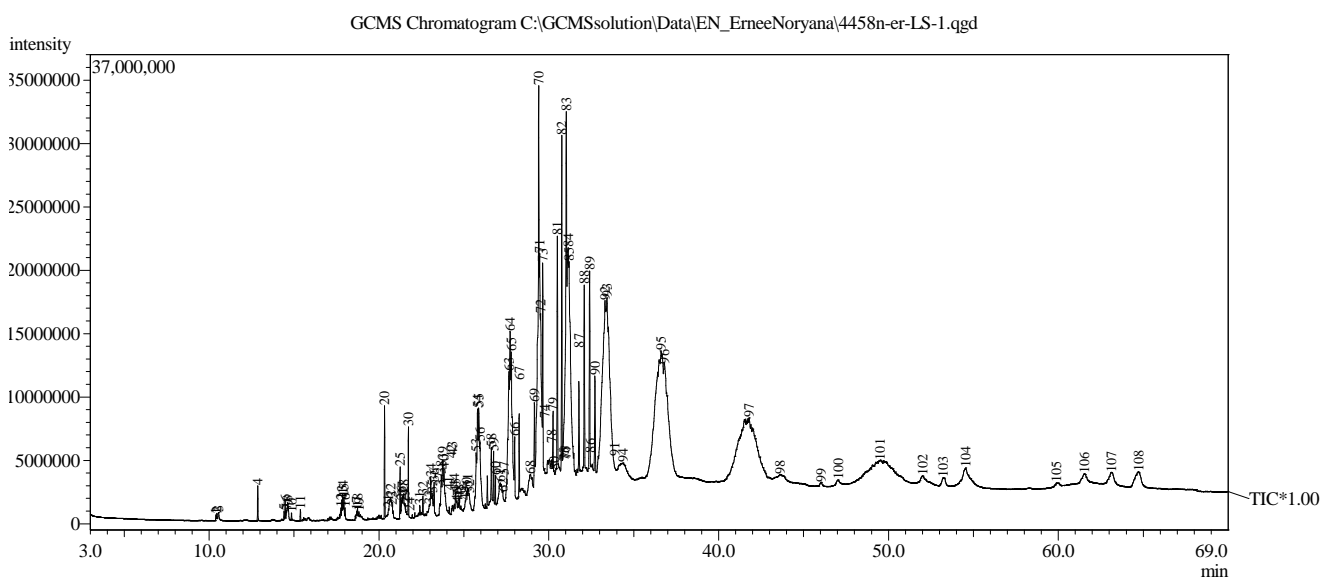


MAKMAL GCMS G39 BSF, JABATAN KIMIA, FAKULTI SAINS, UPM, SERDANG, SELANGOR
GCMS QP2010 Plus SHIMADZU

GCMS Sample Information

Data Acquired by : Admin
Acquisition Date : 1/11/2018 11:41:09 AM
Sample Type : Unknown
Level # : 1
Sample Name : SealerLiquid
Sample ID : Dr.ErneeNoryana
IS Amount : [1]=1
Sample Amount : 5
Dilution Factor : 100
Vial # : 8
Injection Volume : 0.3
Data File : C:\GCMSsolution\Data\EN_ErneeNoryana\4458n-er-LS-1.qgd
Method File : C:\GCMSsolution\Data\Qbio_FaridahQ\Q-ZB5ms50320.qgm
Report File : gcpotrait-ZB-5ms.qgr
Tuning File : C:\GCMSsolution\System\Tune1\241117EIF2zb5ms.qgt
Modified by : Admin
Modified : 1/11/2018 1:00:31 PM



Peak#	R.Time	Area	Height	A/H	Mark	Base m/z	Area%	Height%
1	10.400	893143	430902	2.07		59.10	0.03	0.09
2	10.458	1293598	541945	2.39	V	59.10	0.04	0.11
3	10.544	2246573	572585	3.92	V	59.10	0.07	0.12
4	12.863	5336798	2759148	1.93		57.10	0.17	0.57
5	14.416	1348773	728344	1.85		59.10	0.04	0.15
6	14.480	2937380	1496029	1.96	V	59.10	0.09	0.31
7	14.542	1672069	726509	2.30	V	59.10	0.05	0.15
8	14.624	5227529	1375775	3.80	V	59.10	0.17	0.28
9	14.675	2422081	1062634	2.28	V	59.10	0.08	0.22
10	14.853	973176	549961	1.77	V	73.05	0.03	0.11
11	15.374	1618891	909420	1.78		205.10	0.05	0.19
12	17.761	1818487	763998	2.38		59.10	0.06	0.16
13	17.813	3098822	1420951	2.18	V	59.10	0.10	0.29
14	17.871	4506786	1870307	2.41	V	59.10	0.14	0.38
15	17.938	4536080	1672413	2.71	V	59.10	0.14	0.34
16	18.005	1894743	873098	2.17	V	59.10	0.06	0.18
17	18.692	1789494	618583	2.89	V	59.10	0.06	0.13
18	18.741	2070753	823274	2.52	V	59.10	0.07	0.17
19	18.823	999389	464466	2.15	V	59.10	0.03	0.10
20	20.332	15809336	8872448	1.78		216.95	0.50	1.83
21	20.558	1276150	510794	2.50		59.10	0.04	0.11
22	20.694	10391518	1470123	7.07	V	59.10	0.33	0.30
23	20.783	4047491	873449	4.63	V	59.10	0.13	0.18
24	20.842	979176	487157	2.01	V	59.10	0.03	0.10
25	21.243	7828428	4087874	1.92		216.95	0.25	0.84
26	21.333	4793252	1415705	3.39	V	59.10	0.15	0.29
27	21.379	3859613	1763727	2.19	V	59.10	0.12	0.36
28	21.463	8321703	1911862	4.35	V	59.10	0.26	0.39
29	21.542	4533932	1145387	3.96	V	59.10	0.14	0.24
30	21.732	15204190	7150662	2.13	V	71.05	0.48	1.47
31	22.408	1663937	757718	2.20		57.10	0.05	0.16
32	22.585	3220793	1486548	2.17	V	71.05	0.10	0.31
33	22.958	1574218	425006	3.70		59.10	0.05	0.09
34	23.052	7950825	2644361	3.01	V	59.10	0.25	0.54
35	23.142	5796484	1385168	4.18	V	59.10	0.18	0.28
36	23.243	5483456	2384748	2.30	V	127.15	0.17	0.49
37	23.575	3135307	863437	3.63		59.10	0.10	0.18
38	23.670	12215871	3124818	3.91	V	59.10	0.39	0.64
39	23.758	17063482	4281336	3.99	V	59.10	0.54	0.88
40	23.830	21291551	3583677	5.94	V	59.10	0.68	0.74

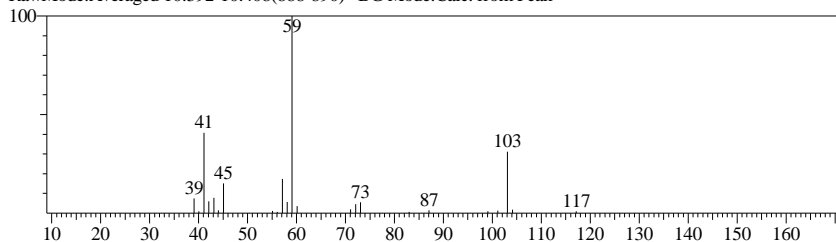
Peak#	R.Time	Area	Height	A/H	Mark	Base m/z	Area%	Height%
41	24.141	2467269	591878	4.17	V	196.10	0.08	0.12
42	24.267	883967	427201	2.07		210.10	0.03	0.09
43	24.318	2863145	691298	4.14	V	224.10	0.09	0.14
44	24.443	2503205	706218	3.54	V	224.10	0.08	0.15
45	24.513	4845492	1587450	3.05	V	177.10	0.15	0.33
46	24.592	2198598	804047	2.73	V	224.10	0.07	0.17
47	24.650	2318739	593875	3.90	V	238.15	0.07	0.12
48	24.723	2386982	946556	2.52	V	210.10	0.08	0.19
49	25.083	1429621	532771	2.68		59.10	0.05	0.11
50	25.172	4926799	1207254	4.08	V	59.10	0.16	0.25
51	25.229	7026071	1444719	4.86	V	59.10	0.22	0.30
52	25.317	2608600	878143	2.97	V	59.10	0.08	0.18
53	25.724	18603316	4273598	4.35		59.10	0.59	0.88
54	25.807	30097445	7755683	3.88	V	59.10	0.96	1.60
55	25.871	31781887	7838114	4.05	V	59.10	1.01	1.61
56	25.950	24243343	5170600	4.69	V	59.10	0.77	1.06
57	26.385	4352604	2407571	1.81		368.15	0.14	0.50
58	26.617	8649384	4222741	2.05		368.15	0.28	0.87
59	26.750	7201635	4081030	1.76	V	85.10	0.23	0.84
60	26.853	4246060	2310975	1.84		368.20	0.14	0.48
61	27.149	17370114	1521703	11.41		59.10	0.55	0.31
62	27.292	2421472	733231	3.30	V	57.10	0.08	0.15
63	27.653	63996092	10086965	6.34	V	59.10	2.03	2.07
64	27.720	50357532	13235167	3.80	V	59.10	1.60	2.72
65	27.785	65894684	11582710	5.69	V	59.10	2.10	2.38
66	27.990	10324648	4780947	2.16	V	85.10	0.33	0.98
67	28.255	11501637	6347382	1.81		85.10	0.37	1.31
68	28.916	14988999	1232091	12.17		59.10	0.48	0.25
69	29.153	15243351	6758957	2.26	V	85.10	0.48	1.39
70	29.408	170190130	31626027	5.38	V	85.10	5.41	6.51
71	29.460	56218647	18355845	3.06	V	59.10	1.79	3.78
72	29.517	58239933	13569084	4.29	V	59.10	1.85	2.79
73	29.643	38753976	17504020	2.21	V	57.10	1.23	3.60
74	29.750	4586077	1196712	3.83	V	57.10	0.15	0.25
75	29.917	10232782	1687958	6.06	V	336.30	0.33	0.35
76	29.972	4744700	1728237	2.75	V	57.10	0.15	0.36
77	30.031	9985333	1760633	5.67	V	336.30	0.32	0.36
78	30.157	7631985	1684210	4.53	V	57.10	0.24	0.35
79	30.248	12496016	5464625	2.29	V	113.10	0.40	1.12
80	30.308	2392214	764818	3.13	V	350.30	0.08	0.16
81	30.504	51592597	19118973	2.70	V	113.10	1.64	3.93
82	30.765	64738659	26945559	2.40	V	57.10	2.06	5.54
83	31.027	156591480	28664935	5.46	V	57.10	4.98	5.90
84	31.140	102245539	17859168	5.73	V	59.10	3.25	3.67
85	31.192	120885787	16715323	7.23	V	59.10	3.84	3.44
86	31.470	3016327	1468251	2.05	V	113.10	0.10	0.30
87	31.776	16101704	7100834	2.27		113.10	0.51	1.46
88	32.088	35817319	14420504	2.48		57.10	1.14	2.97
89	32.402	41188050	15610679	2.64		57.10	1.31	3.21
90	32.718	21408103	7550014	2.84		57.10	0.68	1.55
91	32.900	5352902	1208855	4.43		57.10	0.17	0.25
92	33.310	170990541	13466523	12.70	V	59.10	5.44	2.77
93	33.429	224343198	13652864	16.43	V	59.10	7.13	2.81
94	34.362	20643229	789932	25.76	MI	57.10	0.66	0.16
95	36.618	280998064	9406792	29.87		59.10	8.94	1.94
96	36.805	164103135	8530127	19.24	V	59.10	5.22	1.75
97	41.796	356292521	4839480	73.61	MI	59.10	11.33	1.00
98	43.617	14199508	470204	29.86	MI	57.10	0.45	0.10
99	46.034	2467204	301477	8.19	MI	85.10	0.08	0.06
100	47.043	5597038	394361	13.92	MI	57.10	0.18	0.08
101	49.504	176301260	1791996	98.38	MI	59.10	5.61	0.37
102	52.018	7950027	495504	16.06	MI	57.10	0.25	0.10
103	53.215	8909162	619934	14.38	MI	57.10	0.28	0.13
104	54.550	30898785	1239748	24.92		57.10	0.98	0.26
105	59.891	4336819	279754	15.47	MI	85.10	0.14	0.06
106	61.543	13088671	690043	18.97		57.10	0.42	0.14
107	63.131	17724459	902766	19.63		57.10	0.56	0.19
108	64.718	30723716	1217044	25.24		57.10	0.98	0.25
		3144835566	486131035				100.00	100.00

