinecarboxylic acid $$ o-Pyridinecarboxylic acid $$ 2-Carboxypyridine $$ 2-Picolini



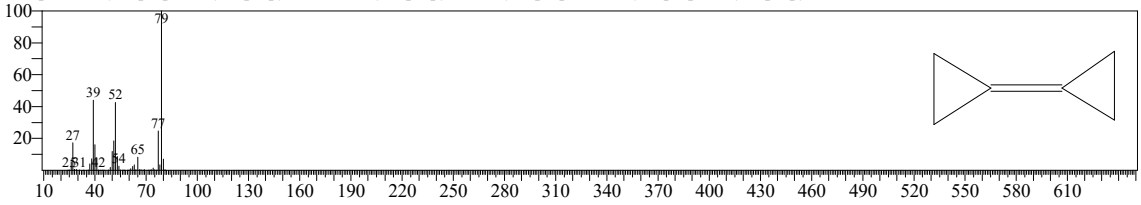
Hit#:3 Entry:2869 Library:NIST20M1.lib
SI:91 Formula:C6H5NO CAS:1121-60-4 MolWeight:107 RetIndex:976
CompName:2-Pyridinecarboxaldehyde $$ Picolinaldehyde $$ 2-Pyridinealdehyde $$ Pyridine-2-carboxaldehyde $$ Pyridine-2-aldehyde $$ 2-Pyridylaldehy



Hit#:4 Entry:1275 Library:NIST20M1.lib
SI:86 Formula:C5H8BN CAS:110-51-0 MolWeight:93 RetIndex:0
CompName:Boron, trihydro(pyridine)-, (T-4)- $$ Pyridine, compd. with borane (1:1) $$ Pyridine borane $$ Borane-pyridine complex $$ Boron, trihydro(py

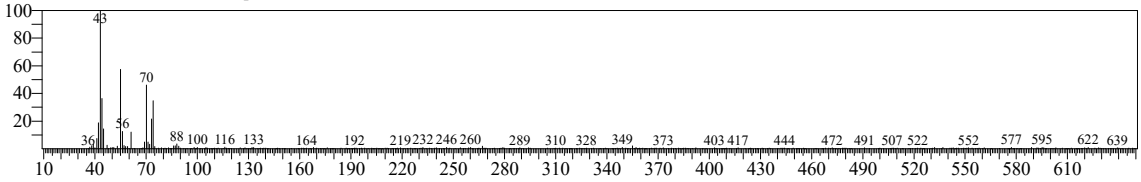


Hit#:5 Entry:539 Library:NIST20M1.lib
SI:85 Formula:C6H8 CAS:27567-82-4 MolWeight:80 RetIndex:661
CompName:Cyclopropane, cyclopropylidene- $$ Cyclopropylidene cyclopropane $$ Cyclopropane cyclopropylidene-

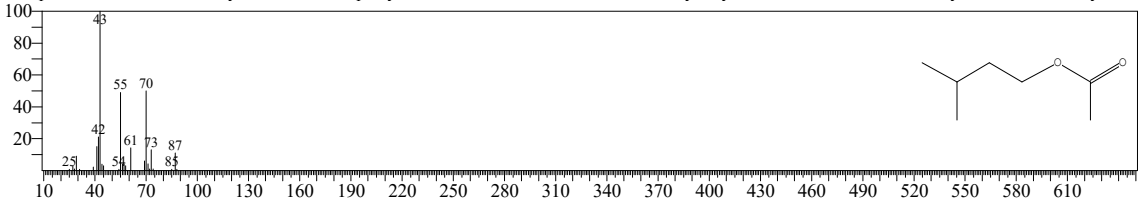


TNAU

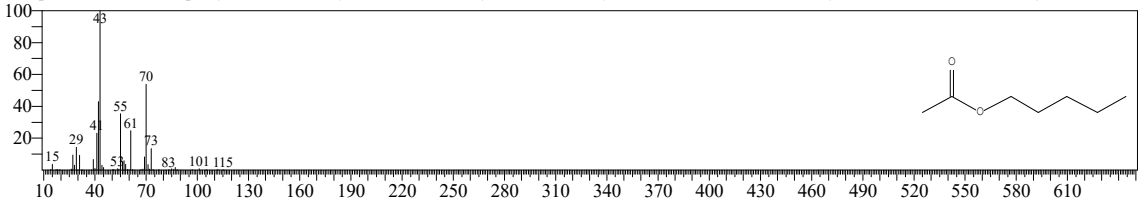
<< Target >>
Line#:2 R.Time:8.920(Scan#:785) MassPeaks:371
RawMode:Averaged 8.915-8.925(784-786) BasePeak:43.05(4273)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



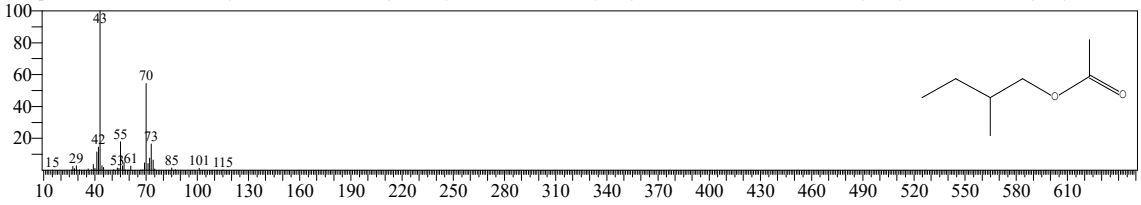
Hit#:1 Entry:6783 Library:NIST20R.lib
SI:85 Formula:C7H14O2 CAS:123-92-2 MolWeight:130 RetIndex:820
CompName:1-Butanol, 3-methyl-, acetate $$ Isopentyl alcohol, acetate $$ Acetic acid, 3-methylbutyl ester $$ Banana oil $$ Isoamyl acetate $$ Isoamyl etha



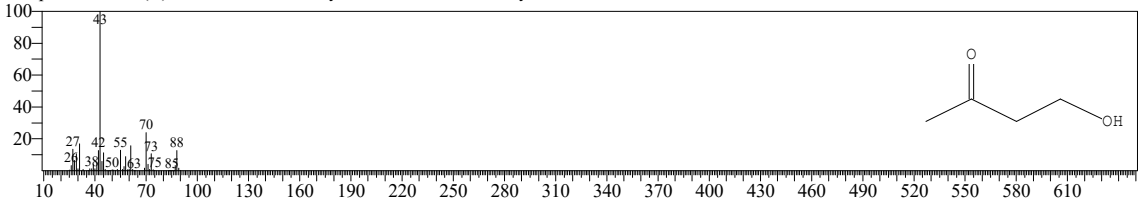
Hit#:2 Entry:8685 Library:NIST20M1.lib
SI:82 Formula:C7H14O2 CAS:628-63-7 MolWeight:130 RetIndex:884
CompName:Acetic acid, pentyl ester $$ n-Amyl acetate $$ n-Pentyl acetate $$ Amyl acetate $$ Birnenoel $$ Pentyl acetate $$ Acetic acid, amyl ester $$ An



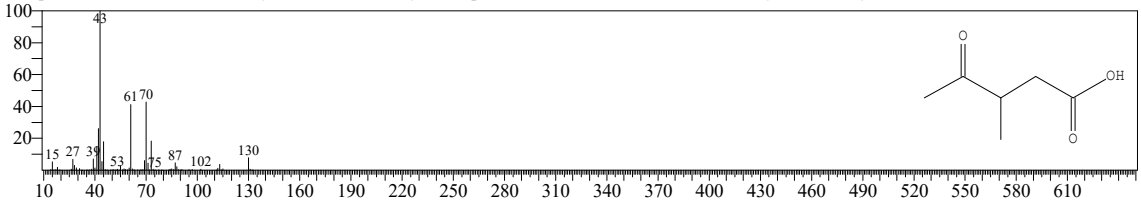
Hit#:3 Entry:6819 Library:NIST20R.lib
SI:82 Formula:C7H14O2 CAS:624-41-9 MolWeight:130 RetIndex:820
CompName:1-Butanol, 2-methyl-, acetate $$ 2-Methyl-1-butyl acetate $$ 2-Methylbutyl acetate $$ Acetic acid 2-methylbutyl ester $$ 2-Methybutyl acetate



Hit#:4 Entry:1220 Library:NIST20R.lib
SI:81 Formula:C4H8O2 CAS:590-90-9 MolWeight:88 RetIndex:798
CompName:CH3C(O)CH2CH2OH $$ Methylolacetone $$ Monomethylolacetone $$ 3-Ketobutan-1-ol $$ 3-Oxo-1-butanol $$ 3-Oxobutanol $$ 4-Butanol-

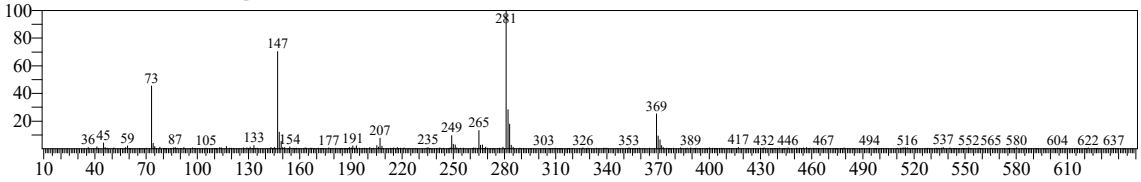


Hit#:5 Entry:8539 Library:NIST20M1.lib
SI:79 Formula:C6H10O3 CAS:6628-79-1 MolWeight:130 RetIndex:1046
CompName:Pentanoic acid, 3-methyl-4-oxo- $$ 3-Methyl-4-oxopentanoic acid $$ Levulinic acid, 3-methyl $$ 3-Methyllevulinic acid

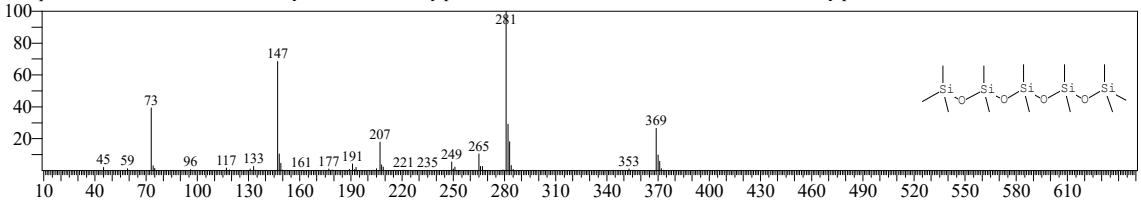


TNAU

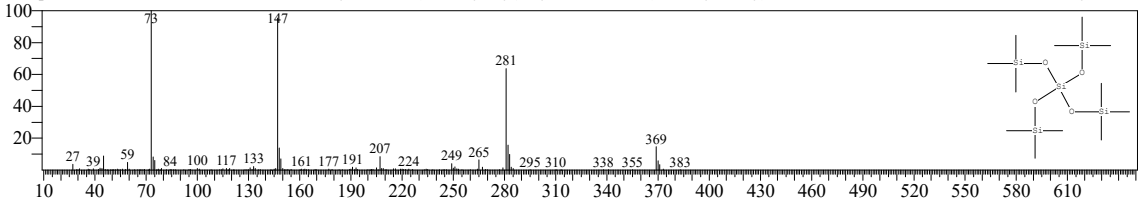
<< Target >>
Line#:3 R.Time:9.770(Scan#:955) MassPeaks:330
RawMode:Averaged 9.765-9.775(954-956) BasePeak:281.00(4059)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



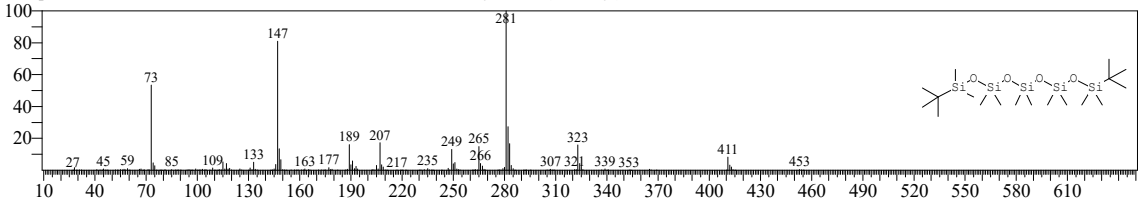
Hit#:1 Entry:40975 Library:NIST20R.lib
SI:94 Formula:C12H36O4Si5 CAS:141-63-9 MolWeight:384 RetIndex:1068
CompName:Pentasiloxane, dodecamethyl- $$ Dodecamethylpentasiloxane $$ 1,1,1,3,3,5,5,7,7,9,9,9-Dodecamethylpentasiloxane #



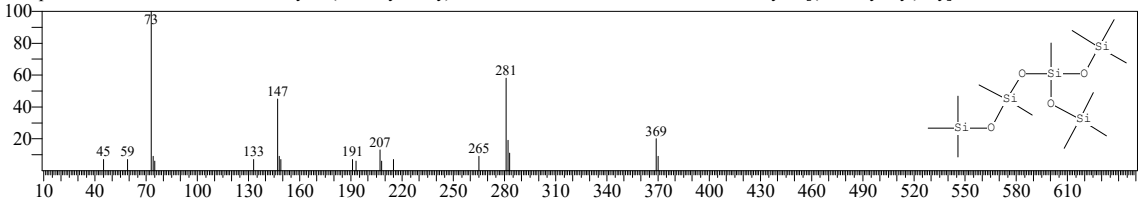
Hit#:2 Entry:249272 Library:NIST20M1.lib
SI:84 Formula:C12H36O4Si5 CAS:3555-47-3 MolWeight:384 RetIndex:1068
CompName:Trisiloxane, 1,1,1,5,5,5-hexamethyl-3,3-bis[(trimethylsilyl)oxy]- $$ Tetrakis(trimethylsiloxy)silane $$ Trisiloxane, 1,1,1,5,5,5-hexamethyl-3,3-b



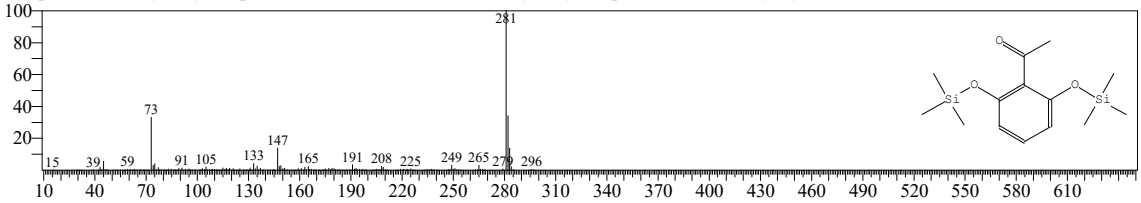
Hit#:3 Entry:27848 Library:NIST20M2.lib
SI:81 Formula:C18H48O4Si5 CAS:0-00-0 MolWeight:468 RetIndex:1495
CompName:1,3,5,7,9-Pentasiloxane, 1,1,3,3,5,5,7,7,9,9-decamethyl-1,9-di(tert.butyl)-



Hit#:4 Entry:249271 Library:NIST20M1.lib
SI:79 Formula:C12H36O4Si5 CAS:38146-99-5 MolWeight:384 RetIndex:1068
CompName:1,1,1,3,5,5,7,7,7-Nonamethyl-3-(trimethylsiloxy)tetrasiloxane $$ 1,1,1,3,3,5,7,7,7-Nonamethyl-5-[(trimethylsilyl)oxy]tetrasiloxane #

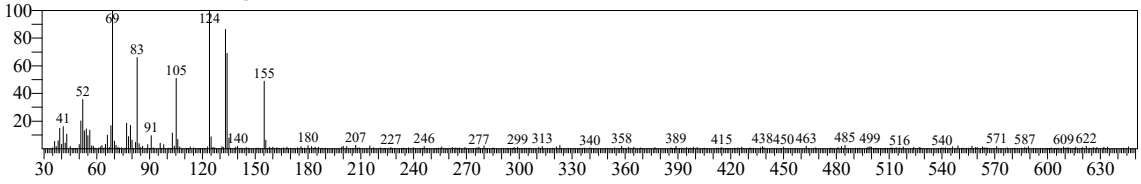


Hit#:5 Entry:158097 Library:NIST20M1.lib
SI:75 Formula:C14H24O3Si2 CAS:0-00-0 MolWeight:296 RetIndex:1625
CompName:2,6-Dihydroxyacetophenone, 2TMS derivative $$ 2',6'-Dihydroxyacetophenone, bis(trimethylsilyl) ether