GCMS Library

<< Target >>

Line#1 R.Time:10.400(Scan#:889) BasePeak:59.10(113786)

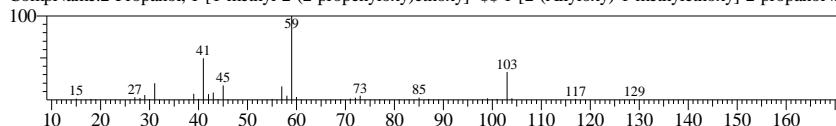
RawMode:Averaged 10.392-10.408(888-890) BG Mode:Calc. from Peak



Hit#1 Entry:12792 Library:NIST08s.LIB

SI:96 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

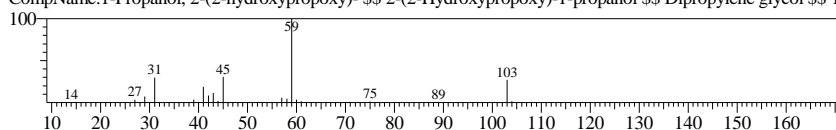
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#2 Entry:5921 Library:NIST08s.LIB

SI:90 Formula:C6H14O3 CAS:106-62-7 MolWeight:134 RetIndex:1034

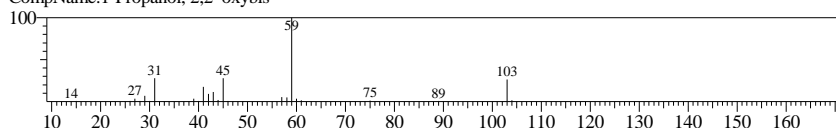
CompName:1-Propanol, 2-(2-hydroxypropoxy)- \$\$ 2-(2-Hydroxypropoxy)-1-propanol \$\$ Dipropylene glycol \$\$ 1



Hit#3 Entry:8889 Library:NIST08.LIB

SI:90 Formula:C6H14O3 CAS:108-61-2 MolWeight:134 RetIndex:1051

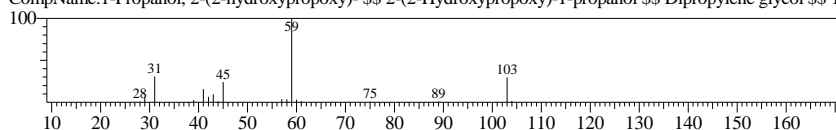
CompName:1-Propanol, 2,2'-oxybis-



Hit#4 Entry:5920 Library:NIST08s.LIB

SI:90 Formula:C6H14O3 CAS:106-62-7 MolWeight:134 RetIndex:1034

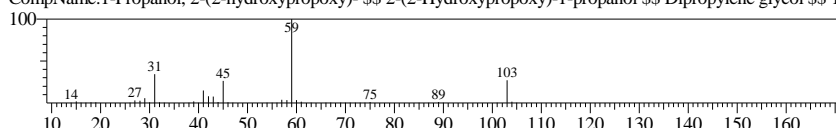
CompName:1-Propanol, 2-(2-hydroxypropoxy)- \$\$ 2-(2-Hydroxypropoxy)-1-propanol \$\$ Dipropylene glycol \$\$ 1



Hit#5 Entry:8891 Library:NIST08.LIB

SI:90 Formula:C6H14O3 CAS:106-62-7 MolWeight:134 RetIndex:1034

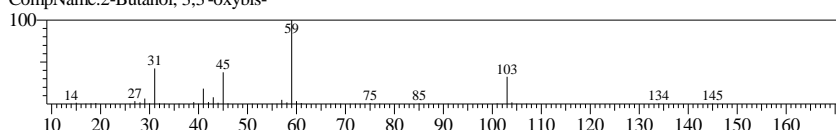
CompName:1-Propanol, 2-(2-hydroxypropoxy)- \$\$ 2-(2-Hydroxypropoxy)-1-propanol \$\$ Dipropylene glycol \$\$ 1



Hit#6 Entry:20866 Library:NIST08.LIB

SI:89 Formula:C8H18O3 CAS:54305-61-2 MolWeight:162 RetIndex:1089

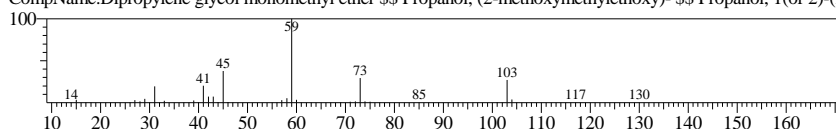
CompName:2-Butanol, 3,3'-oxybis-



Hit#7 Entry:14295 Library:NIST08.LIB

SI:89 Formula:C7H16O3 CAS:34590-94-8 MolWeight:148 RetIndex:983

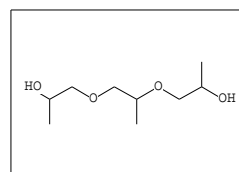
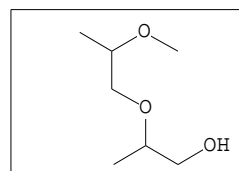
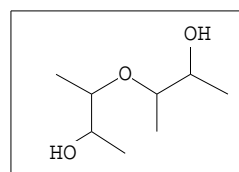
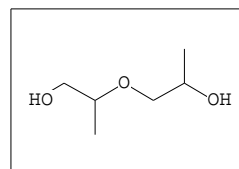
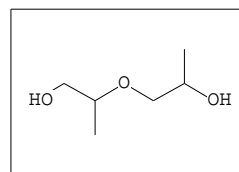
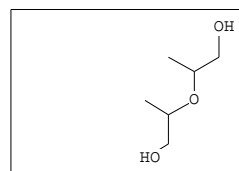
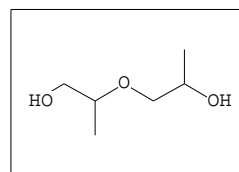
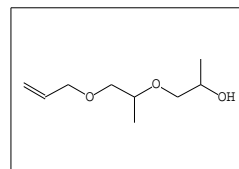
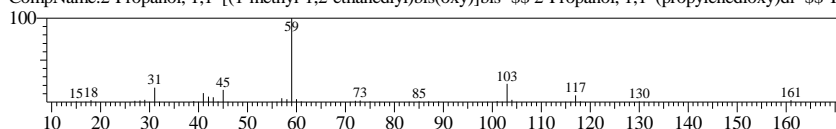
CompName:Dipropylene glycol monomethyl ether \$\$ Propanol, (2-methoxymethylethoxy)- \$\$ Propanol, 1(or 2)-(2-methoxyethoxy)-



Hit#8 Entry:37499 Library:NIST08.LIB

SI:89 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

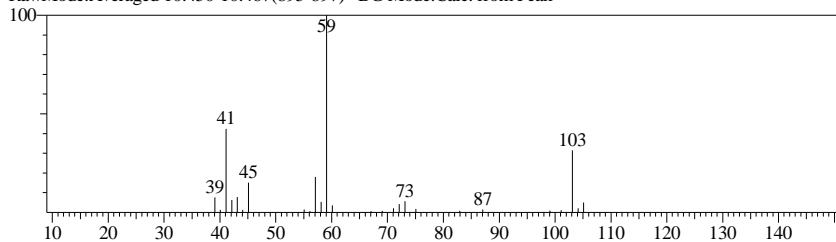
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



<< Target >>

Line# 2 RTime:10.458(Scan#:896) BasePeak:59.10(127392)

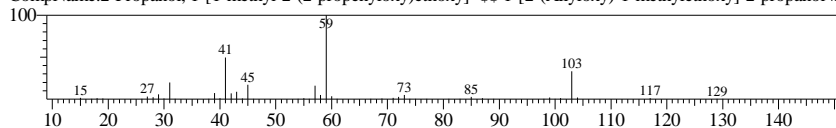
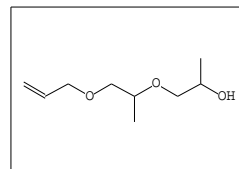
RawMode:Averaged 10.450-10.467(895-897) BG Mode:Calc. from Peak



Hit#:1 Entry:12792 Library:NIST08s.LIB

SI:96 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

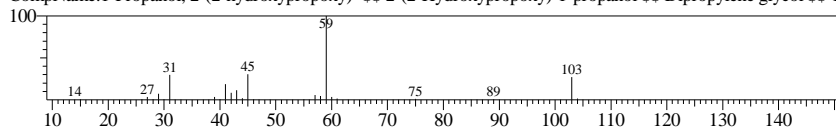
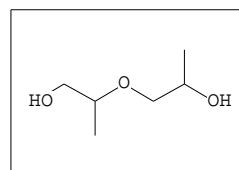
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:5921 Library:NIST08s.LIB

SI:90 Formula:C6H14O3 CAS:106-62-7 MolWeight:134 RetIndex:1034

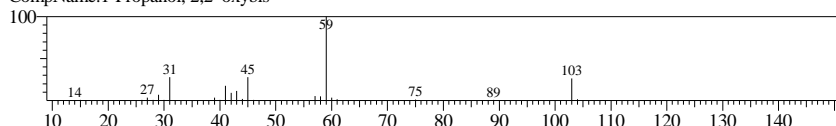
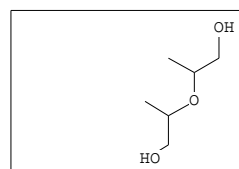
CompName:1-Propanol, 2-(2-hydroxypropoxy)- \$- 2-(2-Hydroxypropoxy)-1-propanol \$- Dipropylene glycol \$- 1



Hit#:3 Entry:8889 Library:NIST08.LIB

SI:90 Formula:C6H14O3 CAS:108-61-2 MolWeight:134 RetIndex:1051

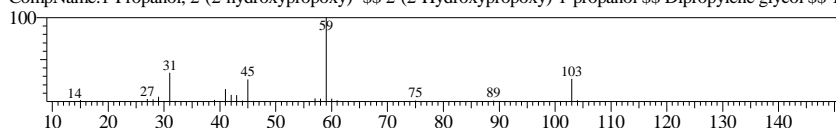
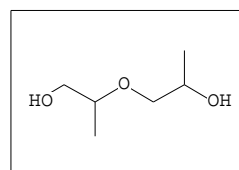
CompName:1-Propanol, 2,2'-oxybis-



Hit#:4 Entry:8891 Library:NIST08.LIB

SI:89 Formula:C6H14O3 CAS:106-62-7 MolWeight:134 RetIndex:1034

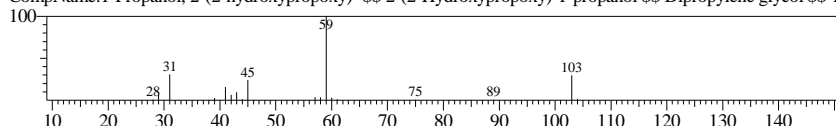
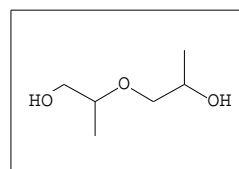
CompName:1-Propanol, 2-(2-hydroxypropoxy)- \$- 2-(2-Hydroxypropoxy)-1-propanol \$- Dipropylene glycol \$- 1



Hit#:5 Entry:5920 Library:NIST08s.LIB

SI:89 Formula:C6H14O3 CAS:106-62-7 MolWeight:134 RetIndex:1034

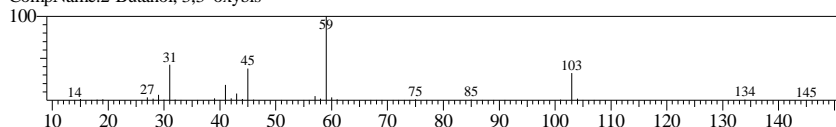
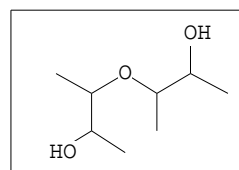
CompName:1-Propanol, 2-(2-hydroxypropoxy)- \$- 2-(2-Hydroxypropoxy)-1-propanol \$- Dipropylene glycol \$- 1