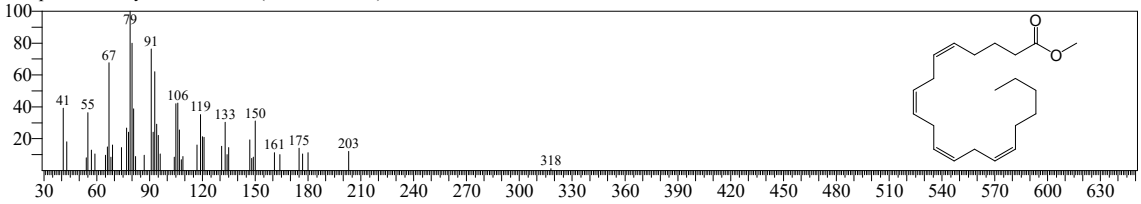


TNAU

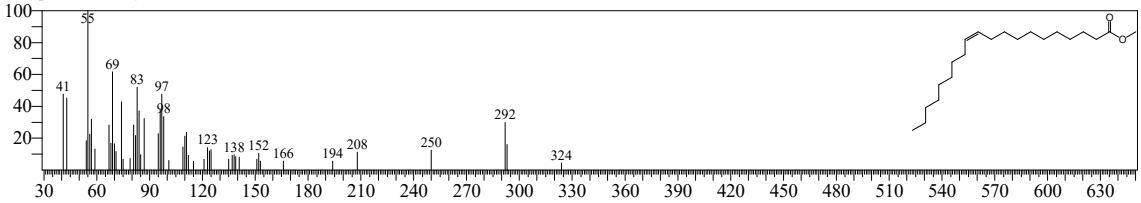
<< Target >>
Line#:4 R.Time:11.135(Scan#:1228) MassPeaks:278
RawMode:Averaged 11.130-11.140(1227-1229) BasePeak:69.00(1635)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



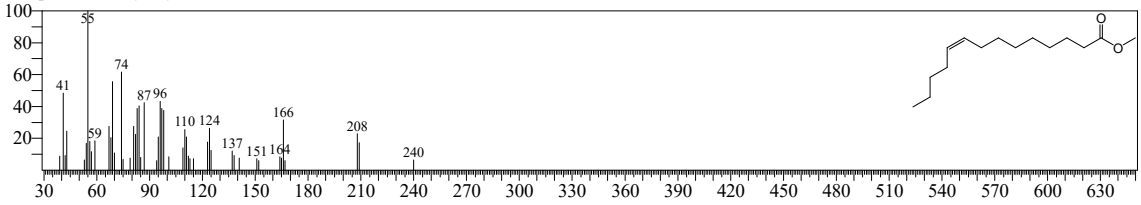
Hit#:1 Entry:33 Library:FA_ME_SP2560_EI_V3.lib
SI:39 Formula:C21H34O2 CAS:506-32-1 MolWeight:318 RetIndex:3109
CompName:Methyl arachidonate ; (5Z,8Z,11Z,14Z)-icosa-5,8,11,14-tetraenoic acid



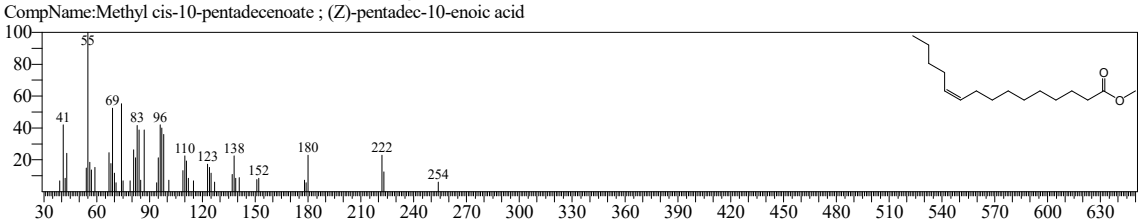
Hit#:2 Entry:24 Library:FA_ME_SP2560_EI_V3.lib
SI:39 Formula:C21H40O2 CAS:5561-99-9 MolWeight:324 RetIndex:2874
CompName:Methyl cis-11-icosenoate ; ; Icos-11-enoic acid



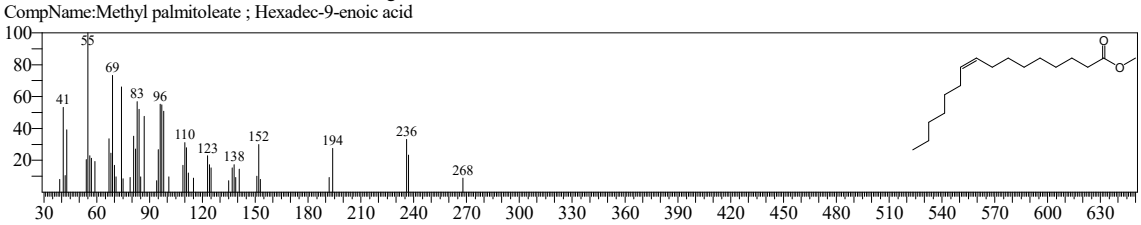
Hit#:3 Entry:9 Library:FA_ME_SP2560_EI_V3.lib
SI:39 Formula:C15H28O2 CAS:544-64-9 MolWeight:240 RetIndex:2283
CompName:Methyl myristoleate ; ; (Z)-tetradec-9-enoic acid



Hit#:4 Entry:11 Library:FA_ME_SP2560_EI_V3.lib
SI:39 Formula:C16H30O2 CAS:84743-29-3 MolWeight:254 RetIndex:2388
CompName:Methyl cis-10-pentadecenoate ; (Z)-pentadec-10-enoic acid

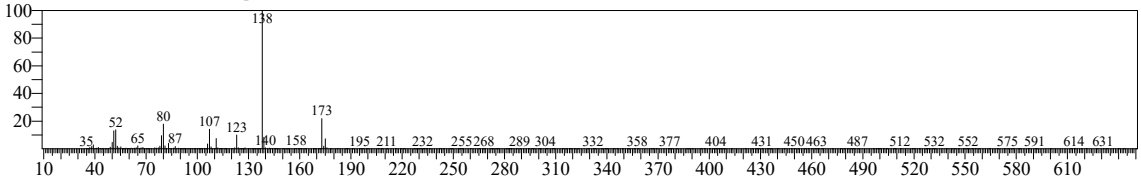


Hit#:5 Entry:13 Library:FA_ME_SP2560_EI_V3.lib
SI:39 Formula:C17H32O2 CAS:373-49-9 MolWeight:268 RetIndex:2478
CompName:Methyl palmitoleate ; Hexadec-9-enoic acid

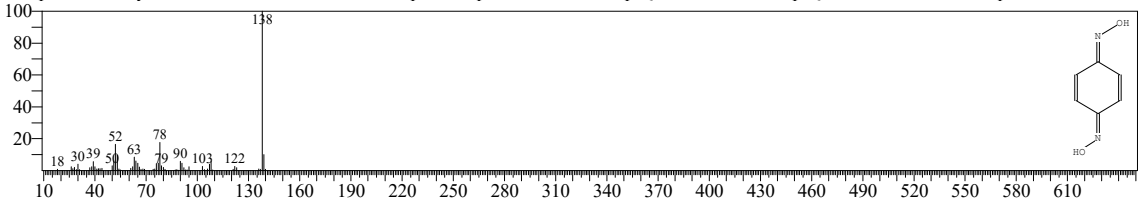


TNAU

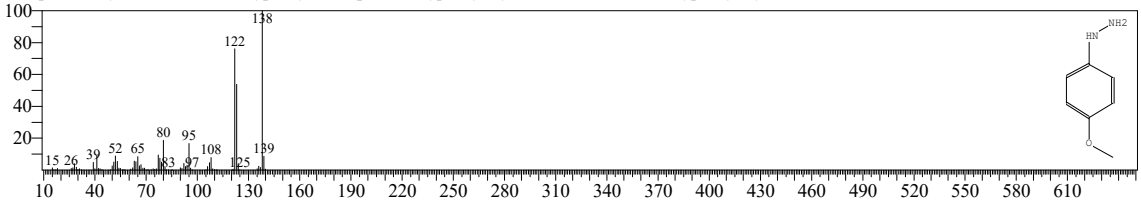
<< Target >>
Line#:5 R.Time:12.490(Scan#:1499) MassPeaks:297
RawMode:Averaged 12.485-12.495(1498-1500) BasePeak:138.05(39351)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



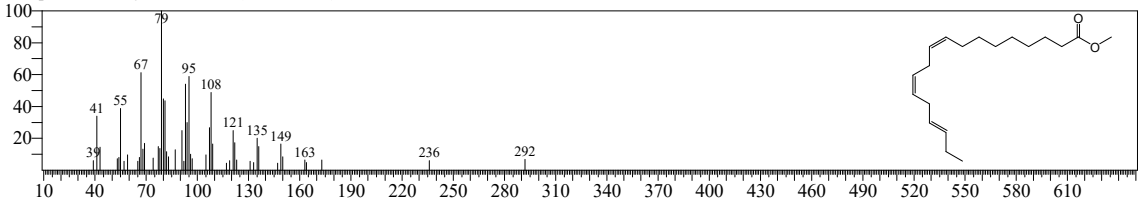
Hit#:1 Entry:8360 Library:NIST20R.lib
SI:73 Formula:C6H6N2O2 CAS:105-11-3 MolWeight:138 RetIndex:1349
CompName:2,5-Cyclohexadiene-1,4-dione, dioxime $$ p-Benzoquinone, dioxime $$ p-Quinone dioxime $$ p-Quinone oxime $$ Benzoquinone dioxime $$



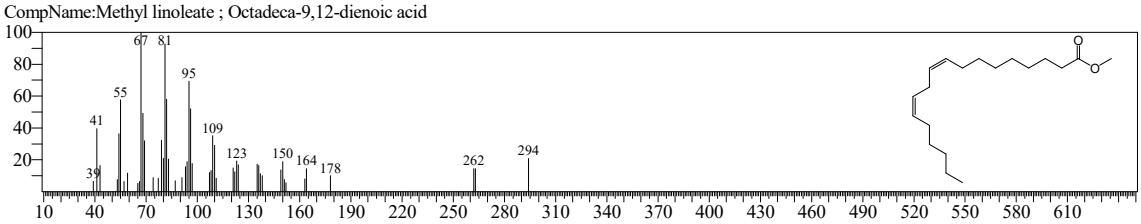
Hit#:2 Entry:11222 Library:NIST20M1.lib
SI:71 Formula:C7H10N2O CAS:3471-32-7 MolWeight:138 RetIndex:1325
CompName:Hydrazine, (4-methoxyphenyl)- $$ (p-Methoxyphenyl)hydrazine $$ 1-(4-Methoxyphenyl)hydrazine #



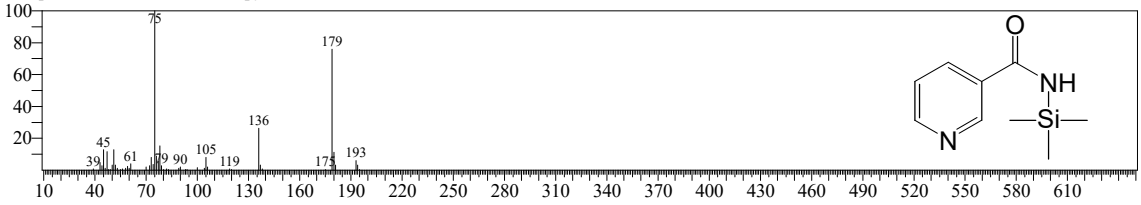
Hit#:3 Entry:25 Library:FA_ME_SP2560_EI_V3.lib
SI:34 Formula:C19H32O2 CAS:463-40-1 MolWeight:292 RetIndex:2892
CompName:Methyl linolenate ; (9Z,12Z,15Z)-octadeca-9,12,15-trienoic acid



Hit#:4 Entry:21 Library:FA_ME_SP2560_EI_V3.lib
SI:34 Formula:C19H34O2 CAS:60-33-3 MolWeight:294 RetIndex:2775
CompName:Methyl linoleate ; Octadeca-9,12-dienoic acid

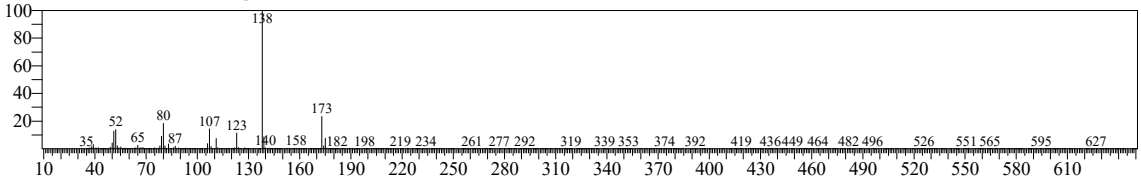


Hit#:5 Entry:137 Library:OA_TMS_DB5_67min_V3.lib
SI:33 Formula:C9H14N2OSi CAS:98-92-0 MolWeight:194 RetIndex:1486
CompName:Niacinamide-TMS ; pyridine-3-carboxamide

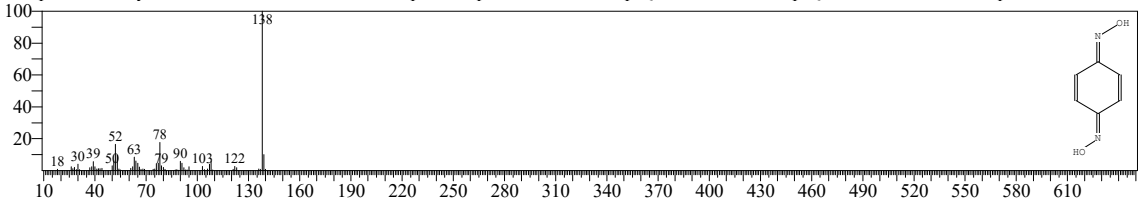


TNAU

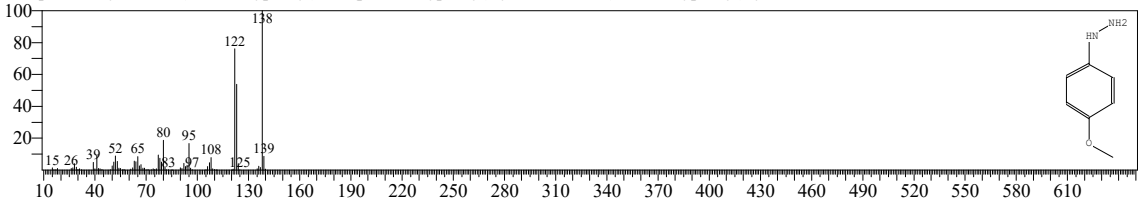
<< Target >>
Line#:6 R.Time:12.820(Scan#:1565) MassPeaks:392
RawMode:Averaged 12.815-12.825(1564-1566) BasePeak:138.05(56730)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



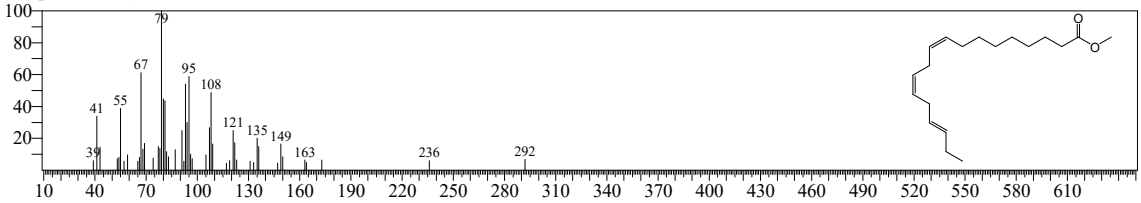
Hit#:1 Entry:8360 Library:NIST20R.lib
SI:73 Formula:C6H6N2O2 CAS:105-11-3 MolWeight:138 RetIndex:1349
CompName:2,5-Cyclohexadiene-1,4-dione, dioxime $$ p-Benzoquinone, dioxime $$ p-Quinone dioxime $$ p-Quinone oxime $$ Benzoquinone dioxime $$



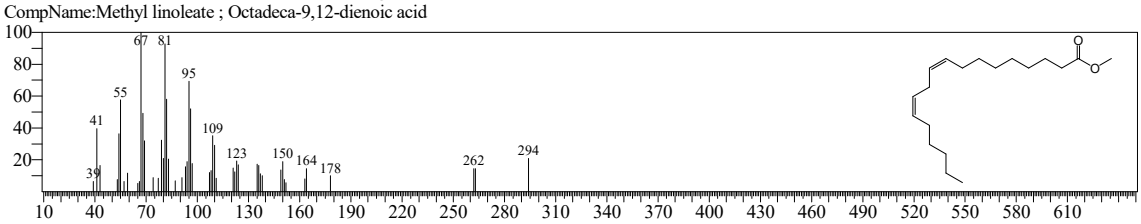
Hit#:2 Entry:11222 Library:NIST20M1.lib
SI:71 Formula:C7H10N2O CAS:3471-32-7 MolWeight:138 RetIndex:1325
CompName:Hydrazine, (4-methoxyphenyl)- $$ (p-Methoxyphenyl)hydrazine $$ 1-(4-Methoxyphenyl)hydrazine #



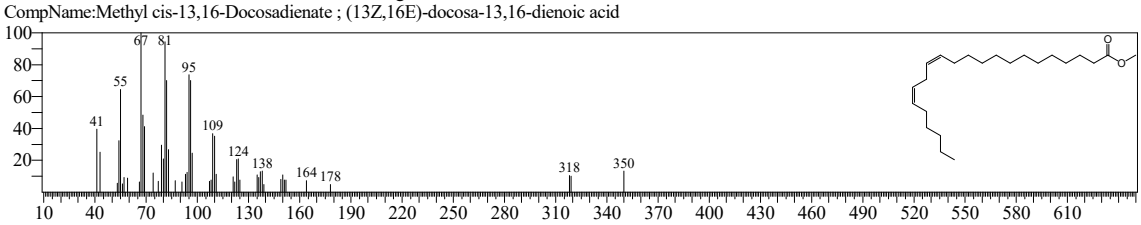
Hit#:3 Entry:25 Library:FA_ME_SP2560_EI_V3.lib
SI:35 Formula:C19H32O2 CAS:463-40-1 MolWeight:292 RetIndex:2892
CompName:Methyl linolenate ; (9Z,12Z,15Z)-octadeca-9,12,15-trienoic acid



Hit#:4 Entry:21 Library:FA_ME_SP2560_EI_V3.lib
SI:34 Formula:C19H34O2 CAS:60-33-3 MolWeight:294 RetIndex:2775
CompName:Methyl linoleate ; Octadeca-9,12-dienoic acid

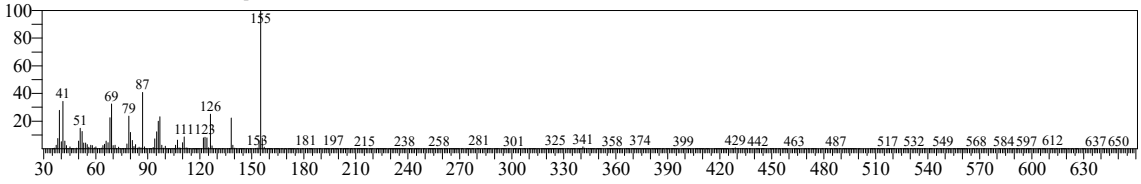


Hit#:5 Entry:34 Library:FA_ME_SP2560_EI_V3.lib
SI:34 Formula:C23H42O2 CAS:7370-49-2 MolWeight:350 RetIndex:3169
CompName:Methyl cis-13,16-Docosadienate ; (13Z,16E)-docosa-13,16-dienoic acid

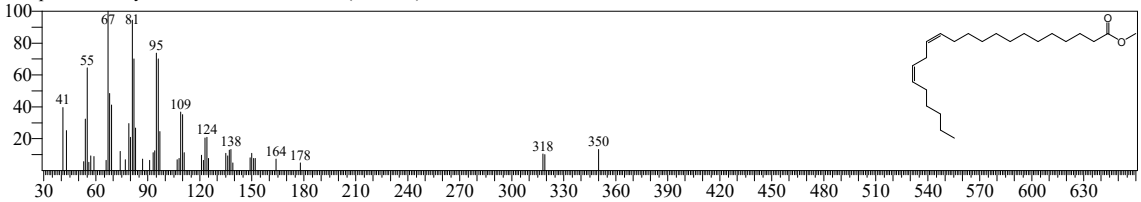


TNAU

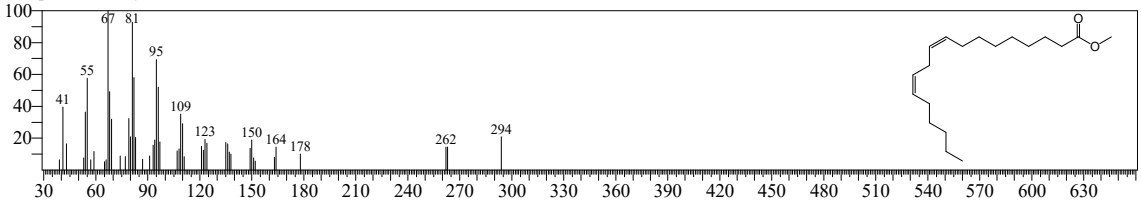
<< Target >>
Line#:7 R.Time:13.390(Scan#:1679) MassPeaks:409
RawMode:Averaged 13.385-13.395(1678-1680) BasePeak:155.05(34056)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



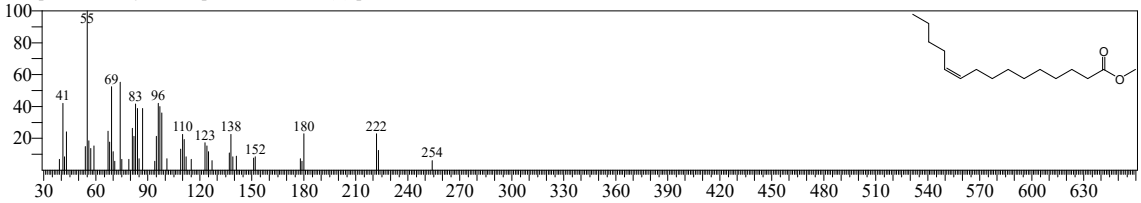
Hit#:1 Entry:34 Library:FA_ME_SP2560_EI_V3.lib
SI:53 Formula:C23H42O2 CAS:7370-49-2 MolWeight:350 RetIndex:3169
CompName:Methyl cis-13,16-Docosadienate ; (13Z,16E)-docosa-13,16-dienoic acid



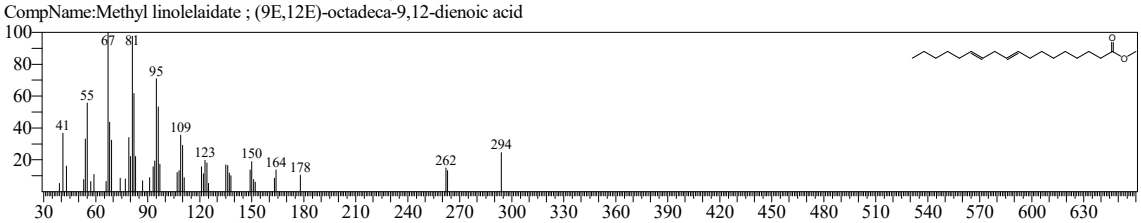
Hit#:2 Entry:21 Library:FA_ME_SP2560_EI_V3.lib
SI:52 Formula:C19H34O2 CAS:60-33-3 MolWeight:294 RetIndex:2775
CompName:Methyl linoleate ; Octadeca-9,12-dienoic acid



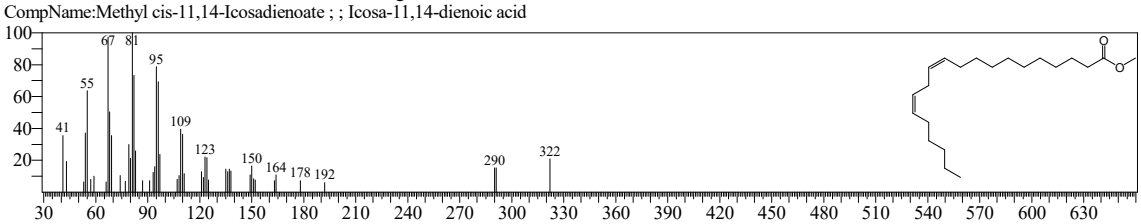
Hit#:3 Entry:11 Library:FA_ME_SP2560_EI_V3.lib
SI:52 Formula:C16H30O2 CAS:84743-29-3 MolWeight:254 RetIndex:2388
CompName:Methyl cis-10-pentadecenoate ; (Z)-pentadec-10-enoic acid



Hit#:4 Entry:20 Library:FA_ME_SP2560_EI_V3.lib
SI:52 Formula:C19H34O2 CAS:506-21-8 MolWeight:294 RetIndex:2727
CompName:Methyl linolelaidate ; (9E,12E)-octadeca-9,12-dienoic acid

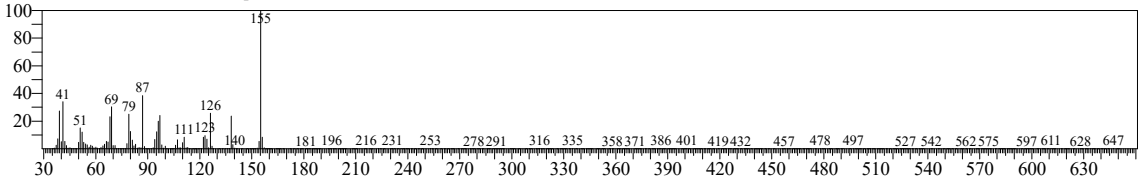


Hit#:5 Entry:27 Library:FA_ME_SP2560_EI_V3.lib
SI:51 Formula:C21H38O2 CAS:2091-39-6 MolWeight:322 RetIndex:2973
CompName:Methyl cis-11,14-Icosadienoate ; ; Icosa-11,14-dienoic acid

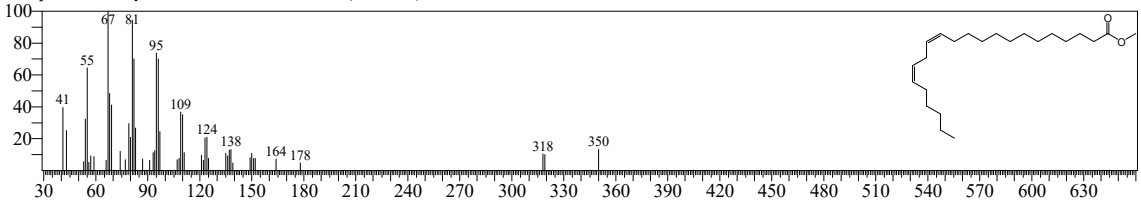


TNAU

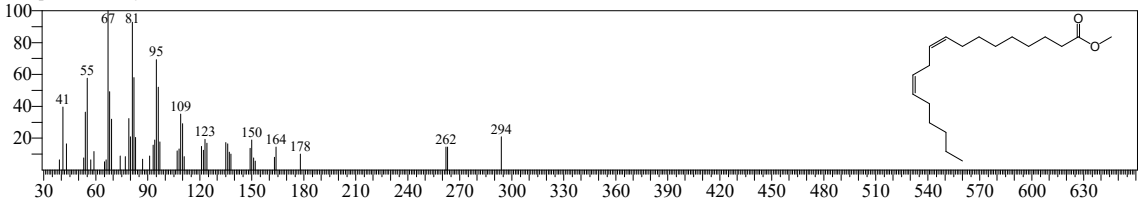
<< Target >>
Line#:8 R.Time:13.575(Scan#:1716) MassPeaks:407
RawMode:Averaged 13.570-13.580(1715-1717) BasePeak:155.05(28048)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



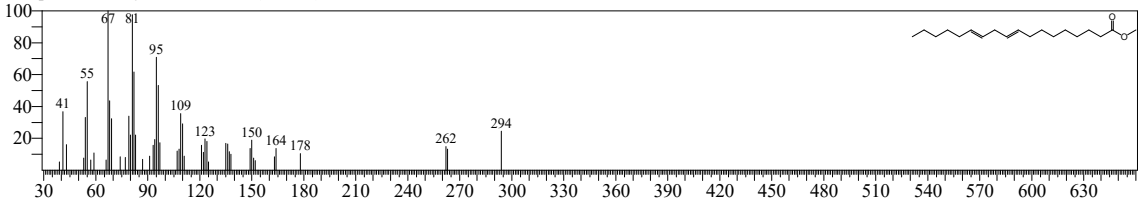
Hit#:1 Entry:34 Library:FA_ME_SP2560_EI_V3.lib
SI:53 Formula:C23H42O2 CAS:7370-49-2 MolWeight:350 RetIndex:3169
CompName:Methyl cis-13,16-Docosadienate ; (13Z,16E)-docosa-13,16-dienoic acid



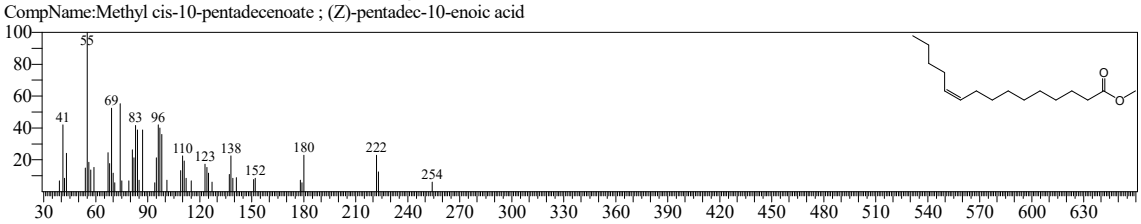
Hit#:2 Entry:21 Library:FA_ME_SP2560_EI_V3.lib
SI:52 Formula:C19H34O2 CAS:60-33-3 MolWeight:294 RetIndex:2775
CompName:Methyl linoleate ; Octadeca-9,12-dienoic acid



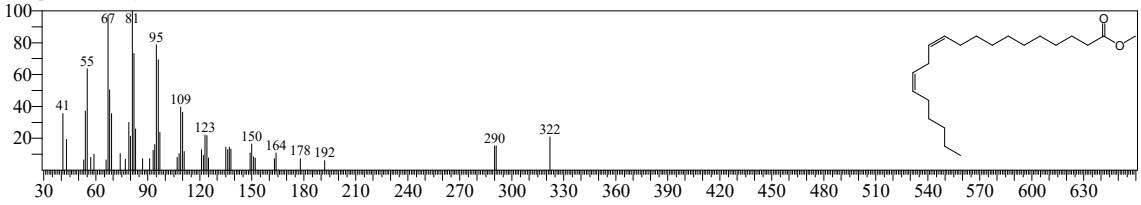
Hit#:3 Entry:20 Library:FA_ME_SP2560_EI_V3.lib
SI:52 Formula:C19H34O2 CAS:506-21-8 MolWeight:294 RetIndex:2727
CompName:Methyl linolelaidate ; (9E,12E)-octadeca-9,12-dienoic acid



Hit#:4 Entry:11 Library:FA_ME_SP2560_EI_V3.lib
SI:52 Formula:C16H30O2 CAS:84743-29-3 MolWeight:254 RetIndex:2388
CompName:Methyl cis-10-pentadecenoate ; (Z)-pentadec-10-enoic acid

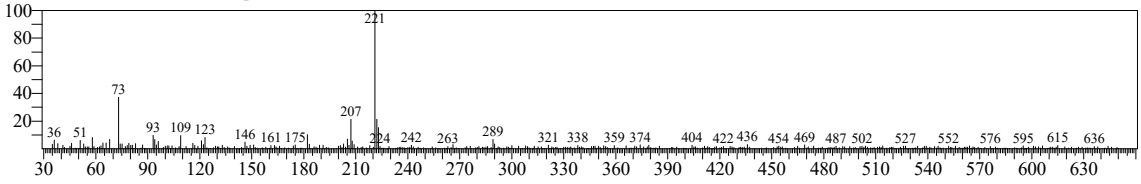


Hit#:5 Entry:27 Library:FA_ME_SP2560_EI_V3.lib
SI:51 Formula:C21H38O2 CAS:2091-39-6 MolWeight:322 RetIndex:2973
CompName:Methyl cis-11,14-Icosadienoate ; ; Icosa-11,14-dienoic acid

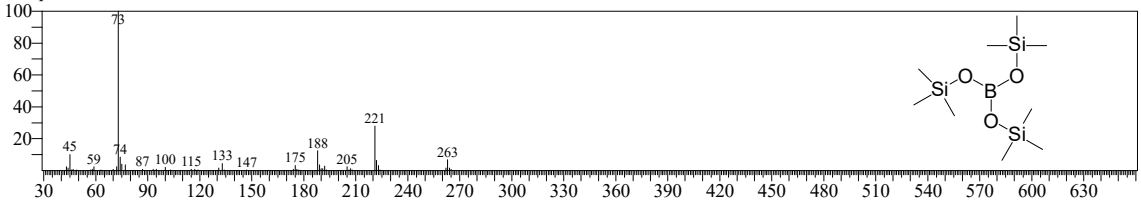


TNAU

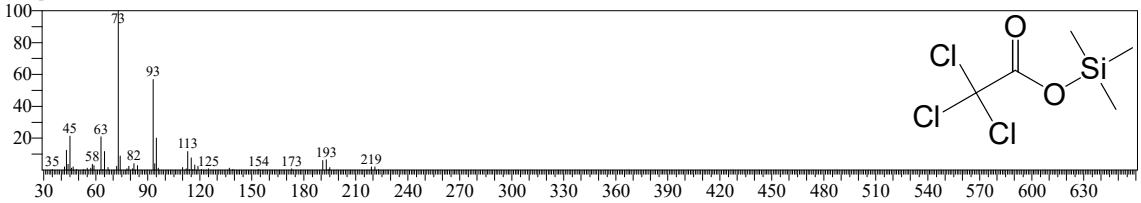
<< Target >>
Line#:9 R.Time:13.800(Scan#:1761) MassPeaks:397
RawMode:Averaged 13.795-13.805(1760-1762) BasePeak:221.05(1429)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