Hit#:6 Entry:20866 Library:NIST08.LIB

SI:89 Formula:C8H18O3 CAS:54305-61-2 MolWeight:162 RetIndex:1089

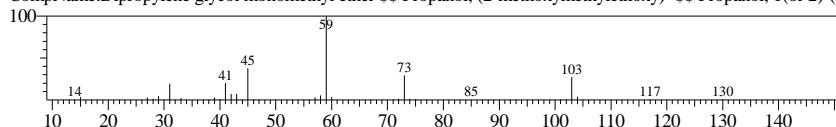
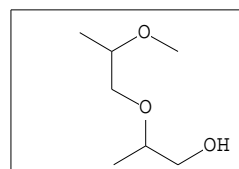
CompName:2-Butanol, 3,3'-oxybis-



Hit#:7 Entry:14295 Library:NIST08.LIB

SI:88 Formula:C7H16O3 CAS:34590-94-8 MolWeight:148 RetIndex:983

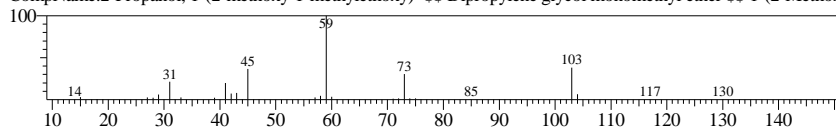
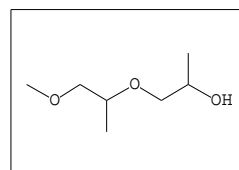
CompName:Dipropylene glycol monomethyl ether \$- Propanol, (2-methoxymethylethoxy)- \$- Propanol, 1-(or 2)-(2-methoxy-1-methylethoxy)-



Hit#:8 Entry:14298 Library:NIST08.LIB

SI:88 Formula:C7H16O3 CAS:20324-32-7 MolWeight:148 RetIndex:967

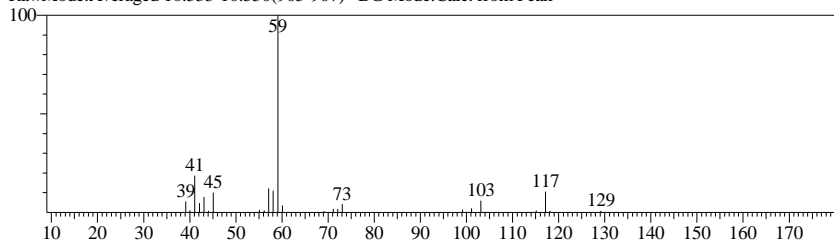
CompName:2-Propanol, 1-(2-methoxy-1-methylethoxy)- \$- Dipropylene glycol monomethyl ether \$- 1-(2-Methoxy-1-methylethoxy)-2-propanol



<< Target >>

Line#3 RTime:10.542(Scan#:906) BasePeak:59.10(213173)

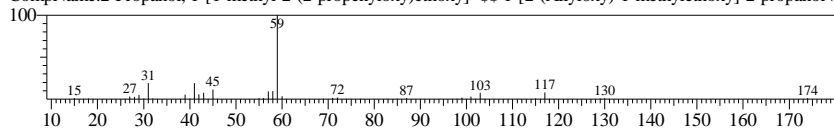
RawMode:Averaged 10.533-10.550(905-907) BG Mode:Calc. from Peak



Hit#1 Entry:27550 Library:NIST08.LIB

SI:95 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

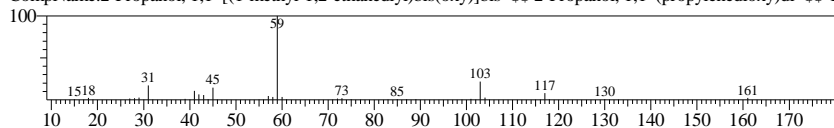
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#2 Entry:37499 Library:NIST08.LIB

SI:89 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

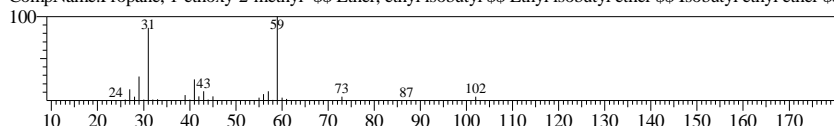
CompName:2-Propanol, 1,1'-[(1-methyl-2-ethanedyl)bis(oxy)]bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#3 Entry:2087 Library:NIST08.LIB

SI:89 Formula:C6H14O CAS:627-02-1 MolWeight:102 RetIndex:629

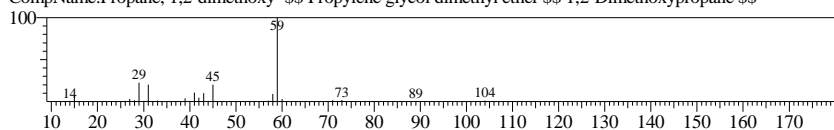
CompName:Propane, 1-ethoxy-2-methyl- \$- Ether, ethyl isobutyl \$- Ethyl isobutyl ether \$- Isobutyl ethyl ether \$-



Hit#4 Entry:2529 Library:NIST08.LIB

SI:89 Formula:C5H12O2 CAS:7778-85-0 MolWeight:104 RetIndex:606

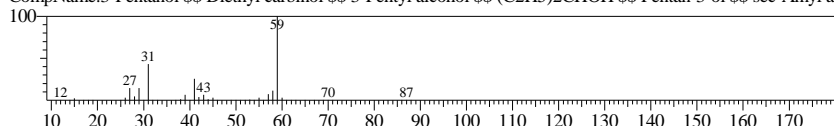
CompName:Propane, 1,2-dimethoxy- \$- Propylene glycol dimethyl ether \$- 1,2-Dimethoxypropane \$-



Hit#5 Entry:1044 Library:NIST08.LIB

SI:88 Formula:C5H12O CAS:584-02-1 MolWeight:88 RetIndex:681

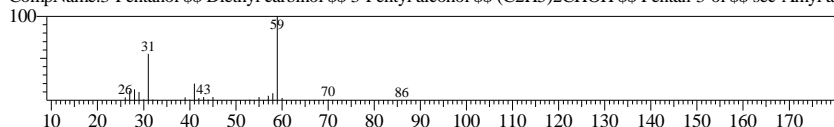
CompName:3-Pentanol \$- Diethyl carbinol \$- 3-Pentyl alcohol \$- (C2H5)2CHOH \$- Pentan-3-ol \$- sec-Amyl al



Hit#6 Entry:1068 Library:NIST08.LIB

SI:88 Formula:C5H12O CAS:584-02-1 MolWeight:88 RetIndex:681

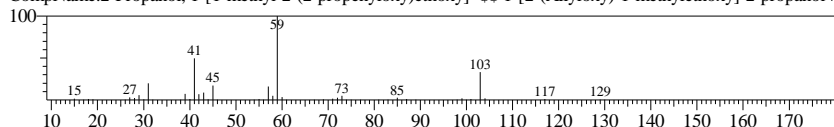
CompName:3-Pentanol \$- Diethyl carbinol \$- 3-Pentyl alcohol \$- (C2H5)2CHOH \$- Pentan-3-ol \$- sec-Amyl al



Hit#7 Entry:12792 Library:NIST08.LIB

SI:88 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

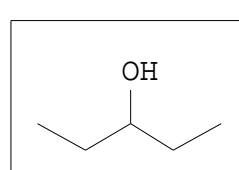
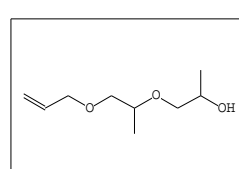
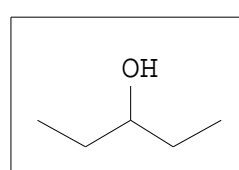
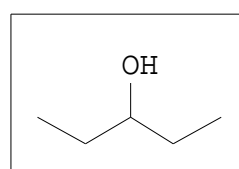
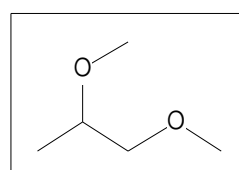
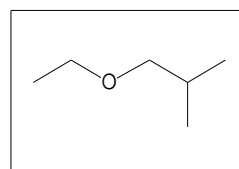
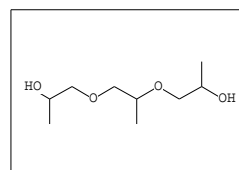
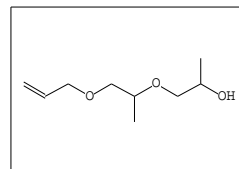
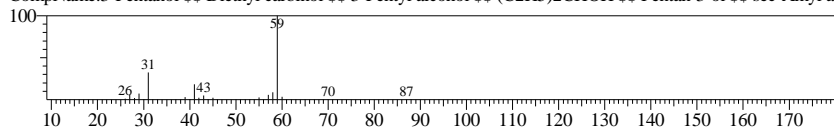
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#8 Entry:1070 Library:NIST08.LIB

SI:88 Formula:C5H12O CAS:584-02-1 MolWeight:88 RetIndex:681

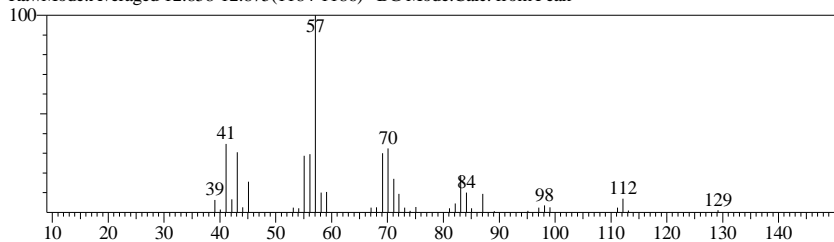
CompName:3-Pentanol \$- Diethyl carbinol \$- 3-Pentyl alcohol \$- (C2H5)2CHOH \$- Pentan-3-ol \$- sec-Amyl al



<< Target >>

Line#:4 R.Time:12.867(Scan#:1185) BasePeak:57.10(523067)

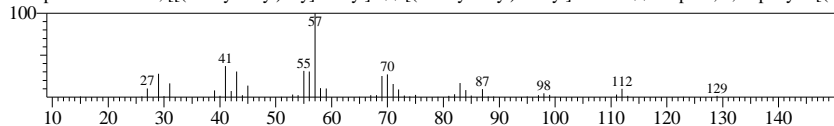
RawMode:Averaged 12.858-12.875(1184-1186) BG Mode:Calc. from Peak



Hit#:1 Entry:34522 Library:NIST08.LIB

SI:98 Formula:C11H22O2 CAS:2461-15-6 MolWeight:186 RetIndex:1217

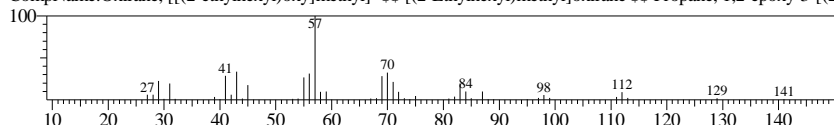
CompName:Oxirane, [[(2-ethylhexyl)oxy]methyl]- \$\$ [(2-Ethylhexyl)methyl]oxirane \$\$ Propane, 1,2-epoxy-3-[(2-



Hit#:2 Entry:14651 Library:NIST08s.LIB

SI:97 Formula:C11H22O2 CAS:2461-15-6 MolWeight:186 RetIndex:1217

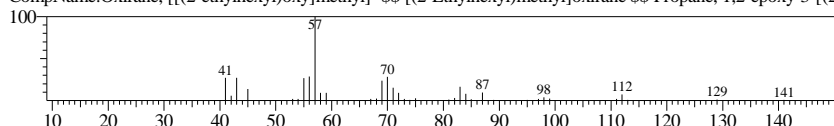
CompName:Oxirane, [[(2-ethylhexyl)oxy]methyl]- \$\$ [(2-Ethylhexyl)methyl]oxirane \$\$ Propane, 1,2-epoxy-3-[(2-