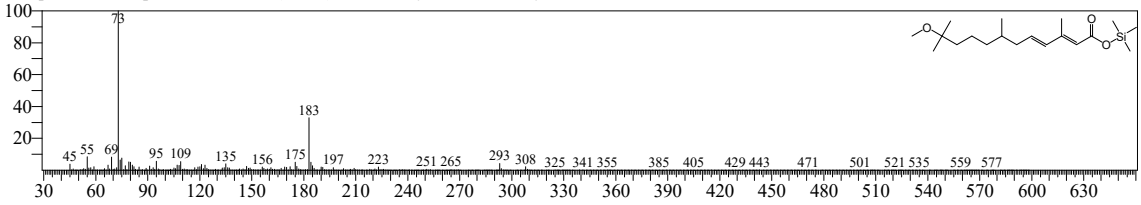
Hit#:1 Entry:3 Library:OA_TMS_DB5_67min_V3.lib
SI:47 Formula:C9H27BO3Si3 CAS:10043-35-3 MolWeight:278 RetIndex:992
CompName:Boric acid-3TMS ; boric acid



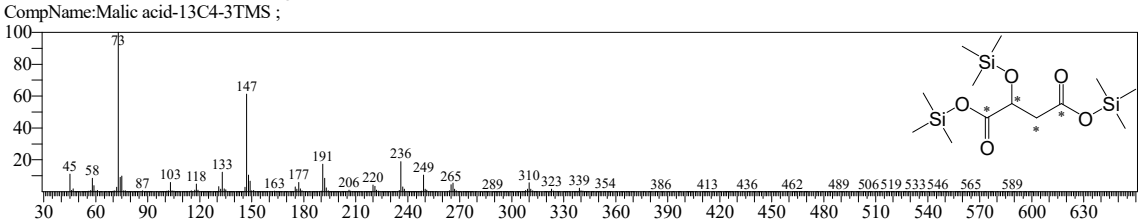
Hit#:2 Entry:6 Library:OA_TMS_DB5_67min_V3.lib
SI:39 Formula:C5H9Cl3O2Si CAS:76-03-9 MolWeight:234 RetIndex:1059
CompName:Trichloroacetic acid-TMS ; 2,2,2-trichloroacetic acid



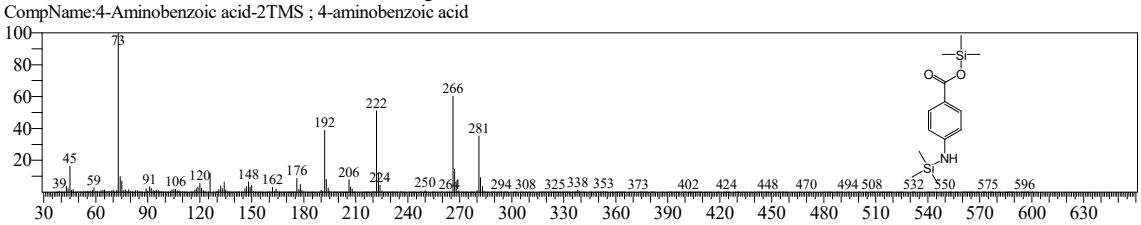
Hit#:3 Entry:466 Library:OA_TMS_DB5_67min_V3.lib
SI:33 Formula:C19H36O3Si CAS:53092-52-7 MolWeight:340 RetIndex:2133
CompName:Methoprene acid-TMS ; (2E,4E)-11-methoxy-3,7,11-trimethyldodeca-2,4-dienoic acid



Hit#:4 Entry:143 Library:OA_TMS_DB5_67min_V3.lib
SI:33 Formula: CAS:0-00-0 MolWeight:354 RetIndex:1495
CompName:Malic acid-13C4-3TMS ;

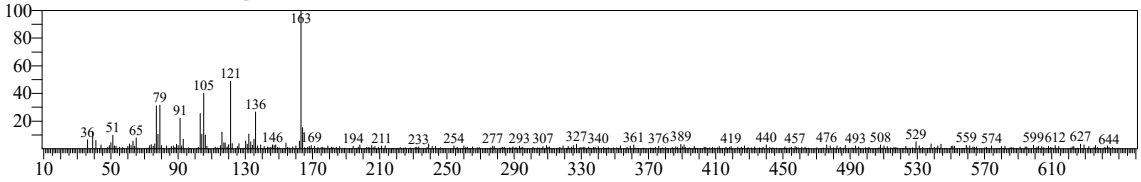


Hit#:5 Entry:328 Library:OA_TMS_DB5_67min_V3.lib
SI:32 Formula:C13H23NO2Si2 CAS:150-13-0 MolWeight:281 RetIndex:1845
CompName:4-Aminobenzoic acid-2TMS ; 4-aminobenzoic acid

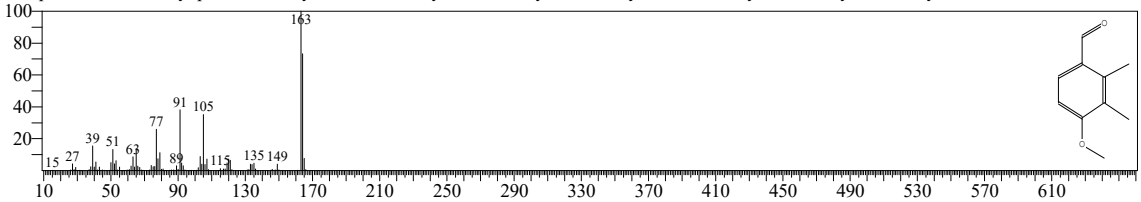


TNAU

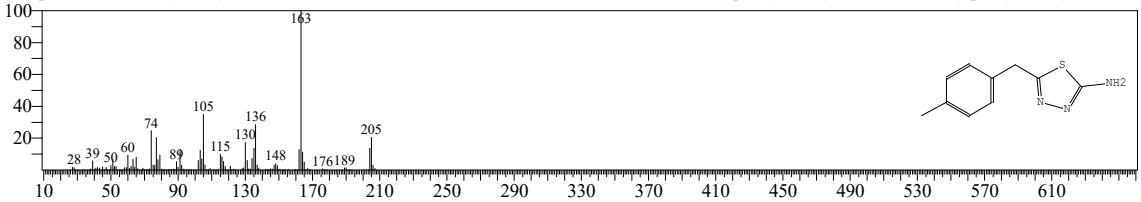
<< Target >>
Line#:10 R.Time:14.465(Scan#:1894) MassPeaks:369
RawMode:Averaged 14.460-14.470(1893-1895) BasePeak:163.10(1523)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



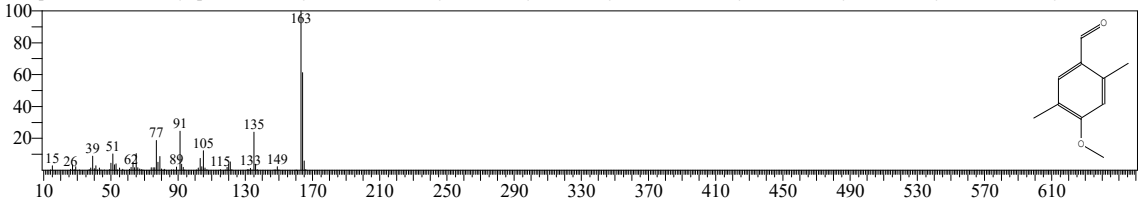
Hit#:1 Entry:25351 Library:NIST20M1.lib
SI:72 Formula:C10H12O2 CAS:38998-17-3 MolWeight:164 RetIndex:1398
CompName:2,3-Dimethyl-para-anisaldehyde $$ Benzaldehyde, 4-methoxy-2,3-dimethyl- $$ 4-Methoxy-2,3-dimethylbenzaldehyde #



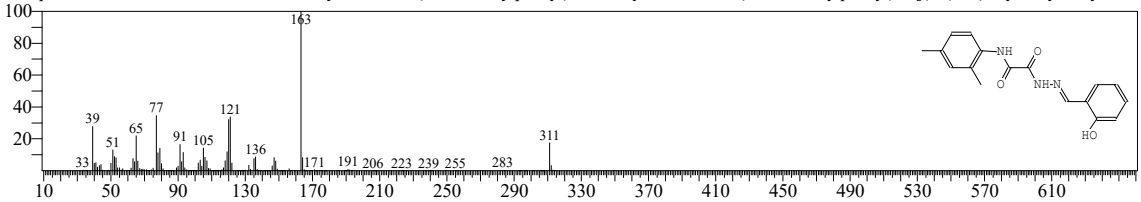
Hit#:2 Entry:57597 Library:NIST20M1.lib
SI:71 Formula:C10H11N3S CAS:39181-45-8 MolWeight:205 RetIndex:1890
CompName:5-(4-Methylbenzyl)-1,3,4-thiadiazol-2-amine $$ 1,3,4-Thiadiazol-2-amine, 5-[(4-methylphenyl)methyl]- $$ 5-[(4-Methylphenyl)methyl]-1,3,4-tl



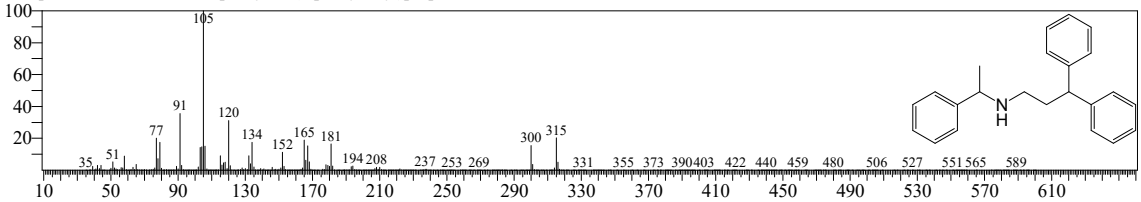
Hit#:3 Entry:25340 Library:NIST20M1.lib
SI:70 Formula:C10H12O2 CAS:6745-75-1 MolWeight:164 RetIndex:1398
CompName:2,5-Dimethyl-para-anisaldehyde $$ 2,5-Dimethyl-4-methoxybenzaldehyde $$ Benzaldehyde, 4-methoxy-2,5-dimethyl- $$ 4-Methoxy-2,5-dimet



Hit#:4 Entry:176186 Library:NIST20M1.lib
SI:70 Formula:C17H17N3O3 CAS:0-00-0 MolWeight:311 RetIndex:3129
CompName:Oxalic acid, monoamide monohydrazide, N-(2,4-dimethylphenyl)-N"-salicylideno- $$ N-(2,4-Dimethylphenyl)-2-[(2E)-2-(2-hydroxybenzyliden

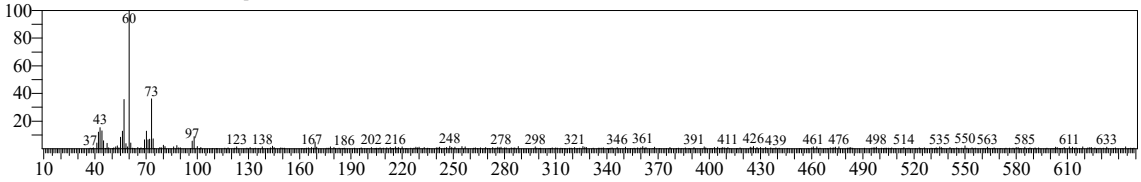


Hit#:5 Entry:530 Library:OA_TMS_DB5_67min_V3.lib
SI:48 Formula:C23H25N CAS:13042-18-7 MolWeight:315 RetIndex:2545
CompName:Fendiline ; 3,3-diphenyl-N-(1-phenylethyl)propan-1-amine

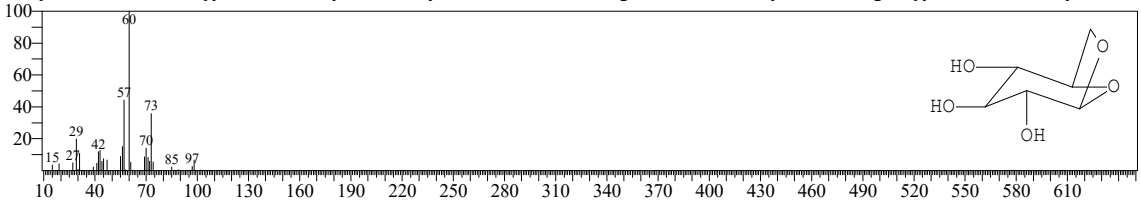


TNAU

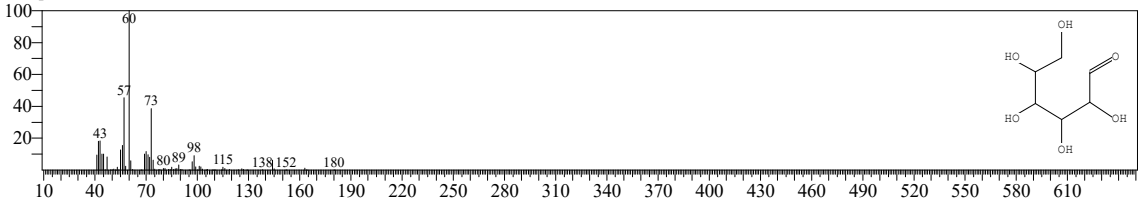
<< Target >>
Line#:11 R.Time:18.110(Scan#:2623) MassPeaks:319
RawMode:Averaged 18.105-18.115(2622-2624) BasePeak:60.00(2583)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



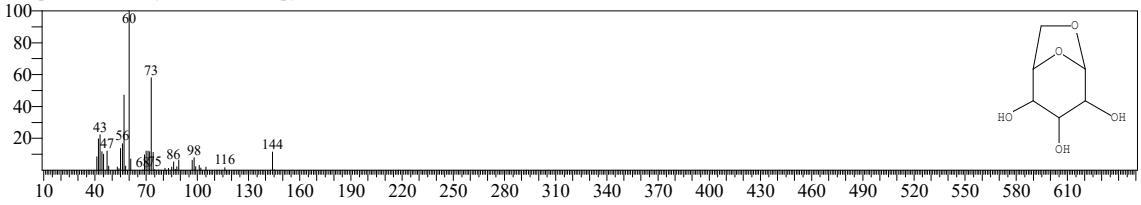
Hit#:1 Entry:23811 Library:NIST20M1.lib
SI:89 Formula:C6H10O5 CAS:498-07-7 MolWeight:162 RetIndex:1404
CompName:.beta.-D-Glucopyranose, 1,6-anhydro- $$ Anhydro-d-mannosan $$ Levoglucosan $$ 1,6-Anhydro-.beta.-D-glucopyranose $$ 1,6-Anhydro-beta-



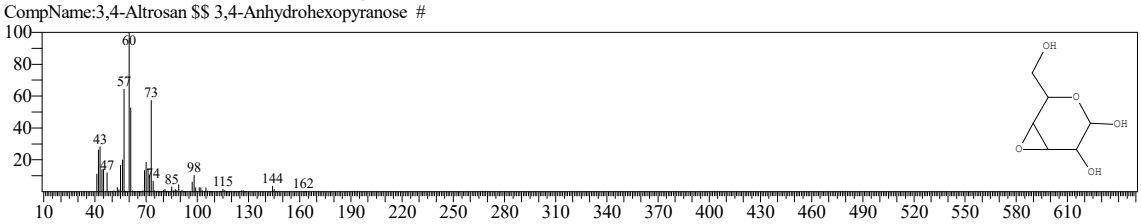
Hit#:2 Entry:36240 Library:NIST20M1.lib
SI:87 Formula:C6H12O6 CAS:2595-97-3 MolWeight:180 RetIndex:1698
CompName:D-Allose $$ .beta.-D-Allose $$ Hexose #



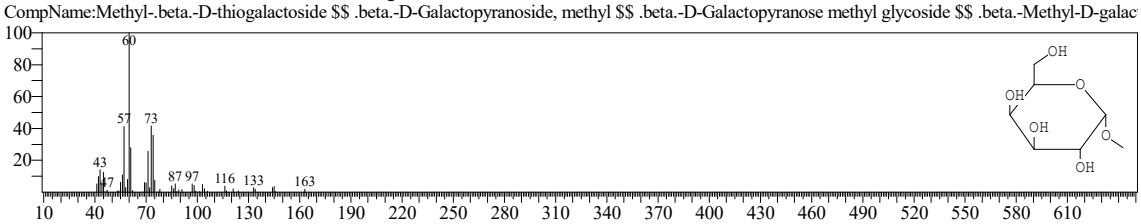
Hit#:3 Entry:23812 Library:NIST20M1.lib
SI:85 Formula:C6H10O5 CAS:0-00-0 MolWeight:162 RetIndex:1404
CompName:1,6-Anhydro-.beta.-d-talopyranose



Hit#:4 Entry:23808 Library:NIST20M1.lib
SI:81 Formula:C6H10O5 CAS:0-00-0 MolWeight:162 RetIndex:1400
CompName:3,4-Altrosan $$ 3,4-Anhydrohexopyranose #

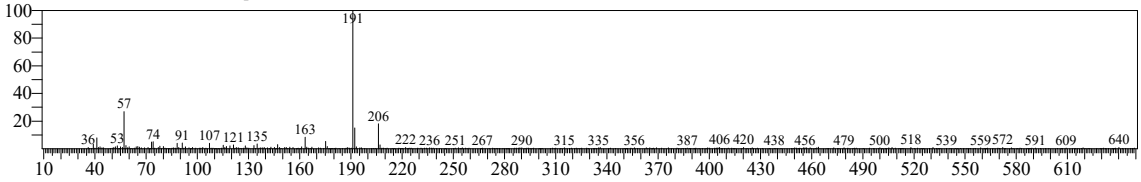


Hit#:5 Entry:47349 Library:NIST20M1.lib
SI:80 Formula:C7H14O6 CAS:1824-94-8 MolWeight:194 RetIndex:1714
CompName:Methyl-.beta.-D-thiogalactoside $$ .beta.-D-Galactopyranoside, methyl $$ .beta.-D-Galactopyranose methyl glycoside $$ .beta.-Methyl-D-galac

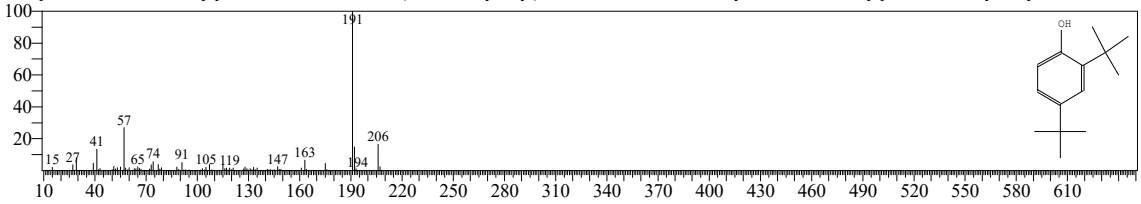


TNAU

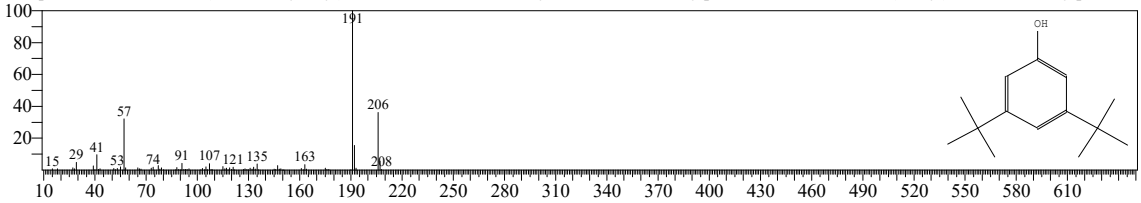
<< Target >>
Line#:12 R.Time:18.495(Scan#:2700) MassPeaks:378
RawMode:Averaged 18.490-18.500(2699-2701) BasePeak:191.10(4239)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



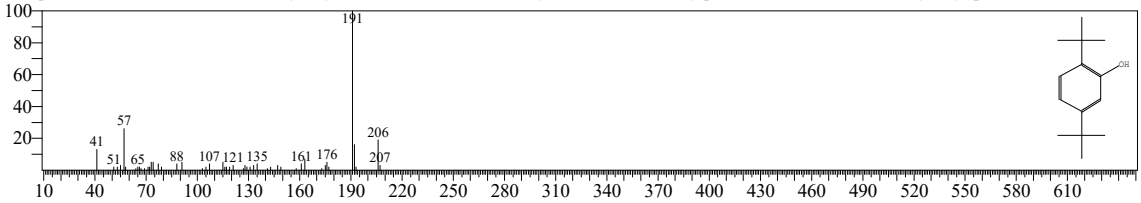
Hit#:1 Entry:24086 Library:NIST20R.lib
SI:91 Formula:C14H22O CAS:96-76-4 MolWeight:206 RetIndex:1555
CompName:2,4-Di-tert-butylphenol $$ Phenol, 2,4-bis(1,1-dimethylethyl)- $$ Phenol, 2,4-di-tert-butyl- $$ 2,4-di-t-Butylphenol $$ 1-Hydroxy-2,4-di-tert-bu



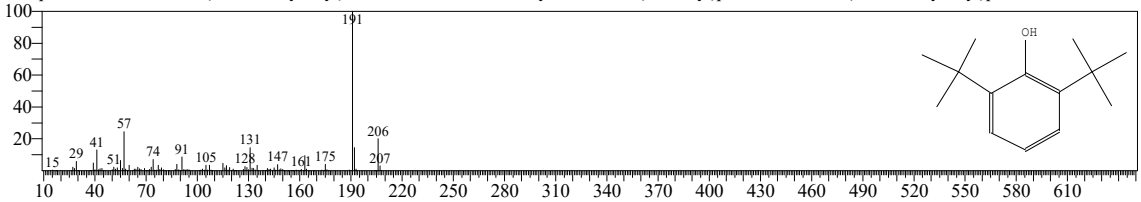
Hit#:2 Entry:24110 Library:NIST20R.lib
SI:88 Formula:C14H22O CAS:1138-52-9 MolWeight:206 RetIndex:1555
CompName:Phenol, 3,5-bis(1,1-dimethylethyl)- $$ Phenol, 3,5-di-tert-butyl- $$ 3,5-Di-tert-butylphenol $$ Phenol, 3,5-bis(t-butyl) $$ 3,5-Di-t-butylphenol $



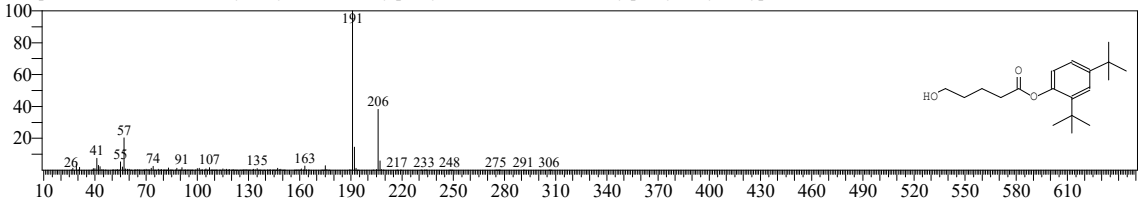
Hit#:3 Entry:24098 Library:NIST20R.lib
SI:88 Formula:C14H22O CAS:5875-45-6 MolWeight:206 RetIndex:1555
CompName:Phenol, 2,5-bis(1,1-dimethylethyl)- $$ Phenol, 2,5-di-tert-butyl- $$ 2,5-Di-tert-butylphenol $$ 2,5-bis(1,1-Dimethylethyl)phenol



Hit#:4 Entry:59031 Library:NIST20M1.lib
SI:86 Formula:C14H22O 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

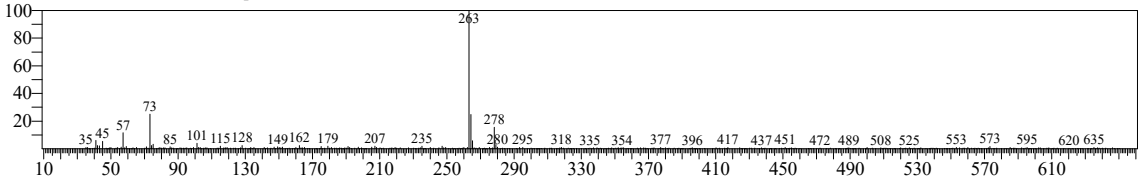


Hit#:5 Entry:170993 Library:NIST20M1.lib
SI:84 Formula:C19H30O3 CAS:166273-38-7 MolWeight:306 RetIndex:2255
CompName:Pentanoic acid, 5-hydroxy-, 2,4-di-t-butylphenyl esters $$ 2,4-Ditert-butylphenyl 5-hydroxypentanoate #

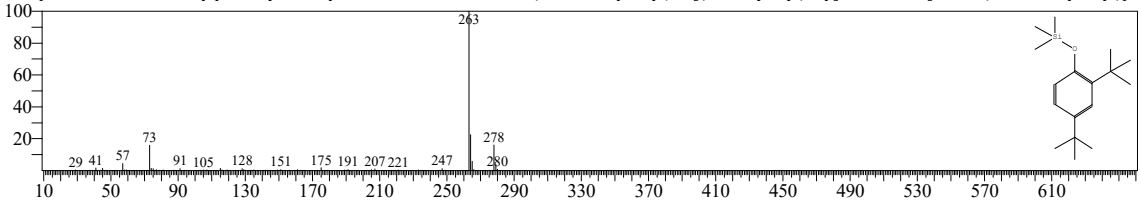


TNAU

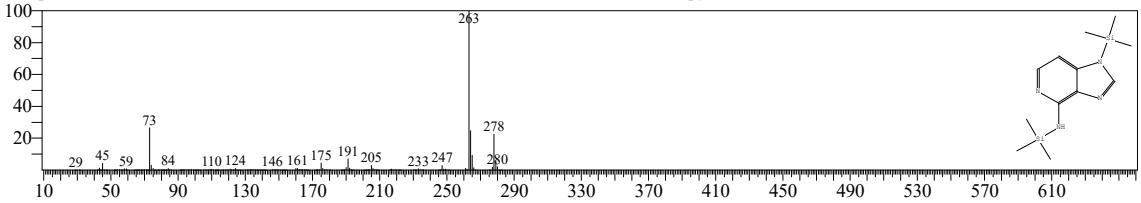
<< Target >>
Line#:13 R.Time:19.190(Scan#:2839) MassPeaks:404
RawMode:Averaged 19.185-19.195(2838-2840) BasePeak:263.15(4303)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



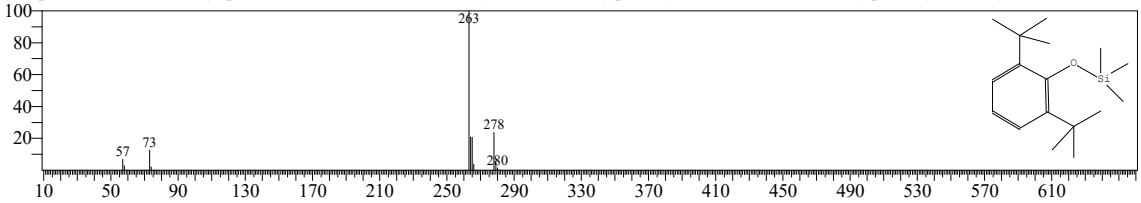
Hit#:1 Entry:33874 Library:NIST20R.lib
SI:86 Formula:C17H30OSi CAS:53925-65-8 MolWeight:278 RetIndex:1632
CompName:2,4-Di-tert-butylphenoxytrimethylsilane $$ Benzene, 2,4-bis(1,1-dimethylethyl)-1-[(trimethylsilyl)oxy]- $$ Silane, [2,4-bis(1,1-dimethylethyl)ph



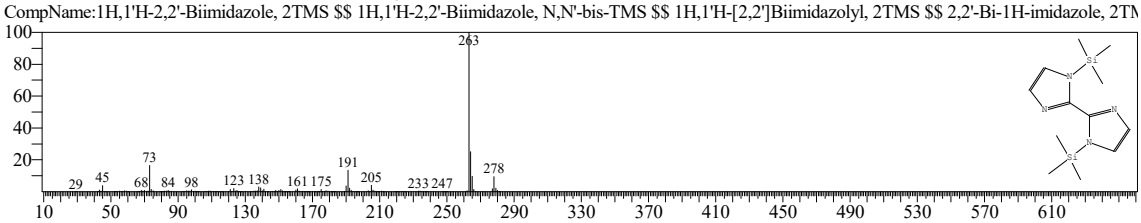
Hit#:2 Entry:136557 Library:NIST20M1.lib
SI:80 Formula:C12H22N4Si2 CAS:0-00-0 MolWeight:278 RetIndex:1703
CompName:3-Deazaadenine, 2TMS $$ 3-Deazaadenine, N,N'-bis-TMS $$ 1H-Imidazo[4,5-c]pyridin-4-amine, N,N'-bis-TMS



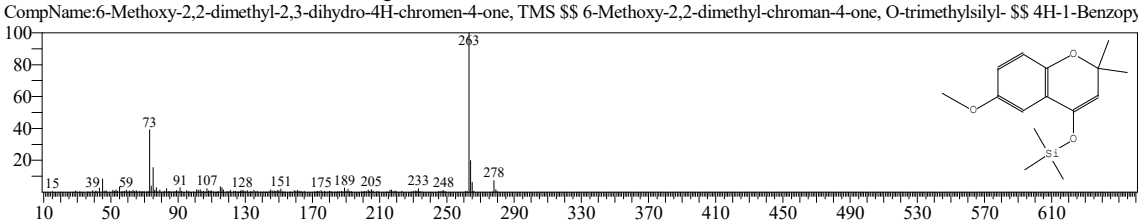
Hit#:3 Entry:33871 Library:NIST20R.lib
SI:78 Formula:C17H30OSi CAS:10416-73-6 MolWeight:278 RetIndex:1632
CompName:2,6-Bis(tert-butyl)phenol, TMS derivative $$ Trimethyl(2,6 ditert.-butylphenoxy)silane $$ (2,6-Ditert-butylphenoxy)(trimethyl)silane #



Hit#:4 Entry:136556 Library:NIST20M1.lib
SI:76 Formula:C12H22N4Si2 CAS:0-00-0 MolWeight:278 RetIndex:1606
CompName:1H,1'H-2,2'-Biimidazole, 2TMS $$ 1H,1'H-2,2'-Biimidazole, N,N'-bis-TMS $$ 1H,1'H-[2,2']Biimidazolyl, 2TMS $$ 2,2'-Bi-1H-imidazole, 2TM

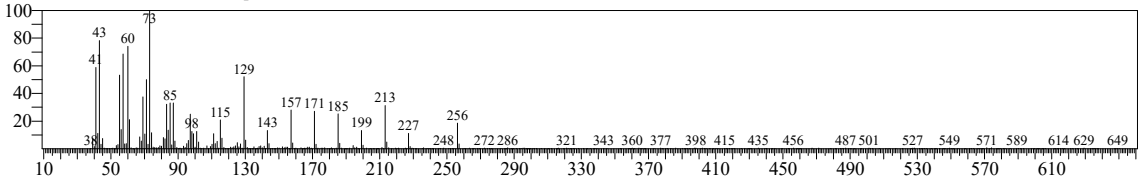


Hit#:5 Entry:137020 Library:NIST20M1.lib
SI:75 Formula:C15H22O3Si CAS:0-00-0 MolWeight:278 RetIndex:1736
CompName:6-Methoxy-2,2-dimethyl-2,3-dihydro-4H-chromen-4-one, TMS $$ 6-Methoxy-2,2-dimethyl-chroman-4-one, O-trimethylsilyl- $$ 4H-1-Benzopy

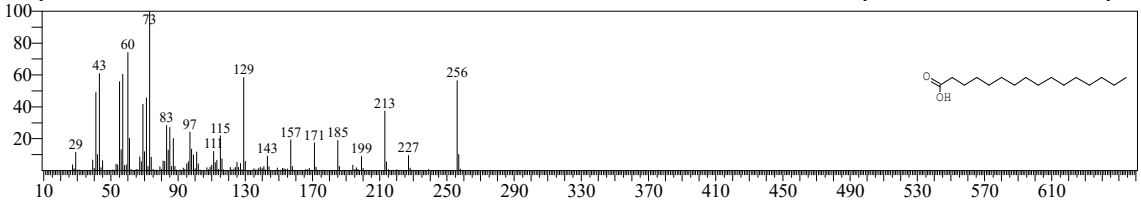


TNAU

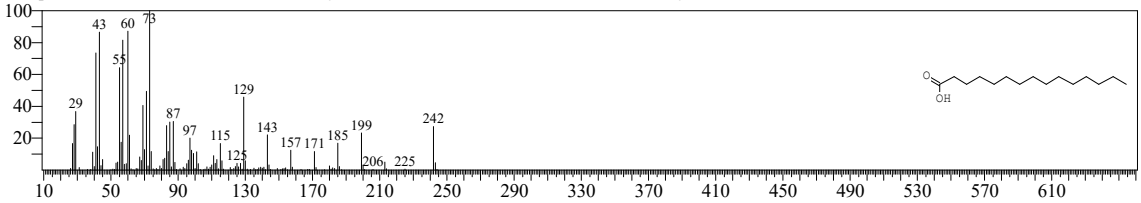
<< Target >>
Line#:14 R.Time:28.295(Scan#:4660) MassPeaks:379
RawMode:Averaged 28.290-28.300(4659-4661) BasePeak:73.05(6539)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