Hit#:3 Entry:14652 Library:NIST08s.LIB

SI:96 Formula:C11H22O2 CAS:2461-15-6 MolWeight:186 RetIndex:1217

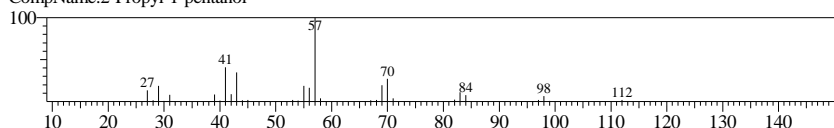
CompName:Oxirane, [[(2-ethylhexyl)oxy]methyl]- \$\$ [(2-Ethylhexyl)methyl]oxirane \$\$ Propane, 1,2-epoxy-3-[(2-



Hit#:4 Entry:8011 Library:NIST08.LIB

SI:89 Formula:C8H18O CAS:58175-57-8 MolWeight:130 RetIndex:995

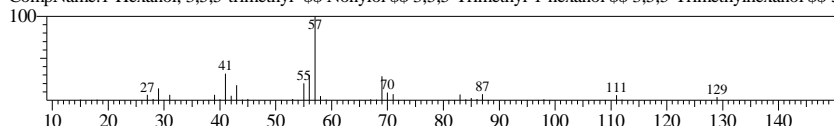
CompName:2-Propyl-1-pentanol



Hit#:5 Entry:13056 Library:NIST08.LIB

SI:89 Formula:C9H20O CAS:3452-97-9 MolWeight:144 RetIndex:1010

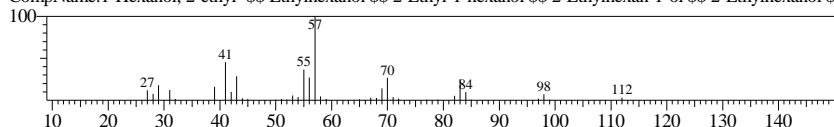
CompName:1-Hexanol, 3,5,5-trimethyl- \$\$ Nonyl-ol \$\$ 3,5,5-Trimethyl-1-hexanol \$\$ 3,5,5-Trimethylhexanol \$\$ 3



Hit#:6 Entry:5471 Library:NIST08.LIB

SI:89 Formula:C8H18O CAS:104-76-7 MolWeight:130 RetIndex:995

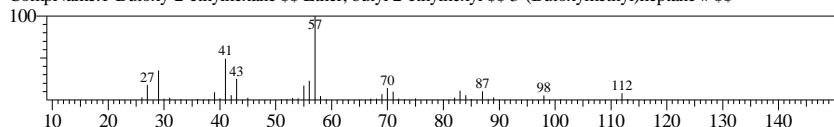
CompName:1-Hexanol, 2-ethyl- \$\$ Ethylhexanol \$\$ 2-Ethyl-1-hexanol \$\$ 2-Ethylhexanol-1-ol \$\$ 2-Ethylhexanol \$



Hit#:7 Entry:34667 Library:NIST08.LIB

SI:89 Formula:C12H26O CAS:62625-25-6 MolWeight:186 RetIndex:1226

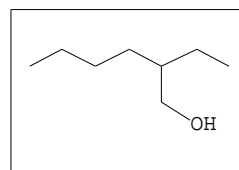
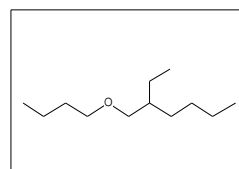
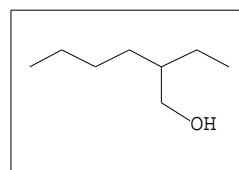
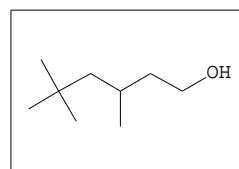
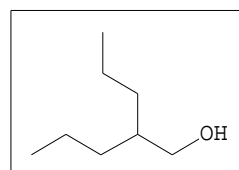
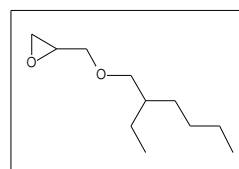
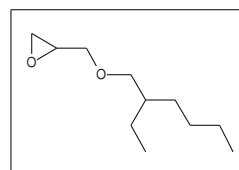
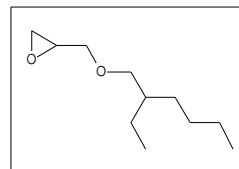
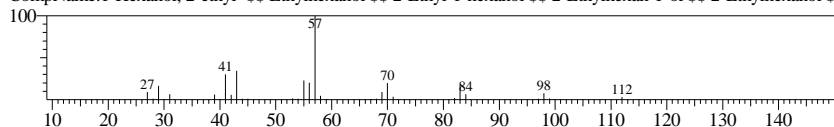
CompName:1-Butoxy-2-ethylhexane \$\$ Ether, butyl 2-ethylhexyl \$\$ 3-(Butoxymethyl)heptane # \$\$



Hit#:8 Entry:5475 Library:NIST08.LIB

SI:88 Formula:C8H18O CAS:104-76-7 MolWeight:130 RetIndex:995

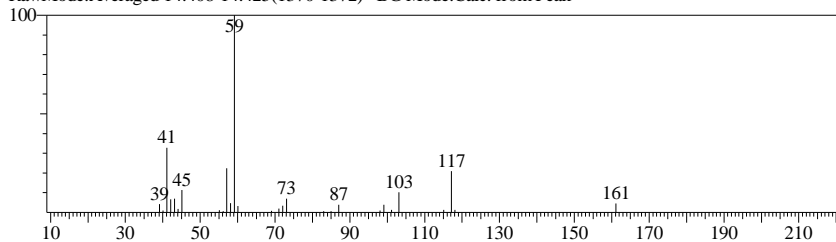
CompName:1-Hexanol, 2-ethyl- \$\$ Ethylhexanol \$\$ 2-Ethyl-1-hexanol \$\$ 2-Ethylhexanol-1-ol \$\$ 2-Ethylhexanol \$



<< Target >>

Line# 5 RTime:14.417(Scan#:1371) BasePeak:59.10(209944)

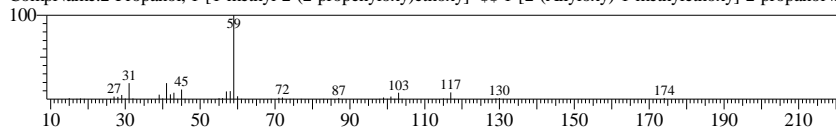
RawMode:Averaged 14.408-14.425(1370-1372) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:90 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

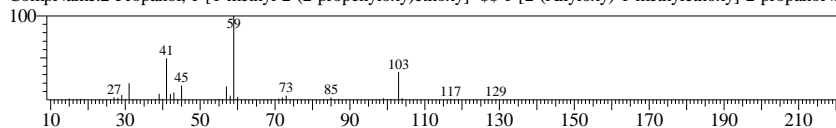
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:12792 Library:NIST08s.LIB

SI:89 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

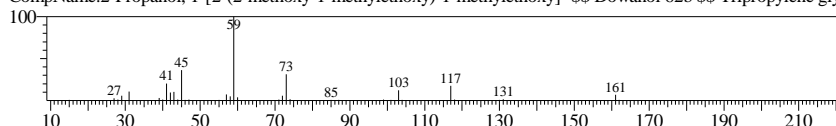
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:17306 Library:NIST08s.LIB

SI:88 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

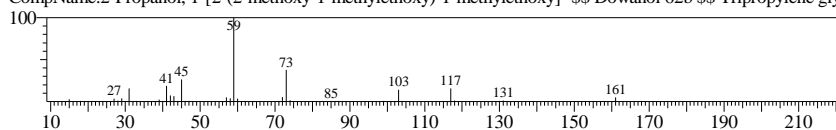
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#:4 Entry:46317 Library:NIST08.LIB

SI:88 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

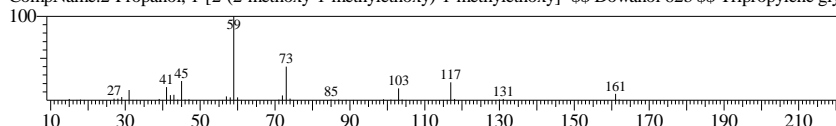
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#:5 Entry:17307 Library:NIST08s.LIB

SI:88 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

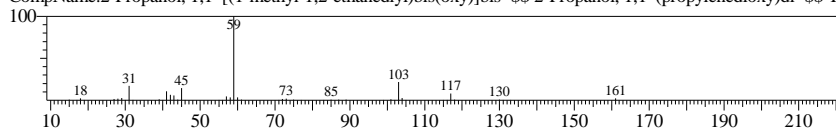
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#:6 Entry:37499 Library:NIST08.LIB

SI:88 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

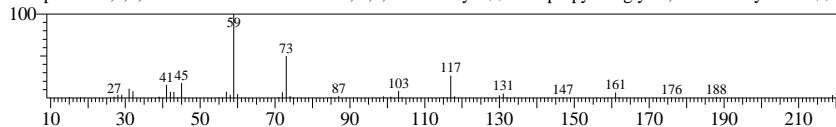
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#:7 Entry:85908 Library:NIST08.LIB

SI:86 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

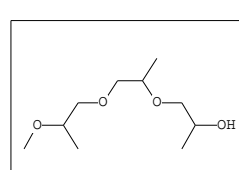
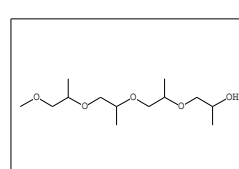
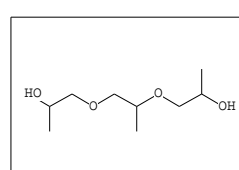
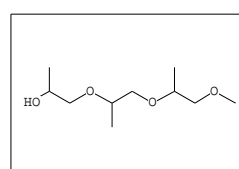
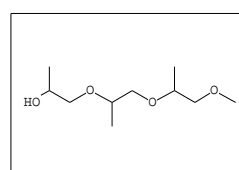
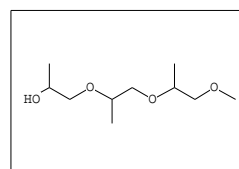
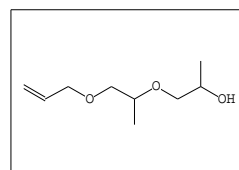
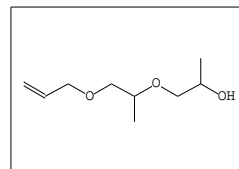
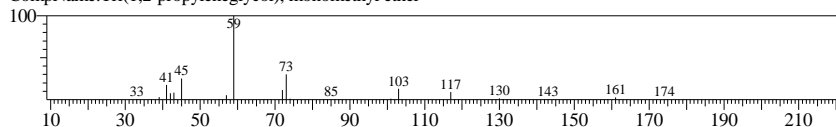
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$- Tetrapropylene glycol, monomethyl ether \$- 4



Hit#:8 Entry:46316 Library:NIST08.LIB

SI:85 Formula:C10H22O4 CAS:0-00-0 MolWeight:206 RetIndex:1277

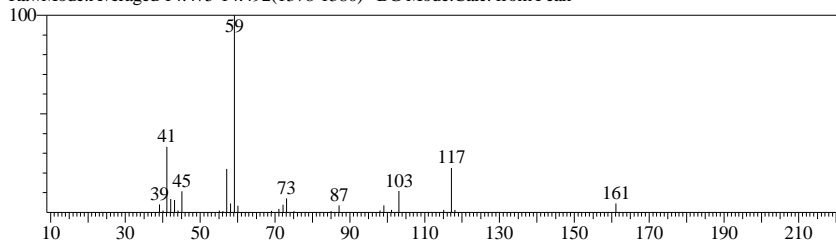
CompName:Tri(1,2-propyleneglycol), monomethyl ether



<< Target >>

Line# 6 RTime:14.483(Scan#:1379) BasePeak:59.10(440649)

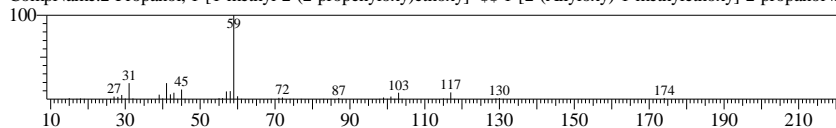
RawMode:Averaged 14.475-14.492(1378-1380) BG Mode:Calc. from Peak



Hit#1 Entry:27550 Library:NIST08.LIB

SI:89 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

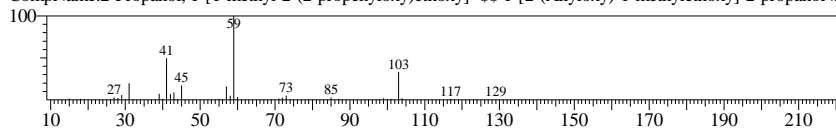
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#2 Entry:12792 Library:NIST08s.LIB

SI:89 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

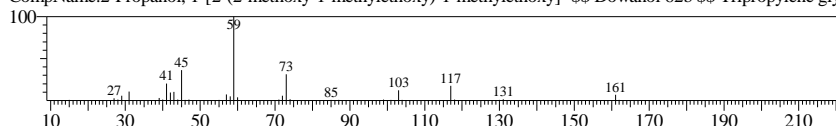
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#3 Entry:17306 Library:NIST08s.LIB

SI:89 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

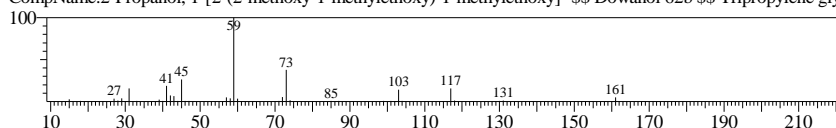
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#4 Entry:46317 Library:NIST08.LIB

SI:89 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

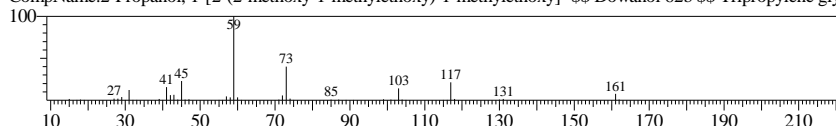
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#5 Entry:17307 Library:NIST08s.LIB

SI:88 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

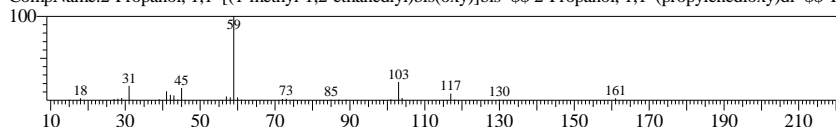
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#6 Entry:37499 Library:NIST08.LIB

SI:88 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

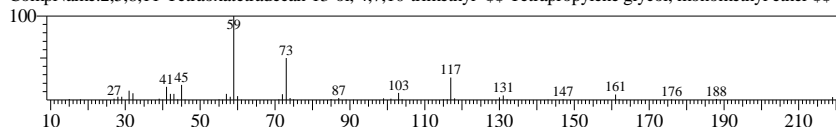
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#7 Entry:85908 Library:NIST08.LIB

SI:86 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

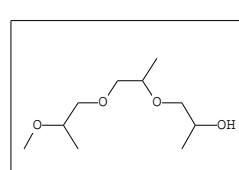
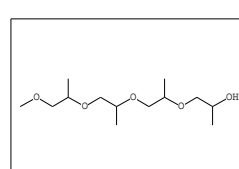
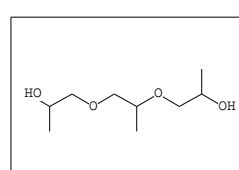
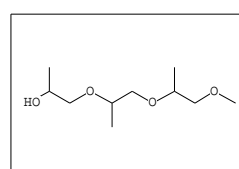
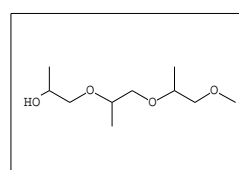
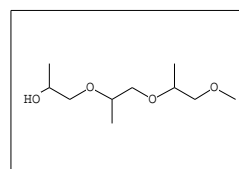
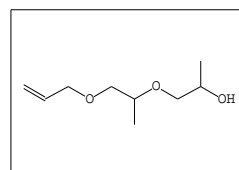
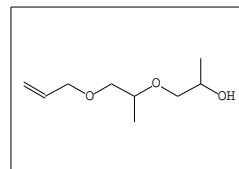
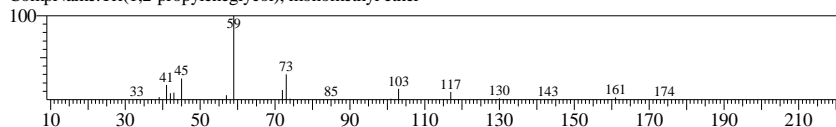
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$- Tetrapropylene glycol, monomethyl ether \$- 4



Hit#8 Entry:46316 Library:NIST08.LIB

SI:85 Formula:C10H22O4 CAS:0-00-0 MolWeight:206 RetIndex:1277

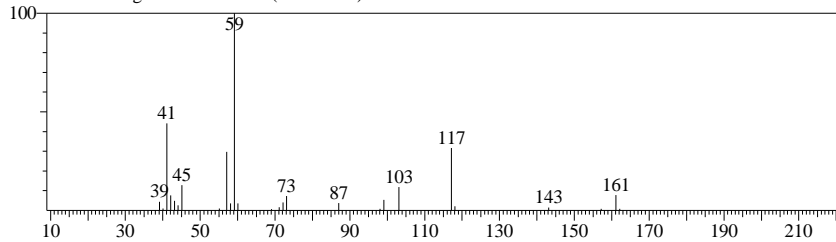
CompName:Tri(1,2-propyleneglycol), monomethyl ether



<< Target >>

Line# 7 RTime:14.542(Scan#:1386) BasePeak:59.10(101326)

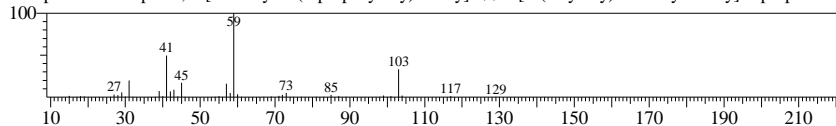
RawMode:Averaged 14.533-14.550(1385-1387) BG Mode:Calc. from Peak



Hit#1 Entry:12792 Library:NIST08s.LIB

SI:87 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

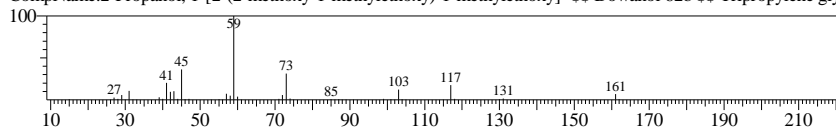
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#2 Entry:17306 Library:NIST08s.LIB

SI:86 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

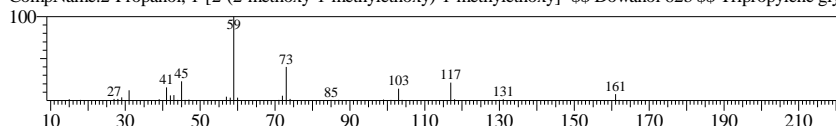
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#3 Entry:17307 Library:NIST08s.LIB

SI:86 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

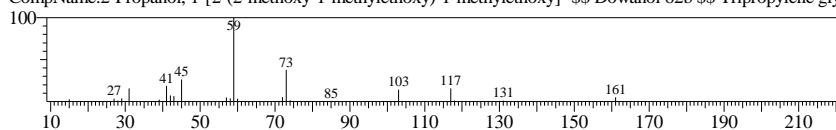
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#4 Entry:46317 Library:NIST08.LIB

SI:86 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

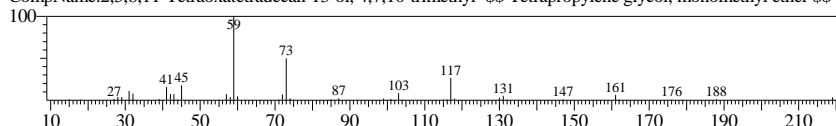
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#5 Entry:85908 Library:NIST08.LIB

SI:85 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

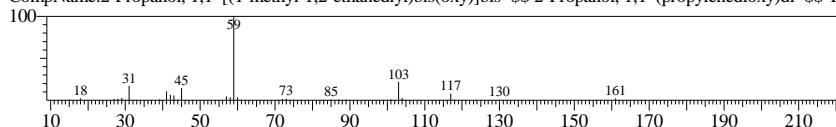
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$- Tetrapropylene glycol, monomethyl ether \$- 4



Hit#6 Entry:37499 Library:NIST08.LIB

SI:84 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

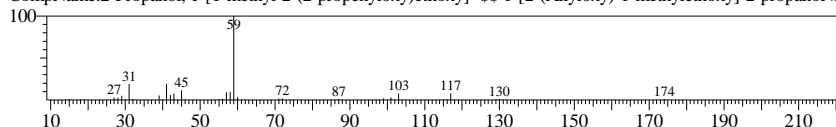
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#7 Entry:27550 Library:NIST08.LIB

SI:84 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

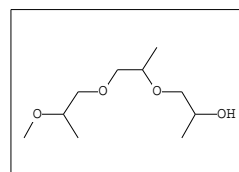
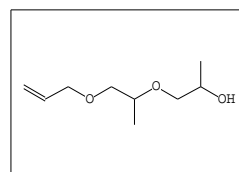
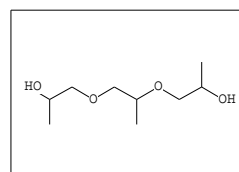
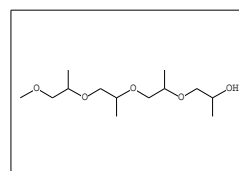
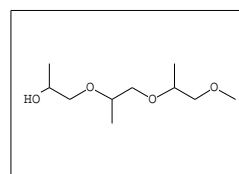
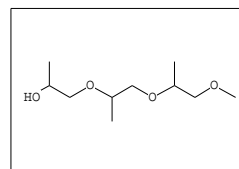
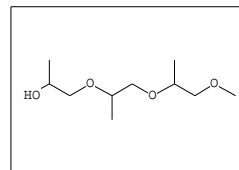
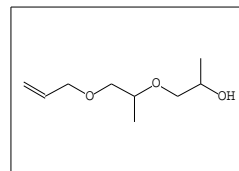
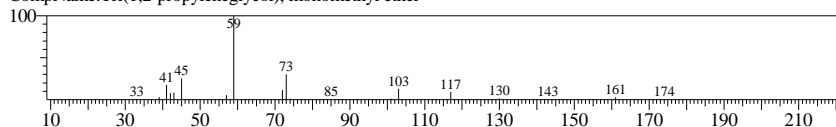
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#8 Entry:46316 Library:NIST08.LIB

SI:83 Formula:C10H22O4 CAS:0-00-0 MolWeight:206 RetIndex:1277

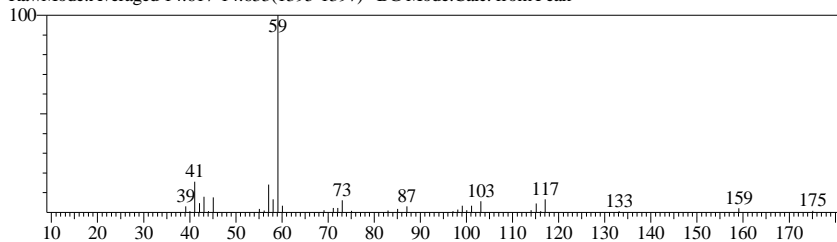
CompName:Tri(1,2-propyleneglycol), monomethyl ether



<< Target >>

Line# 8 RTime: 14.625 (Scan#: 1396) BasePeak: 59.10 (289372)

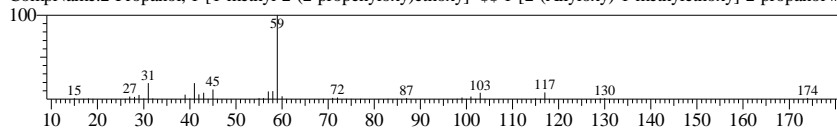
RawMode: Averaged 14.617-14.633 (1395-1397) BG Mode: Calc. from Peak