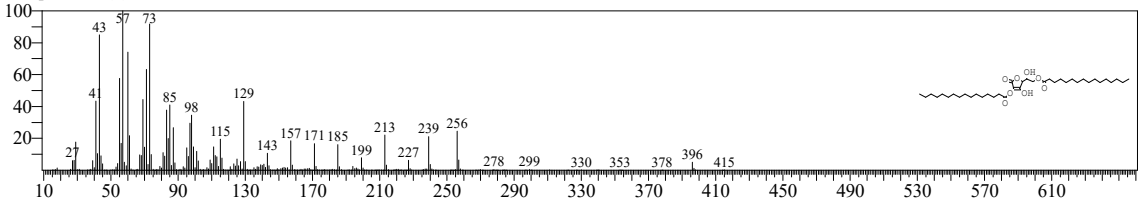
Hit#:1 Entry:31600 Library:NIST20R.lib
SI:95 Formula:C16H32O2 CAS:57-10-3 MolWeight:256 RetIndex:1968
CompName:n-Hexadecanoic acid $$ Hexadecanoic acid $$ n-Hexadecoic acid $$ Palmitic acid $$ Pentadecanecarboxylic acid $$ 1-Pentadecanecarboxylic



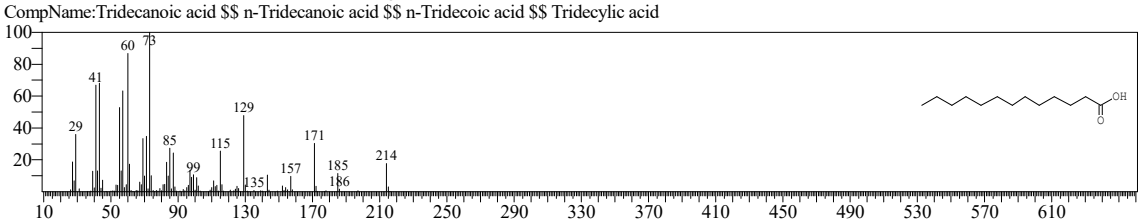
Hit#:2 Entry:29890 Library:NIST20R.lib
SI:91 Formula:C15H30O2 CAS:1002-84-2 MolWeight:242 RetIndex:1869
CompName:Pentadecanoic acid $$ Pentadecylic acid $$ n-Pentadecanoic acid $$ n-Pentadecylic acid



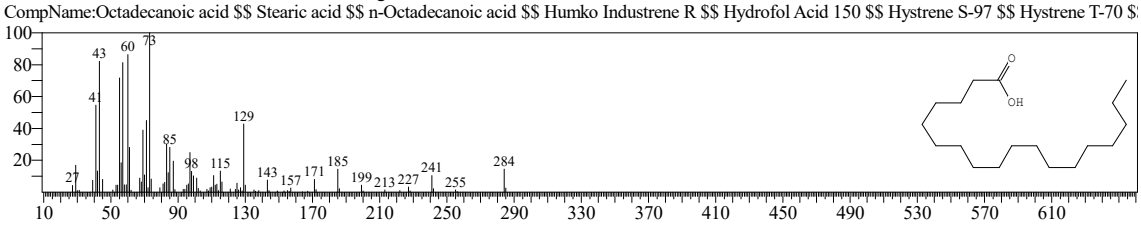
Hit#:3 Entry:44286 Library:NIST20M2.lib
SI:91 Formula:C38H68O8 CAS:28474-90-0 MolWeight:652 RetIndex:4765
CompName:l-(+)-Ascorbic acid 2,6-dihexadecanoate



Hit#:4 Entry:25643 Library:NIST20R.lib
SI:90 Formula:C13H26O2 CAS:638-53-9 MolWeight:214 RetIndex:1670
CompName:Tridecanoic acid $$ n-Tridecanoic acid $$ n-Tridecoic acid $$ Tridecylic acid

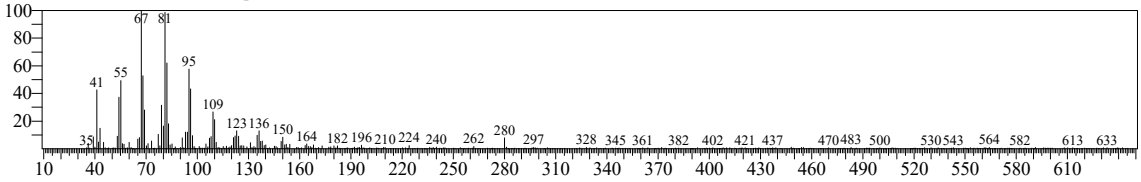


Hit#:5 Entry:144781 Library:NIST20M1.lib
SI:89 Formula:C18H36O2 CAS:57-11-4 MolWeight:284 RetIndex:2167
CompName:Octadecanoic acid $$ Stearic acid $$ n-Octadecanoic acid $$ Humko Industrene R $$ Hydrofol Acid 150 $$ Hystrene S-97 $$ Hystrene T-70 $

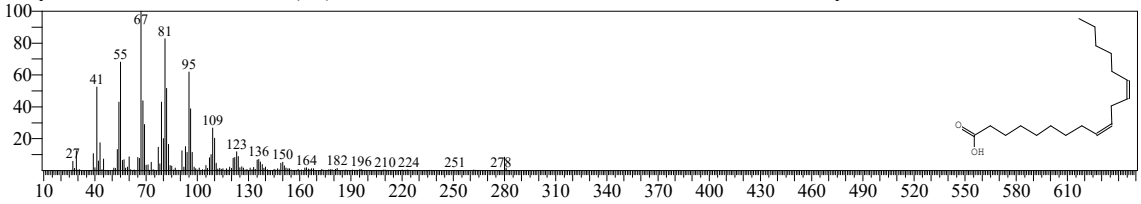


TNAU

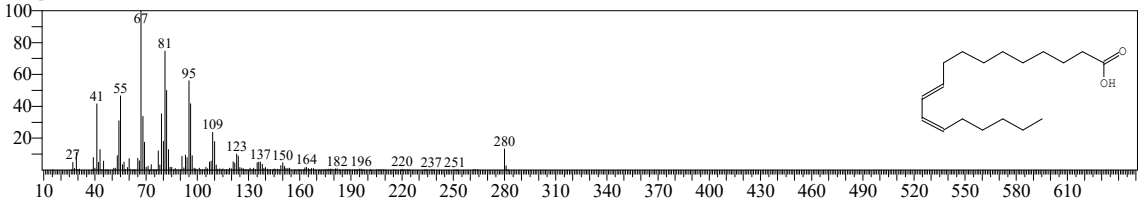
<< Target >>
Line#:15 R.Time:31.470(Scan#:5295) MassPeaks:377
RawMode:Averaged 31.465-31.475(5294-5296) BasePeak:67.05(3894)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



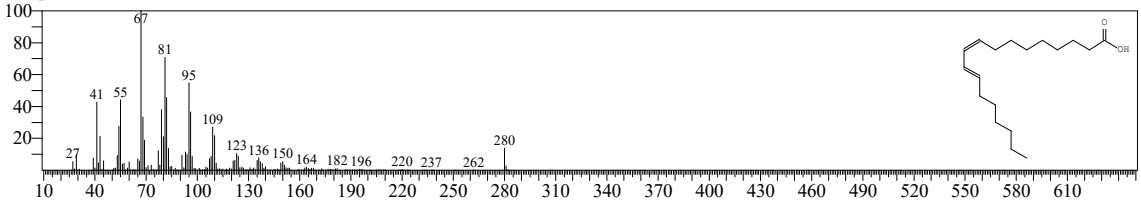
Hit#:1 Entry:34035 Library:NIST20R.lib
SI:94 Formula:C18H32O2 CAS:60-33-3 MolWeight:280 RetIndex:2183
CompName:9,12-Octadecadienoic acid (Z,Z)- $$ cis-9,cis-12-Octadecadienoic acid $$ cis,cis-Linoleic acid $$ Grape seed oil $$ Linoleic $$ Linoleic acid $$



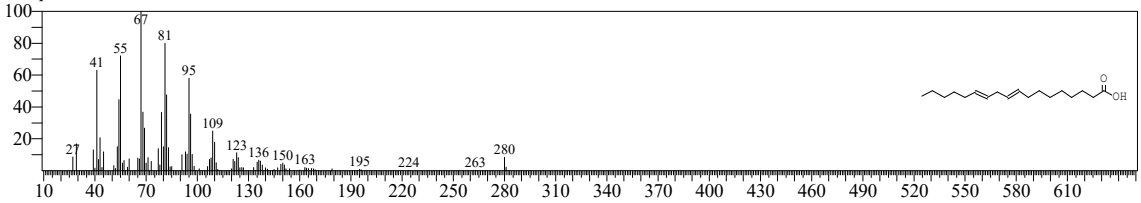
Hit#:2 Entry:139646 Library:NIST20M1.lib
SI:94 Formula:C18H32O2 CAS:2420-56-6 MolWeight:280 RetIndex:2183
CompName:10E,12Z-Octadecadienoic acid



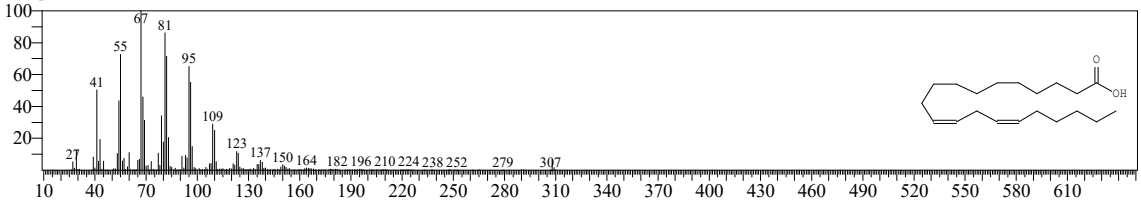
Hit#:3 Entry:139651 Library:NIST20M1.lib
SI:94 Formula:C18H32O2 CAS:544-71-8 MolWeight:280 RetIndex:2183
CompName:(9E,11E)-Octadecadienoic acid



Hit#:4 Entry:139661 Library:NIST20M1.lib
SI:93 Formula:C18H32O2 CAS:506-21-8 MolWeight:280 RetIndex:2183
CompName:Linoelaidic acid

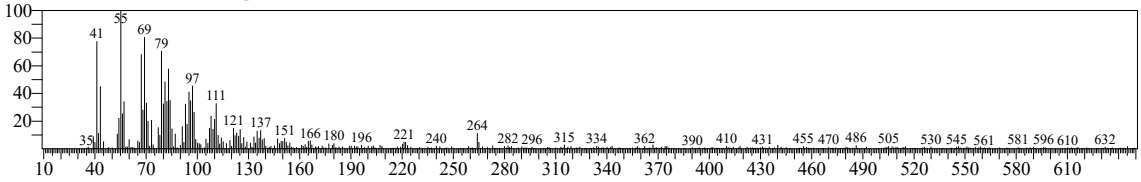


Hit#:5 Entry:173215 Library:NIST20M1.lib
SI:93 Formula:C20H36O2 CAS:2091-39-6 MolWeight:308 RetIndex:2382
CompName:11,14-Eicosadienoic acid

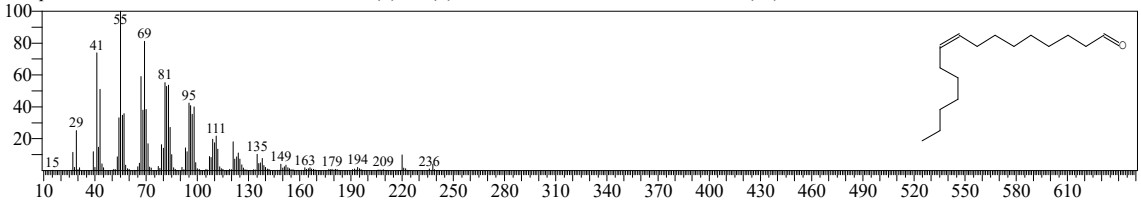


TNAU

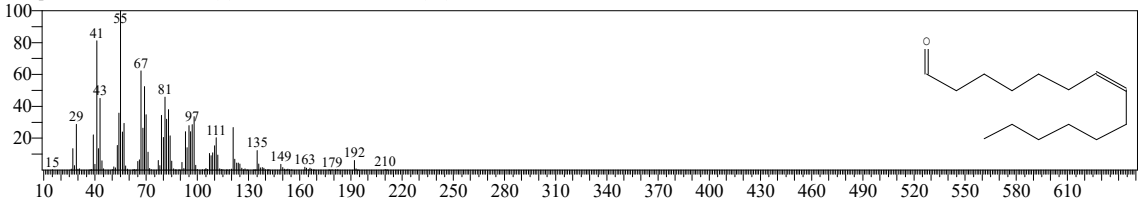
<< Target >>
Line#:16 R.Time:31.580(Scan#:5317) MassPeaks:372
RawMode:Averaged 31.575-31.585(5316-5318) BasePeak:55.10(1892)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



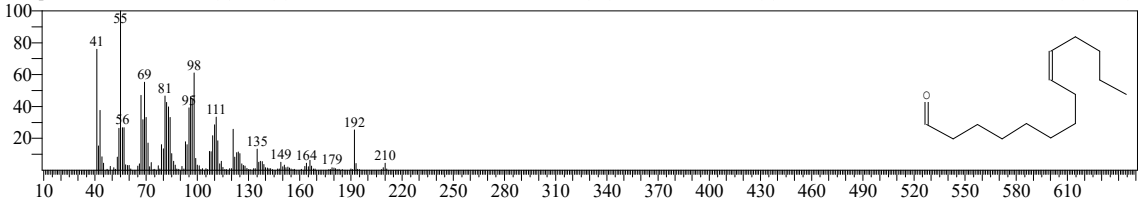
Hit#:1 Entry:92244 Library:NIST20M1.lib
SI:88 Formula:C16H30O CAS:56219-04-6 MolWeight:238 RetIndex:1808
CompName:cis-9-Hexadecenal $$ 9-Hexadecenal, (Z)- $$ (Z)-9-Hexadecenal $$ Z-9-Hexadecenal $$ (9Z)-9-Hexadecenal #



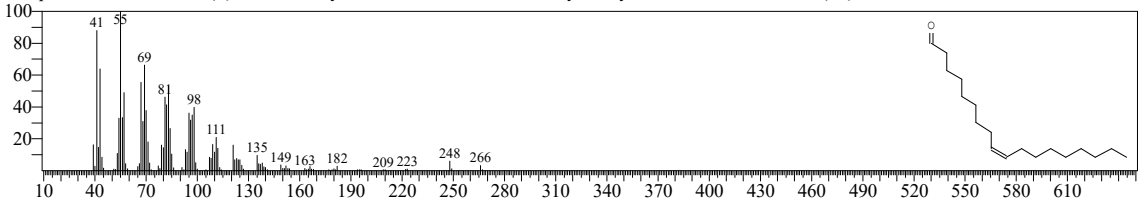
Hit#:2 Entry:63176 Library:NIST20M1.lib
SI:88 Formula:C14H26O CAS:65128-96-3 MolWeight:210 RetIndex:1609
CompName:7-Tetradecenal, (Z)- $$ Z-7-Tetradecenal $$ (7Z)-7-Tetradecenal #



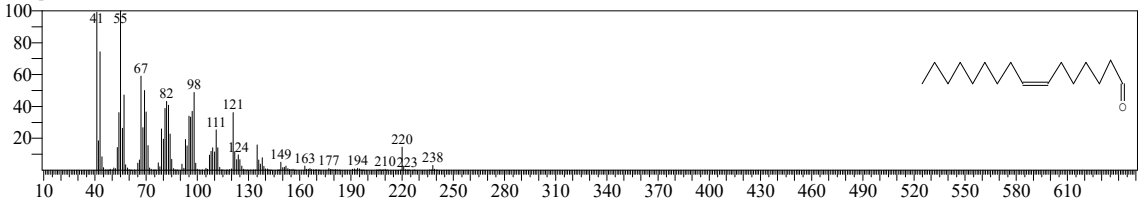
Hit#:3 Entry:24862 Library:NIST20R.lib
SI:87 Formula:C14H26O CAS:53939-27-8 MolWeight:210 RetIndex:1609
CompName:9-Tetradecenal, (Z)- $$ (Z)-9-Tetradecenal $$ Z-9-Tetradecenal $$ (9Z)-9-Tetradecenal $$ (Z)-9-Tetradecen-1-al $$ 9-Tetradecenal, (9Z)- $$ (Z