Hit#1 Entry: 27550 Library: NIST08.LIB

SI: 92 Formula: C₉H₁₈O₃ CAS: 55956-25-7 MolWeight: 174 RetIndex: 1156

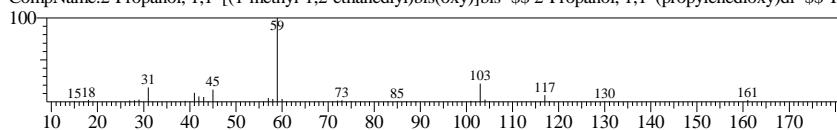
CompName: 2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#2 Entry: 37499 Library: NIST08.LIB

SI: 87 Formula: C₉H₂₀O₄ CAS: 1638-16-0 MolWeight: 192 RetIndex: 1328

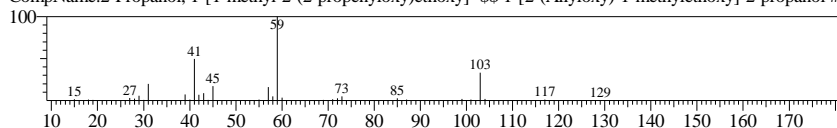
CompName: 2-Propanol, 1,1'-[(1-methyl-2-(2-ethanedyl)bis(oxy))bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#3 Entry: 12792 Library: NIST08.LIB

SI: 87 Formula: C₉H₁₈O₃ CAS: 55956-25-7 MolWeight: 174 RetIndex: 1156

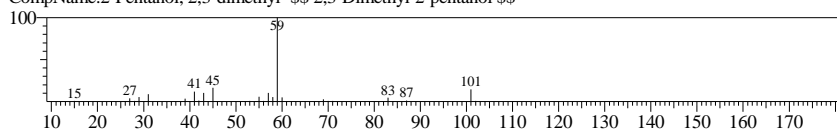
CompName: 2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry: 4678 Library: NIST08.LIB

SI: 86 Formula: C₇H₁₆O CAS: 4911-70-0 MolWeight: 116 RetIndex: 745

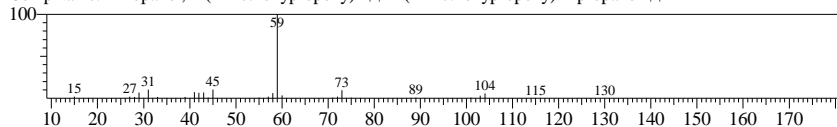
CompName: 2-Pentanol, 2,3-dimethyl- \$- 2,3-Dimethyl-2-pentanol \$-



Hit#5 Entry: 14294 Library: NIST08.LIB

SI: 86 Formula: C₇H₁₆O₃ CAS: 13429-07-7 MolWeight: 148 RetIndex: 967

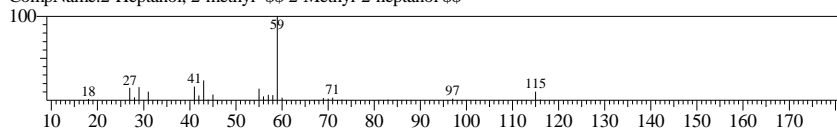
CompName: 2-Propanol, 1-(2-methoxypropoxy)- \$- 1-(2-Methoxypropoxy)-2-propanol \$-



Hit#6 Entry: 5482 Library: NIST08.LIB

SI: 86 Formula: C₈H₁₈O CAS: 625-25-2 MolWeight: 130 RetIndex: 908

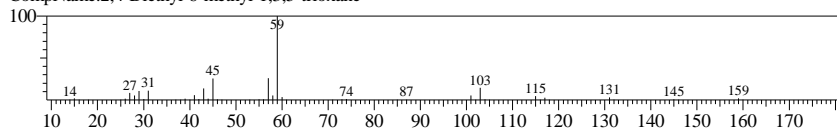
CompName: 2-Heptanol, 2-methyl- \$- 2-Methyl-2-heptanol \$-



Hit#7 Entry: 19968 Library: NIST08.LIB

SI: 86 Formula: C₈H₁₆O₃ CAS: 117888-04-7 MolWeight: 160 RetIndex: 1069

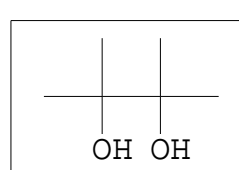
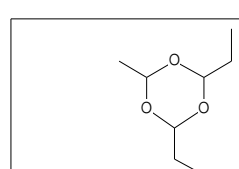
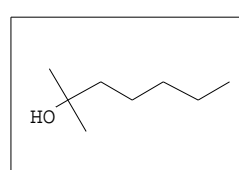
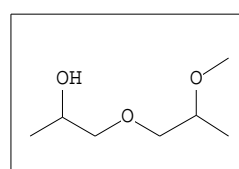
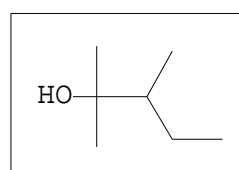
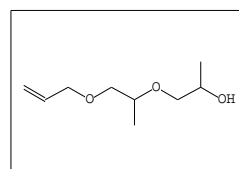
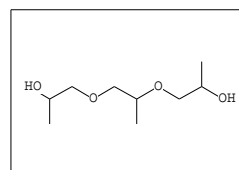
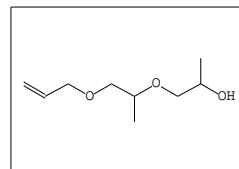
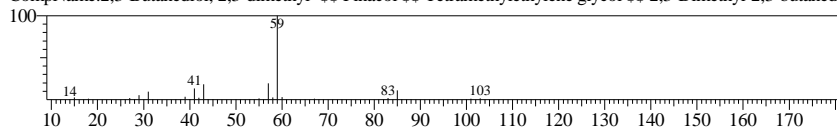
CompName: 2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#8 Entry: 3777 Library: NIST08.LIB

SI: 85 Formula: C₆H₁₄O₂ CAS: 76-09-5 MolWeight: 118 RetIndex: 801

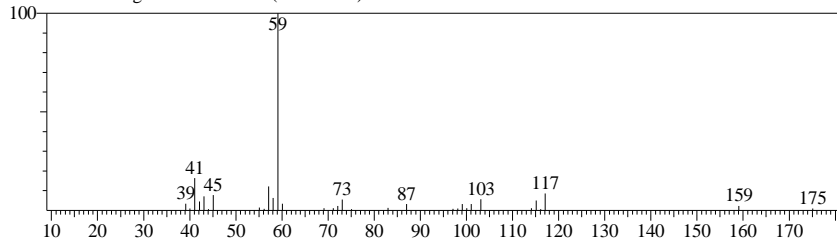
CompName: 2,3-Butanediol, 2,3-dimethyl- \$- Pinacol \$- Tetramethylethylene glycol \$- 2,3-Dimethyl-2,3-butanediol



<< Target >>

Line# 9 RTime: 14.675 (Scan#: 1402) BasePeak: 59.10 (191933)

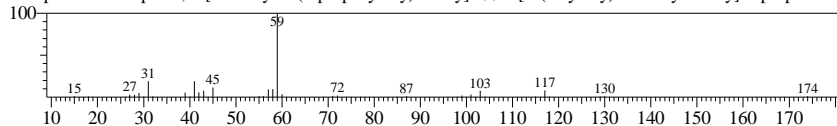
RawMode: Averaged 14.667-14.683 (1401-1403) BG Mode: Calc. from Peak



Hit#1 Entry: 27550 Library: NIST08.LIB

SI: 92 Formula: C₉H₁₈O₃ CAS: 55956-25-7 MolWeight: 174 RetIndex: 1156

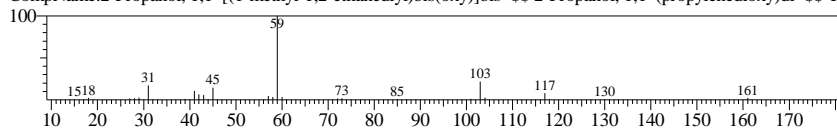
CompName: 2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#2 Entry: 37499 Library: NIST08.LIB

SI: 88 Formula: C₉H₂₀O₄ CAS: 1638-16-0 MolWeight: 192 RetIndex: 1328

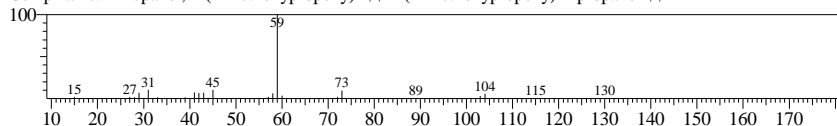
CompName: 2-Propanol, 1,1'-[(1-methyl-2-(2-ethanedyl)bis(oxy))bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#3 Entry: 14294 Library: NIST08.LIB

SI: 86 Formula: C₇H₁₆O₃ CAS: 13429-07-7 MolWeight: 148 RetIndex: 967

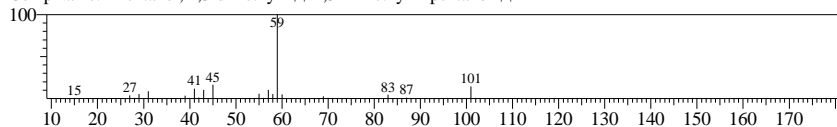
CompName: 2-Propanol, 1-(2-methoxypropoxy)- \$- 1-(2-Methoxypropoxy)-2-propanol \$-



Hit#4 Entry: 4678 Library: NIST08.LIB

SI: 86 Formula: C₇H₁₆O CAS: 4911-70-0 MolWeight: 116 RetIndex: 745

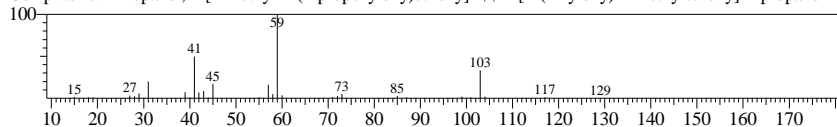
CompName: 2-Pentanol, 2,3-dimethyl- \$- 2,3-Dimethyl-2-pentanol \$-



Hit#5 Entry: 12792 Library: NIST08s.LIB

SI: 86 Formula: C₉H₁₈O₃ CAS: 55956-25-7 MolWeight: 174 RetIndex: 1156

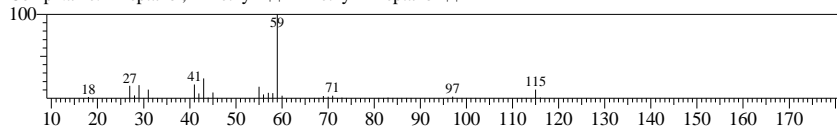
CompName: 2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#6 Entry: 5482 Library: NIST08s.LIB

SI: 86 Formula: C₈H₁₈O CAS: 625-25-2 MolWeight: 130 RetIndex: 908

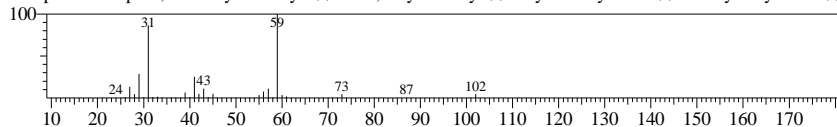
CompName: 2-Heptanol, 2-methyl- \$- 2-Methyl-2-heptanol \$-



Hit#7 Entry: 2087 Library: NIST08s.LIB

SI: 85 Formula: C₆H₁₄O CAS: 627-02-1 MolWeight: 102 RetIndex: 629

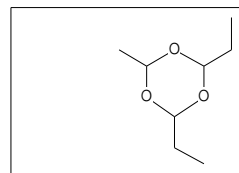
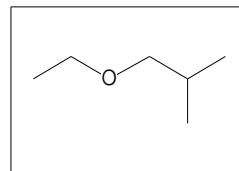
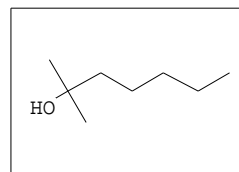
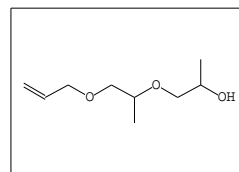
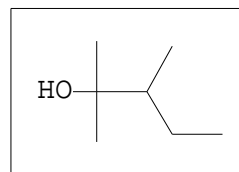
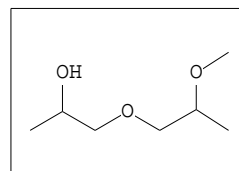
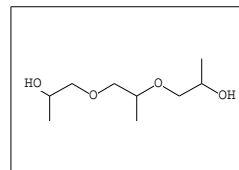
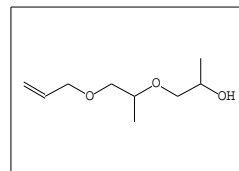
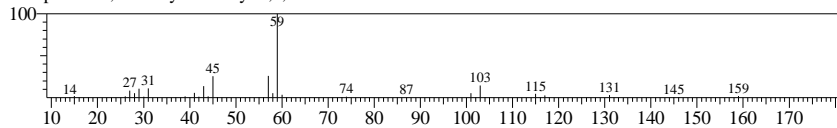
CompName: Propane, 1-ethoxy-2-methyl- \$- Ether, ethyl isobutyl \$- Ethyl isobutyl ether \$- Isobutyl ethyl ether \$-



Hit#8 Entry: 19968 Library: NIST08.LIB

SI: 85 Formula: C₈H₁₆O₃ CAS: 117888-04-7 MolWeight: 160 RetIndex: 1069

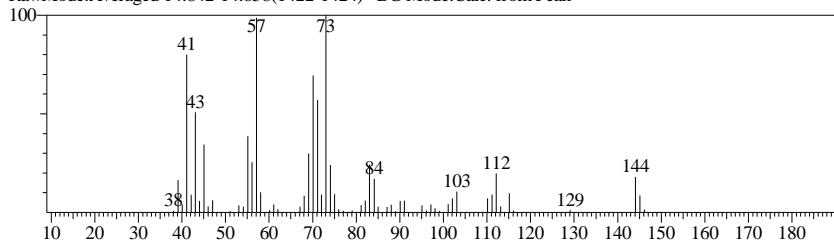
CompName: 2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line# 10 R.Time:14.850(Scan#:1423) BasePeak:73.05(51588)

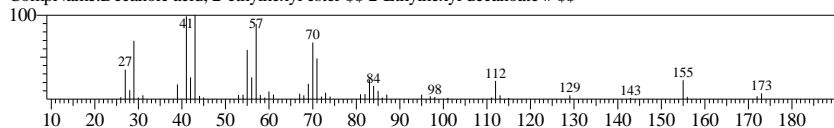
RawMode:Averaged 14.842-14.858(1422-1424) BG Mode:Calc. from Peak



Hit#1 Entry:100788 Library:NIST08.LIB

SI:81 Formula:C18H36O2 CAS:73947-30-5 MolWeight:284 RetIndex:1914

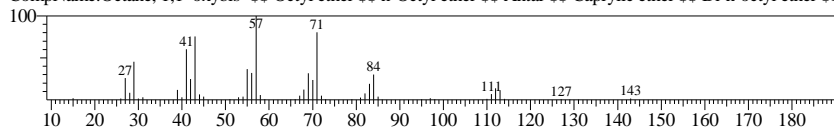
CompName:Decanoic acid, 2-ethylhexyl ester \$ 2-Ethylhexyl decanoate # \$



Hit#2 Entry:20865 Library:NIST08.LIB

SI:80 Formula:C16H34O CAS:629-82-3 MolWeight:242 RetIndex:1688

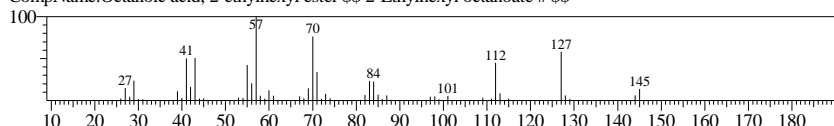
CompName:Octane, 1,1'-oxybis- \$ Octyl ether \$ \$ Antar \$ \$ Caprylic ether \$ \$ Di-n-octyl ether \$ \$



Hit#3 Entry:80752 Library:NIST08.LIB

SI:79 Formula:C16H32O2 CAS:63321-70-0 MolWeight:256 RetIndex:1715

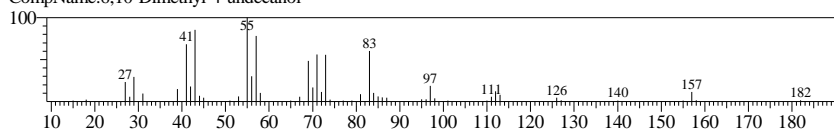
CompName:Octanoic acid, 2-ethylhexyl ester \$ 2-Ethylhexyl octanoate # \$



Hit#4 Entry:43040 Library:NIST08.LIB

SI:79 Formula:C13H28O CAS:92153-95-2 MolWeight:200 RetIndex:1348

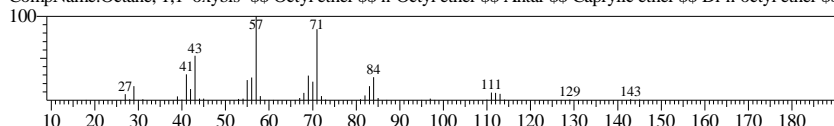
CompName:6,10-Dimethyl-4-undecanol



Hit#5 Entry:20866 Library:NIST08.LIB

SI:78 Formula:C16H34O CAS:629-82-3 MolWeight:242 RetIndex:1688

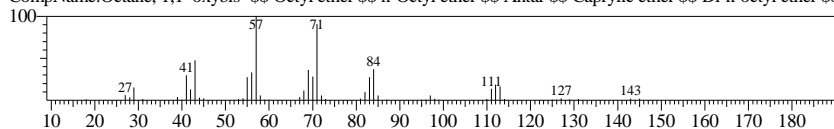
CompName:Octane, 1,1'-oxybis- \$ Octyl ether \$ \$ Antar \$ \$ Caprylic ether \$ \$ Di-n-octyl ether \$ \$



Hit#6 Entry:20864 Library:NIST08.LIB

SI:78 Formula:C16H34O CAS:629-82-3 MolWeight:242 RetIndex:1688

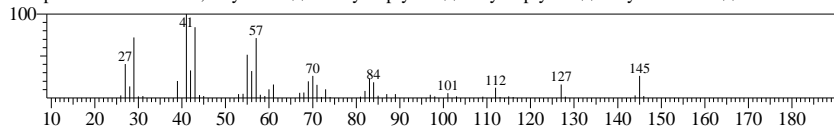
CompName:Octane, 1,1'-oxybis- \$ Octyl ether \$ \$ Antar \$ \$ Caprylic ether \$ \$ Di-n-octyl ether \$ \$



Hit#7 Entry:21855 Library:NIST08.LIB

SI:78 Formula:C16H32O2 CAS:2306-88-9 MolWeight:256 RetIndex:1779

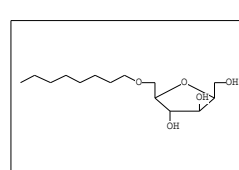
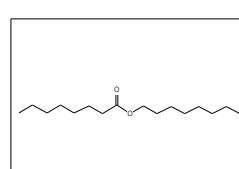
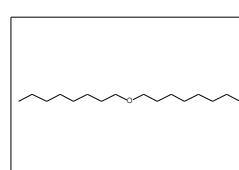
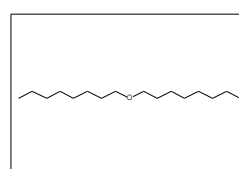
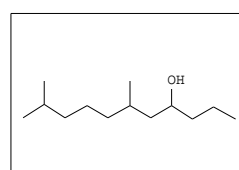
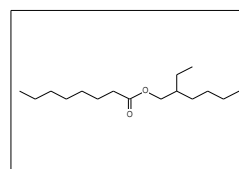
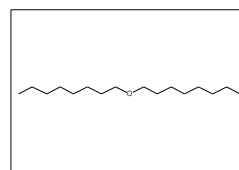
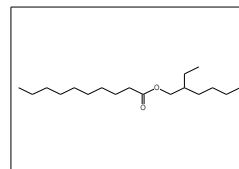
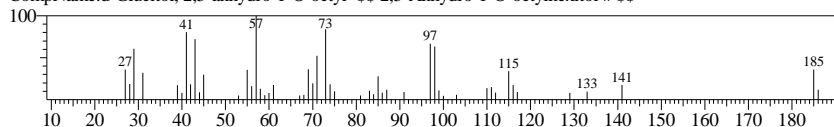
CompName:Octanoic acid, octyl ester \$ n-Octyl caprylate \$ \$ Octyl caprylate \$ \$ Octyl octanoate \$ \$



Hit#8 Entry:94736 Library:NIST08.LIB

SI:78 Formula:C14H28O5 CAS:101221-94-7 MolWeight:276 RetIndex:2201

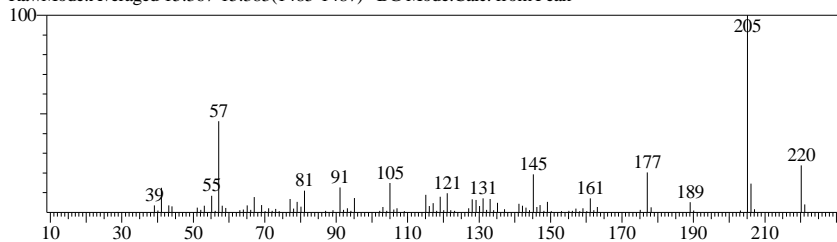
CompName:d-Glucitol, 2,5-anhydro-1-O-octyl- \$ 2,5-Anhydro-1-O-octylhexitol # \$



<< Target >>

Line# 11 R.Time:15.375(Scan#:1486) BasePeak:205.10(146191)

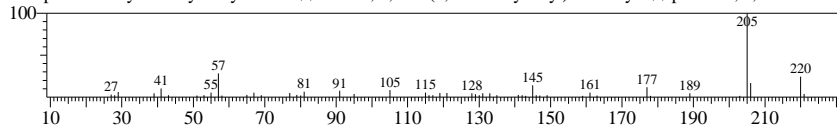
RawMode:Averaged 15.367-15.383(1485-1487) BG Mode:Calc. from Peak



Hit#:1 Entry:56018 Library:NIST08s.LIB

SI:92 Formula:C₁₅H₂₄O CAS:128-37-0 MolWeight:220 RetIndex:1668

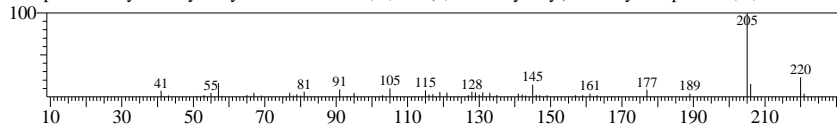
CompName:Butylated Hydroxytoluene \$\$ Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl- \$\$ p-Cresol, 2,6-di-tert-bu



Hit#:2 Entry:18880 Library:NIST08s.LIB

SI:90 Formula:C₁₅H₂₄O CAS:128-37-0 MolWeight:220 RetIndex:1668

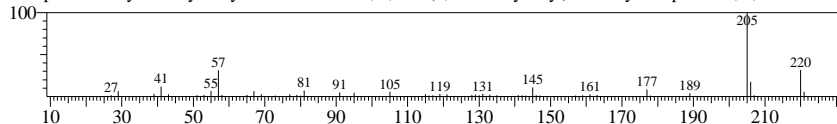
CompName:Butylated Hydroxytoluene \$\$ Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl- \$\$ p-Cresol, 2,6-di-tert-bu



Hit#:3 Entry:18879 Library:NIST08s.LIB

SI:88 Formula:C₁₅H₂₄O CAS:128-37-0 MolWeight:220 RetIndex:1668

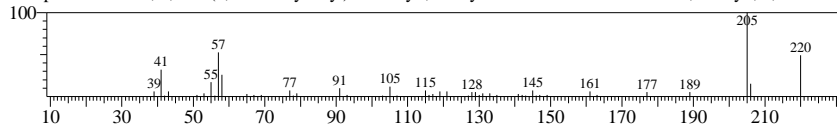
CompName:Butylated Hydroxytoluene \$\$ Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl- \$\$ p-Cresol, 2,6-di-tert-bu