Hit#:4 Entry:123421 Library:NIST20M1.lib
SI:87 Formula:C18H34O CAS:2423-10-1 MolWeight:266 RetIndex:2007
CompName:9-Octadecenal, (Z)- $$ Olealdehyde $$ cis-9-Octadecenal $$ Oleylaldehyde $$ Z-9-Octadecenal $$ (9Z)-9-Octadecenal #

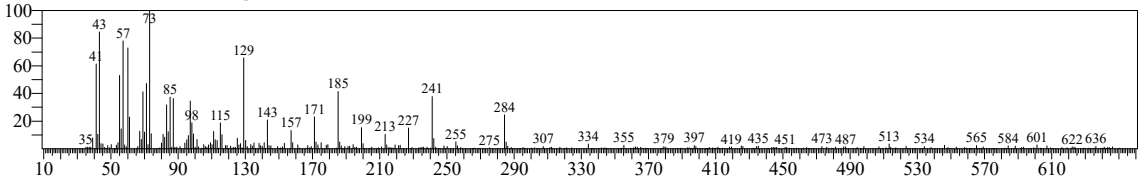


Hit#:5 Entry:92227 Library:NIST20M1.lib
SI:87 Formula:C16H30O CAS:56797-40-1 MolWeight:238 RetIndex:1808
CompName:7-Hexadecenal, (Z)- $$ (Z)-7-Hexadecenal $$ Z-7-Hexadecenal $$ (7Z)-7-Hexadecenal #

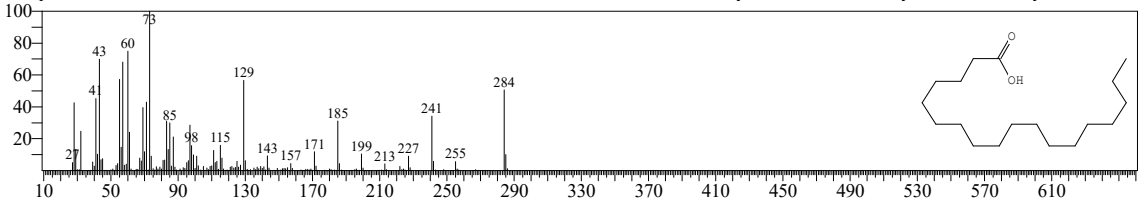


TNAU

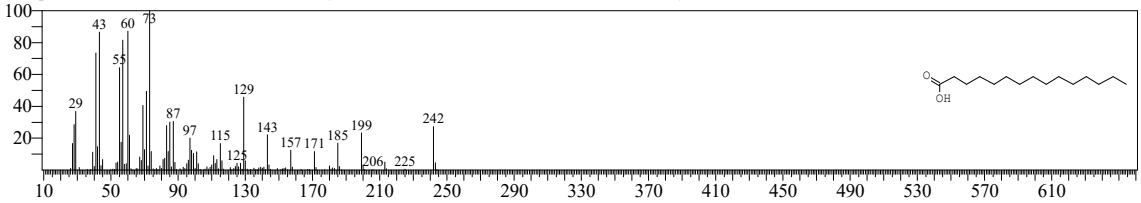
<< Target >>
Line#:17 R.Time:32.030(Scan#:5407) MassPeaks:320
RawMode:Averaged 32.025-32.035(5406-5408) BasePeak:73.05(1442)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



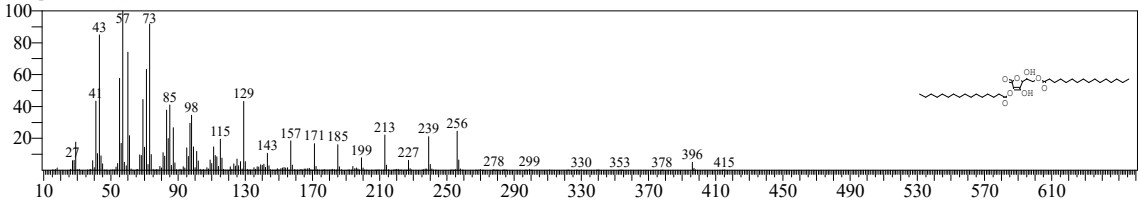
Hit#:1 Entry:34463 Library:NIST20R.lib
SI:93 Formula:C18H36O2 CAS:57-11-4 MolWeight:284 RetIndex:2167
CompName:Octadecanoic acid $$ Stearic acid $$ n-Octadecanoic acid $$ Humko Industrene R $$ Hydrofol Acid 150 $$ Hystrene S-97 $$ Hystrene T-70 $



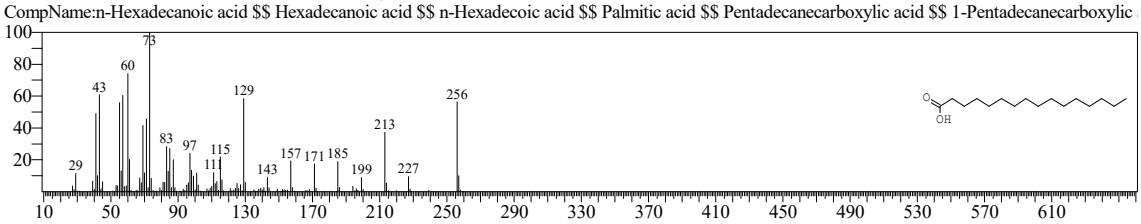
Hit#:2 Entry:29890 Library:NIST20R.lib
SI:90 Formula:C15H30O2 CAS:1002-84-2 MolWeight:242 RetIndex:1869
CompName:Pentadecanoic acid $$ Pentadecylic acid $$ n-Pentadecanoic acid $$ n-Pentadecylic acid



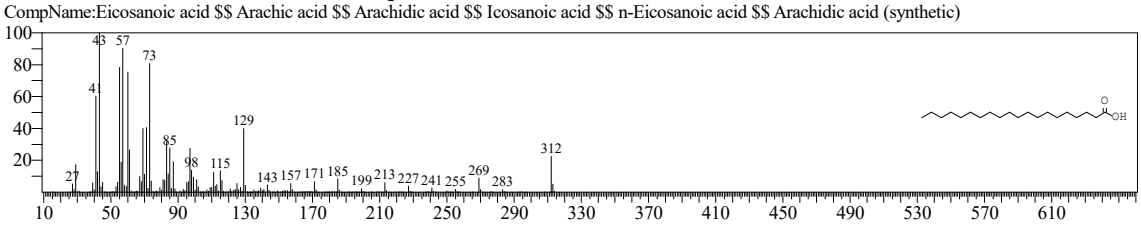
Hit#:3 Entry:44286 Library:NIST20M2.lib
SI:88 Formula:C38H68O8 CAS:28474-90-0 MolWeight:652 RetIndex:4765
CompName:l-(+)-Ascorbic acid 2,6-dihexadecanoate



Hit#:4 Entry:31600 Library:NIST20R.lib
SI:87 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

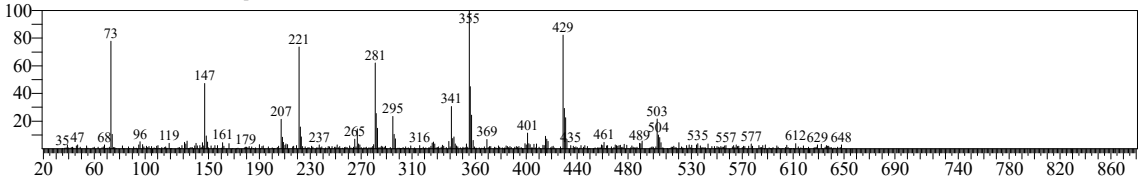


Hit#:5 Entry:36904 Library:NIST20R.lib
SI:87 Formula:C20H40O2 CAS:506-30-9 MolWeight:312 RetIndex:2366
CompName:Eicosanoic acid $$ Arachic acid $$ Arachidic acid $$ Icosanoic acid $$ n-Eicosanoic acid $$ Arachidic acid (synthetic)

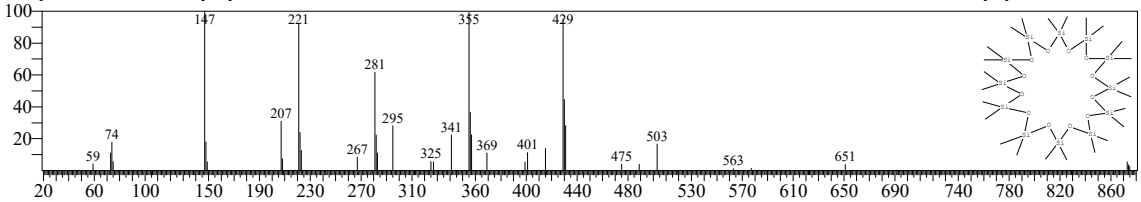


TNAU

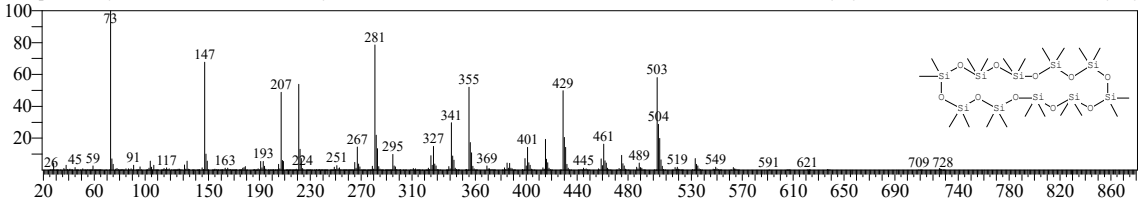
<< Target >>
Line#:18 R.Time:41.600(Scan#:7321) MassPeaks:433
RawMode:Averaged 41.595-41.605(7320-7322) BasePeak:355.10(1160)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



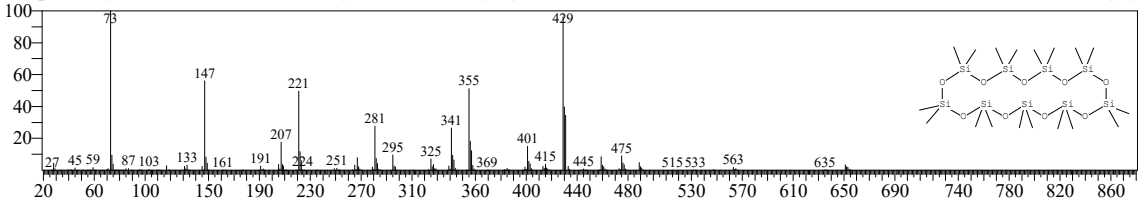
Hit#:1 Entry:46368 Library:NIST20M2.lib
SI:83 Formula:C24H72O12Si12 CAS:18919-94-3 MolWeight:888 RetIndex:2480
CompName:Tetracosamethyl-cyclododecasiloxane $$ 2,2,4,4,6,6,8,8,10,10,12,12,14,14,16,16,18,18,20,20,22,22,24,24-Tetracosamethylcyclododecasiloxan



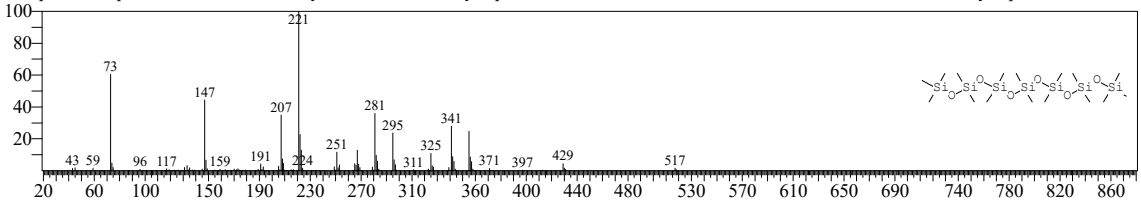
Hit#:2 Entry:45685 Library:NIST20M2.lib
SI:82 Formula:C20H60O10Si10 CAS:18772-36-6 MolWeight:740 RetIndex:2067
CompName:Cyclodecasiloxane, eicosamethyl- $$ 2,2,4,4,6,6,8,8,10,10,12,12,14,14,16,16,18,18,20,20-Icosamethylcyclodecasiloxane # $$ Eicosamethyl-cy



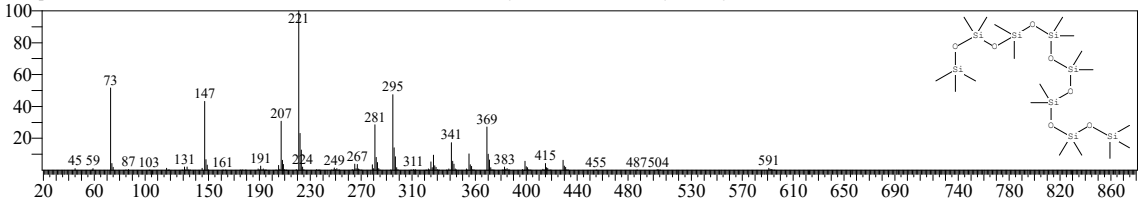
Hit#:3 Entry:44570 Library:NIST20M2.lib
SI:80 Formula:C18H54O9Si9 CAS:556-71-8 MolWeight:666 RetIndex:1860
CompName:Cyclononasiloxane, octadecamethyl- $$ Octadeamethyl-cyclononasiloxane $$ 2,2,4,4,6,6,8,8,10,10,12,12,14,14,16,16,18,18-Octadecamethylcy



Hit#:4 Entry:43312 Library:NIST20R.lib
SI:73 Formula:C16H48O6Si7 CAS:541-01-5 MolWeight:532 RetIndex:1437
CompName:Heptasiloxane, hexadecamethyl- $$ Hexadecamethylheptasiloxane $$ 1,1,1,3,3,5,5,7,7,9,9,11,11,13,13,13-Hexadecamethylheptasiloxane #

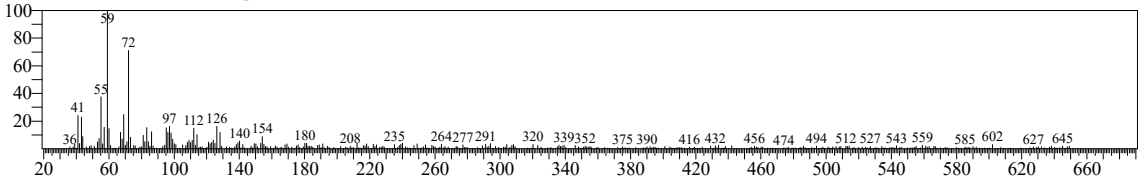


Hit#:5 Entry:42942 Library:NIST20M2.lib
SI:72 Formula:C18H54O7Si8 CAS:556-69-4 MolWeight:606 RetIndex:1622
CompName:1,1,1,3,3,5,5,7,7,9,9,11,11,13,13,15,15,15-octadecamethyloctasiloxane $$ Polydimethylsilane Oil

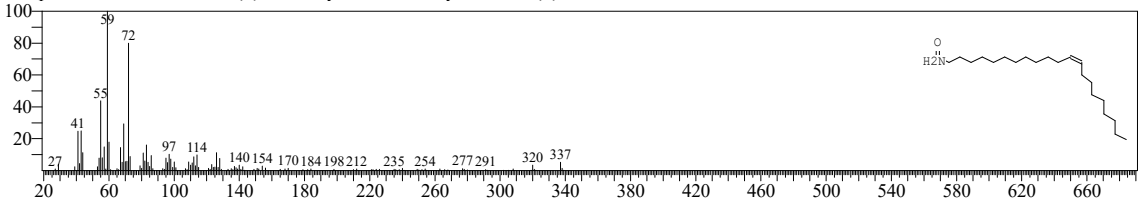


TNAU

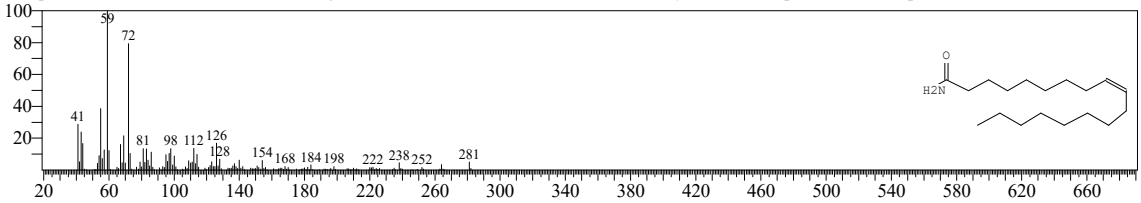
<< Target >>
Line#:19 R.Time:44.155(Scan#:7832) MassPeaks:411
RawMode:Averaged 44.150-44.160(7831-7833) BasePeak:59.05(1584)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



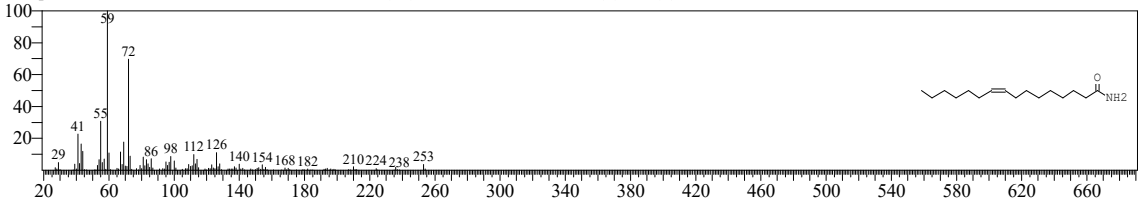
Hit#:1 Entry:38689 Library:NIST20R.lib
SI:88 Formula:C22H43NO CAS:112-84-5 MolWeight:337 RetIndex:2625
CompName:13-Docosenamide, (Z)- $$ Erucylamide $$ Erucyl amide $$ (Z)-13-Docosenamide $$ 13-Docosenamide, cis- $$ Armid E $$ cis-13-Docosenam



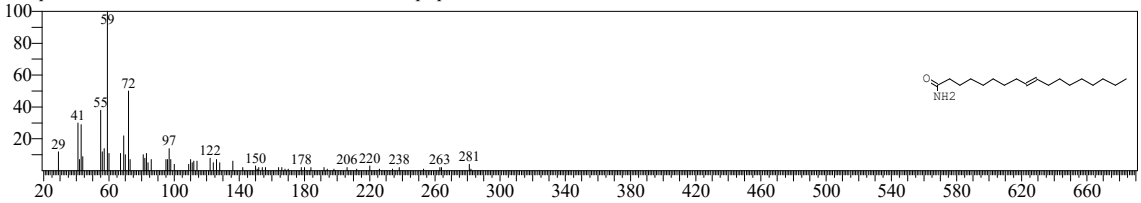
Hit#:2 Entry:34121 Library:NIST20R.lib
SI:88 Formula:C18H35NO CAS:301-02-0 MolWeight:281 RetIndex:2228
CompName:9-Octadecenamide, (Z)- $$ Adogen 73 $$ Oleamide $$ Oleic acid amide $$ Oleyl amide $$ Slip-eze $$ Armoslip CP $$ Crodamide O $$ Croda



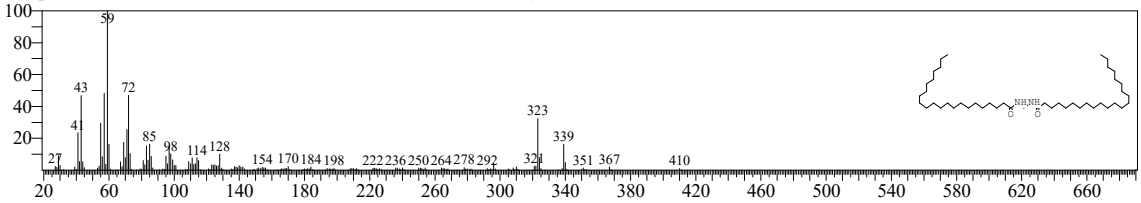
Hit#:3 Entry:108453 Library:NIST20M1.lib
SI:85 Formula:C16H31NO CAS:106010-22-4 MolWeight:253 RetIndex:2029
CompName:Palmitoleamide $$ 9-Hexadecenamide, (9Z)- $$ 9-Hexadecenamide, (Z)- $$ (Z)-Hexadec-9-enamide



Hit#:4 Entry:140706 Library:NIST20M1.lib
SI:82 Formula:C18H35NO CAS:3322-62-1 MolWeight:281 RetIndex:2228
CompName:9-Octadecenamide $$ Armid ow $$ Armoslip cp $$ 9-Oleoamide

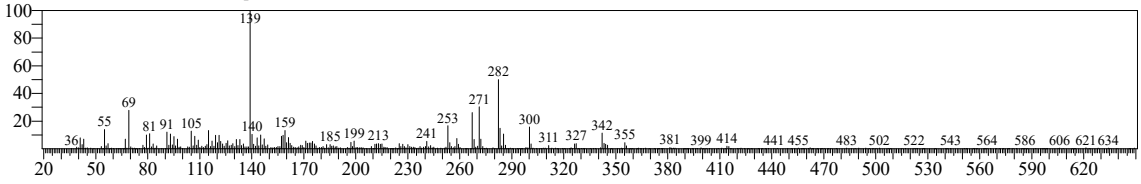


Hit#:5 Entry:45057 Library:NIST20M2.lib
SI:77 Formula:C45H90N2O2 CAS:10436-15-4 MolWeight:690 RetIndex:5295
CompName:Bis(dodecanamido)methane $$ N-[(Docosanoylamino)methyl]docosanamide #

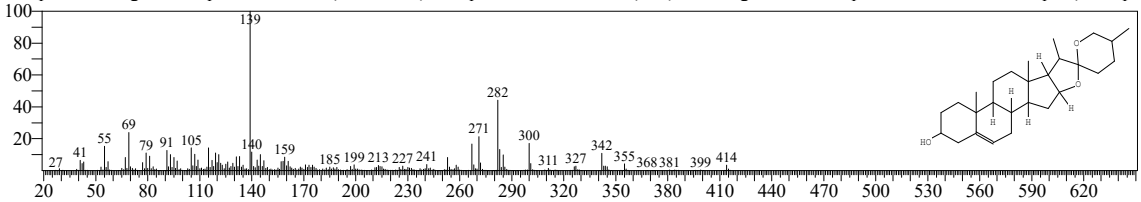


TNAU

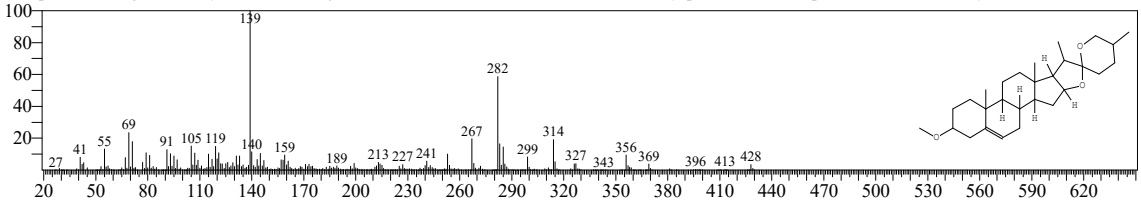
<< Target >>
Line#:20 R.Time:45.600(Scan#:8121) MassPeaks:392
RawMode:Averaged 45.595-45.605(8120-8122) BasePeak:139.10(6556)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



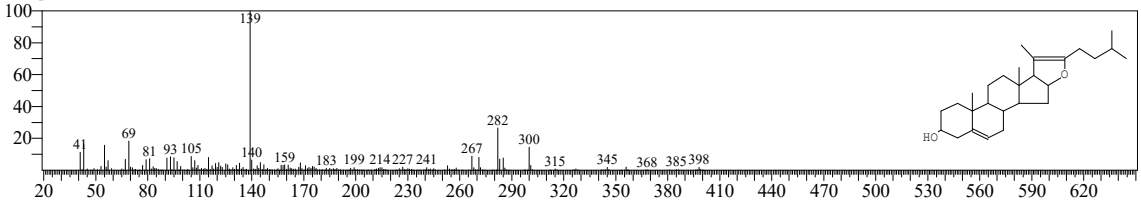
Hit#:1 Entry:8297 Library:NIST20M2.lib
SI:91 Formula:C27H42O3 CAS:512-04-9 MolWeight:414 RetIndex:2844
CompName:Diosgenin $$ Spirost-5-en-3-ol, (3.beta.,25R)- $$ Spirost-5-en-3.beta.-ol, (25R)- $$ Nitogenin $$ 25D-spirost-5-en-3.beta.-ol $$ Spiro(8H-napl



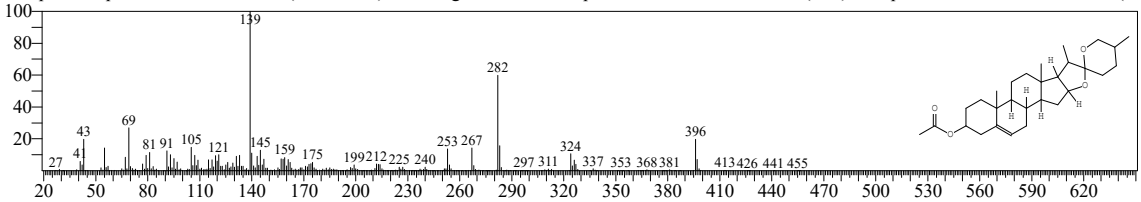
Hit#:2 Entry:14852 Library:NIST20M2.lib
SI:82 Formula:C28H44O3 CAS:116292-24-1 MolWeight:428 RetIndex:2793
CompName:Diosgenin methyl ether $$ Diosgenin Me derivative $$ (3.beta.,25R)-3-Methoxyspirost-5-ene $$ Spirost-5-ene, 3-methoxy-, (3.beta.,25R)-



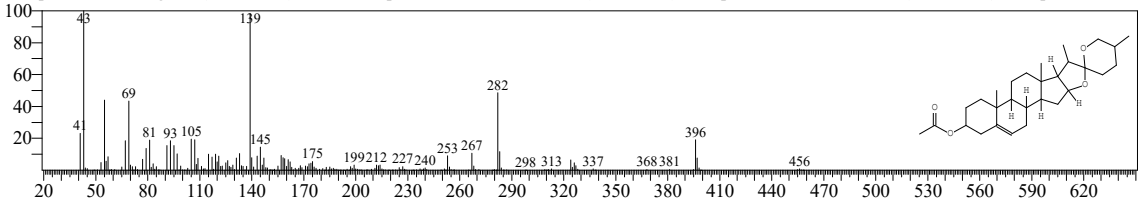
Hit#:3 Entry:41561 Library:NIST20R.lib
SI:81 Formula:C27H42O2 CAS:32277-73-9 MolWeight:398 RetIndex:2713
CompName:Des-N-solasodine $$ Furosta-5,20(22)-dien-3-ol #



Hit#:4 Entry:24878 Library:NIST20M2.lib
SI:79 Formula:C29H44O4 CAS:1061-54-7 MolWeight:456 RetIndex:2984
CompName:Spirost-5-en-3-ol, acetate, (3.beta.,25R)- $$ Diosgenin acetate $$ Spirost-5-en-3.beta.-ol, acetate, (25R)- $$ Spirost-5-en-3-.beta.-ol, acetate, (2

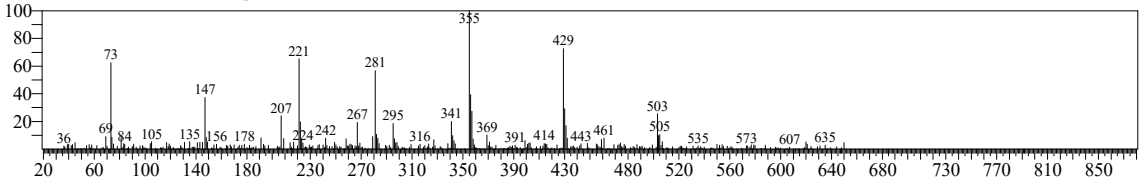


Hit#:5 Entry:42677 Library:NIST20R.lib
SI:75 Formula:C29H44O4 CAS:1180-12-7 MolWeight:456 RetIndex:2984
CompName:Neodiosgenin (3.beta.,25S) acetate $$ Spirost-5-en-3-ol, 3-acetate, (3.beta.,25S)- $$ Spirost-5-en-3-ol, acetate, (3.beta.,25S)- $$ Spirost-5-en-3.

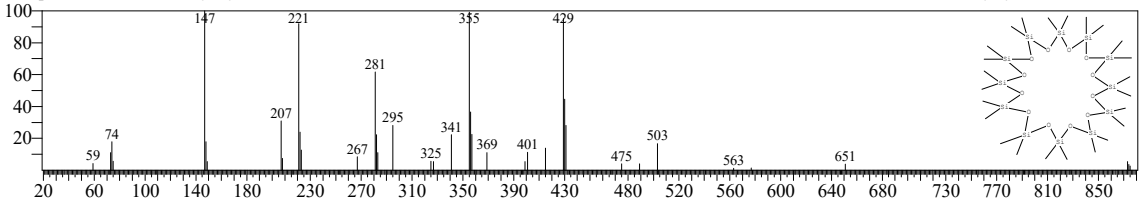


TNAU

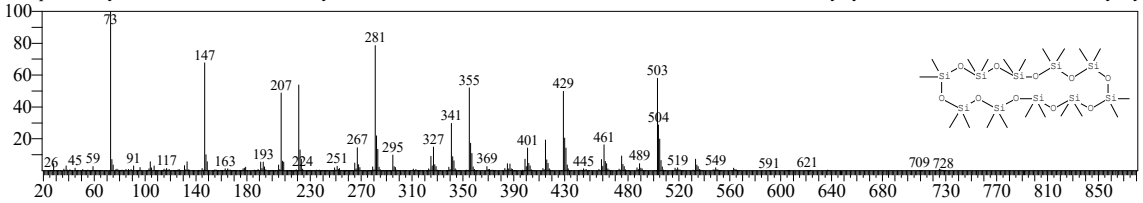
<< Target >>
Line#:21 R.Time:45.765(Scan#:8154) MassPeaks:369
RawMode:Averaged 45.760-45.770(8153-8155) BasePeak:355.00(886)
BG Mode:Calc. from Peak Group 1 - Event 1 Q3 Scan



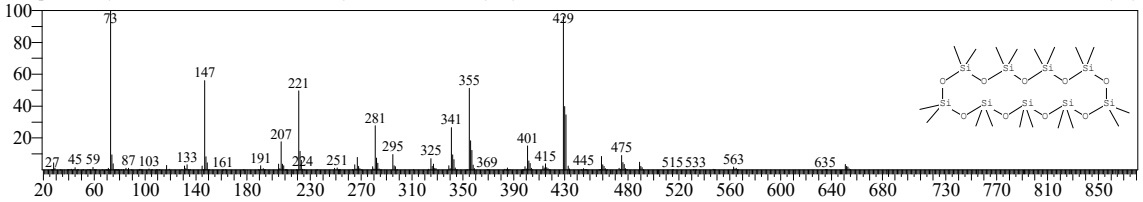
Hit#:1 Entry:46368 Library:NIST20M2.lib
SI:78 Formula:C24H72O12Si12 CAS:18919-94-3 MolWeight:888 RetIndex:2480
CompName:Tetracosamethyl-cyclododecasiloxane $$ 2,2,4,4,6,6,8,8,10,10,12,12,14,14,16,16,18,18,20,20,22,22,24,24-Tetracosamethylcyclododecasiloxan



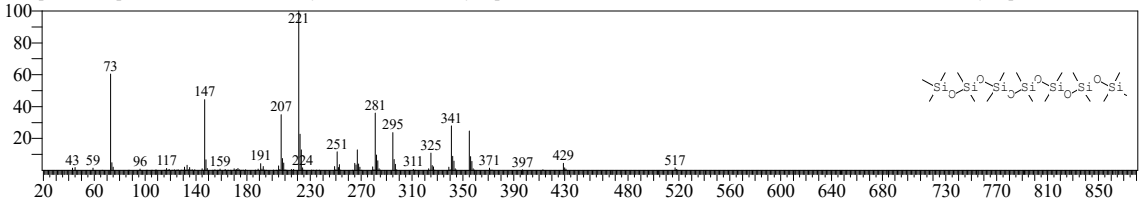
Hit#:2 Entry:45685 Library:NIST20M2.lib
SI:76 Formula:C20H60O10Si10 CAS:18772-36-6 MolWeight:740 RetIndex:2067
CompName:Cyclodecasiloxane, eicosamethyl- $$ 2,2,4,4,6,6,8,8,10,10,12,12,14,14,16,16,18,18,20,20-Icosamethylcyclodecasiloxane # $$ Eicosamethyl-cy



Hit#:3 Entry:44570 Library:NIST20M2.lib
SI:75 Formula:C18H54O9Si9 CAS:556-71-8 MolWeight:666 RetIndex:1860
CompName:Cyclononasiloxane, octadecamethyl- $$ Octadeamethyl-cyclononasiloxane $$ 2,2,4,4,6,6,8,8,10,10,12,12,14,14,16,16,18,18-Octadecamethylcy



Hit#:4 Entry:43312 Library:NIST20R.lib
SI:71 Formula:C16H48O6Si7 CAS:541-01-5 MolWeight:532 RetIndex:1437
CompName:Heptasiloxane, hexadecamethyl- $$ Hexadecamethylheptasiloxane $$ 1,1,1,3,3,5,5,7,7,9,9,11,11,13,13,13-Hexadecamethylheptasiloxane #



Hit#:5 Entry:402 Library:OA_TMS_DB5_67min_V3.lib
SI:59 Formula:C20H42O4Si4 CAS:775-01-9 MolWeight:458 RetIndex:1942
CompName:3,4-Dihydroxymandelic acid-4TMS ; 2-(3,4-dihydroxyphenyl)-2-hydroxyacetic acid