Hit#:4 Entry:95568 Library:NIST08s.LIB

SI:86 Formula:C₁₇H₂₇NO₂ CAS:1918-11-2 MolWeight:277 RetIndex:2026

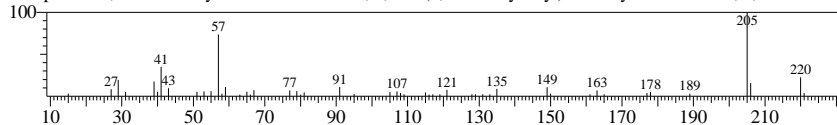
CompName:Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl-, methylcarbamate \$\$ Carbamic acid, methyl-, 2,6-di-tert-



Hit#:5 Entry:56016 Library:NIST08s.LIB

SI:81 Formula:C₁₅H₂₄O CAS:497-39-2 MolWeight:220 RetIndex:1668

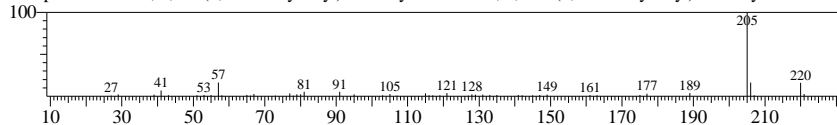
CompName:4,6-di-tert-Butyl-m-cresol \$\$ Phenol, 2,4-bis(1,1-dimethylethyl)-5-methyl- \$\$ m-Cresol, 4,6-di-tert-bu



Hit#:6 Entry:56017 Library:NIST08s.LIB

SI:81 Formula:C₁₅H₂₄O CAS:616-55-7 MolWeight:220 RetIndex:1668

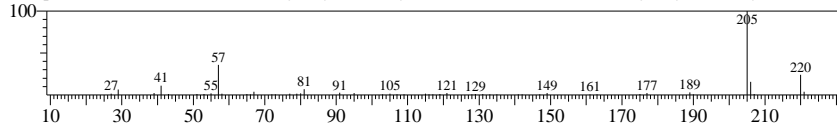
CompName:Phenol, 4,6-di(1,1-dimethylethyl)-2-methyl- \$\$ Phenol, 2,4-bis(1,1-dimethylethyl)-6-methyl- \$\$ 6-Me



Hit#:7 Entry:18878 Library:NIST08s.LIB

SI:81 Formula:C₁₅H₂₄O CAS:616-55-7 MolWeight:220 RetIndex:1668

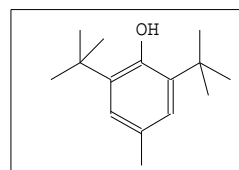
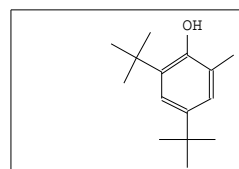
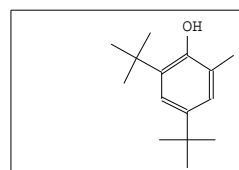
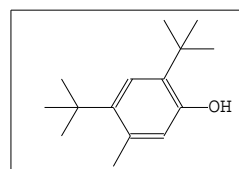
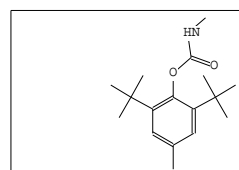
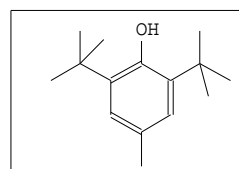
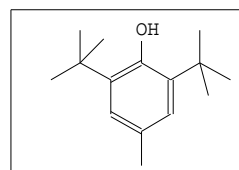
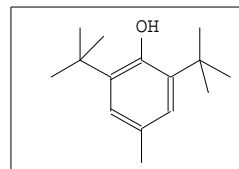
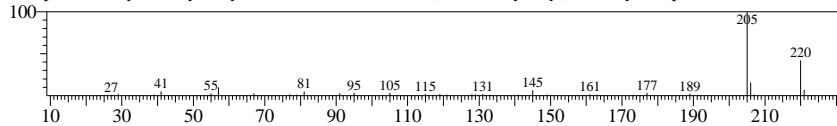
CompName:Phenol, 4,6-di(1,1-dimethylethyl)-2-methyl- \$\$ Phenol, 2,4-bis(1,1-dimethylethyl)-6-methyl- \$\$ 6-Me



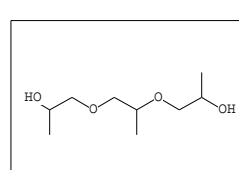
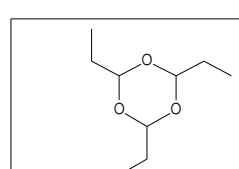
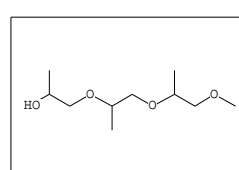
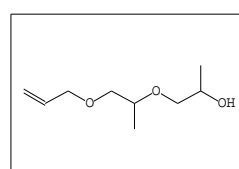
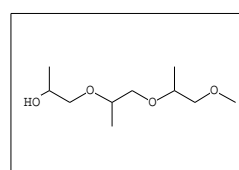
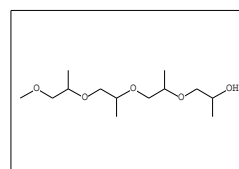
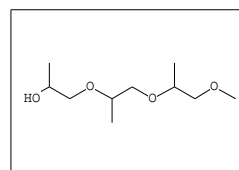
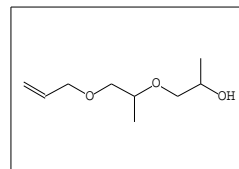
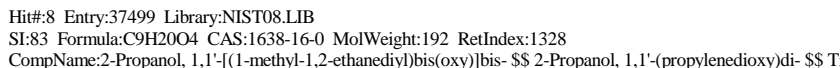
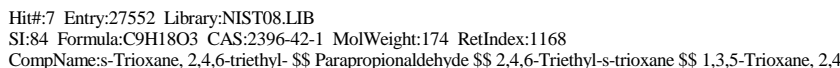
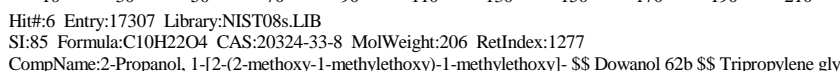
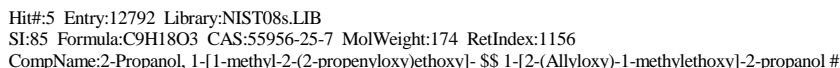
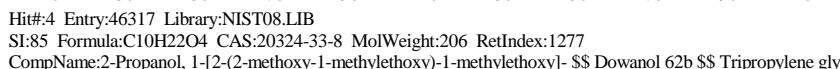
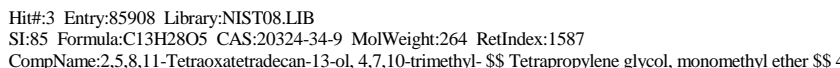
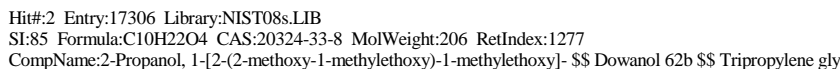
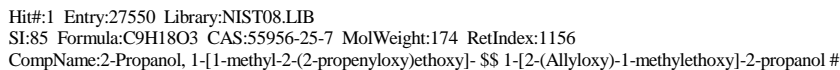
Hit#:8 Entry:18881 Library:NIST08s.LIB

SI:80 Formula:C₁₅H₂₄O CAS:128-37-0 MolWeight:220 RetIndex:1668

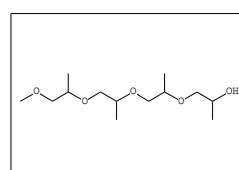
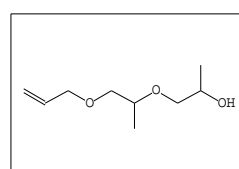
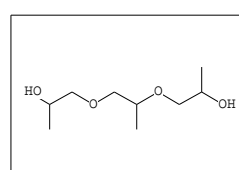
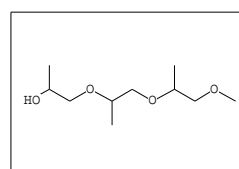
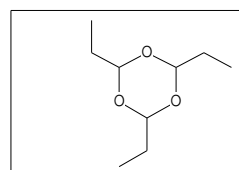
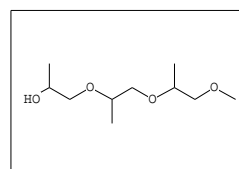
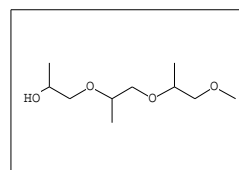
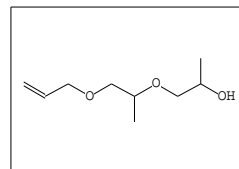
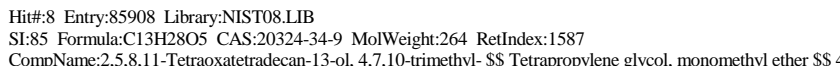
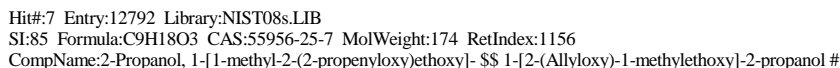
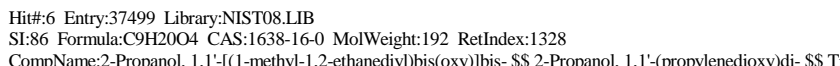
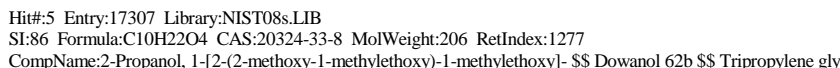
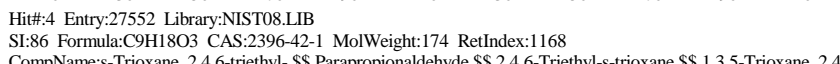
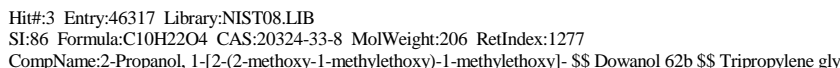
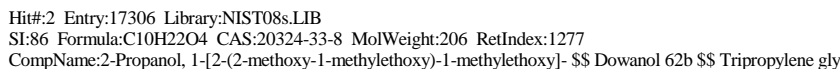
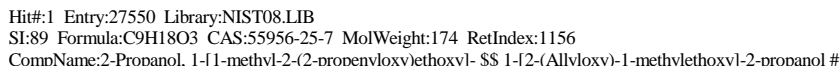
CompName:Butylated Hydroxytoluene \$\$ Phenol, 2,6-bis(1,1-dimethylethyl)-4-methyl- \$\$ p-Cresol, 2,6-di-tert-bu



Line#:12 R.Time:17.758(Scan#:1772) BasePeak:59.10(191095)
RawMode:Averaged 17.750-17.767(1771-1773) BG Mode:Calc. from Peak



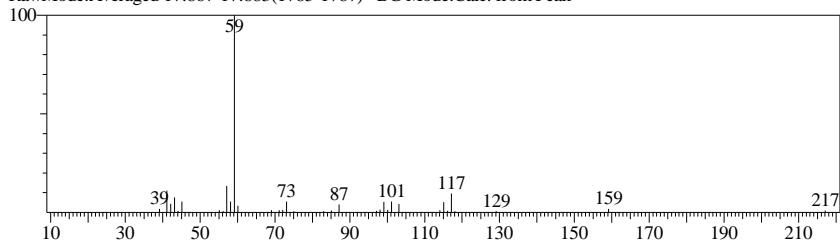
Line#:13 R.Time:17.817(Scan#:1779) BasePeak:59.10(326461)
RawMode:Averaged 17.808-17.825(1778-1780) BG Mode:Calc. from Peak



<< Target >>

Line#:14 R.Time:17.875(Scan#:1786) BasePeak:59.10(478069)

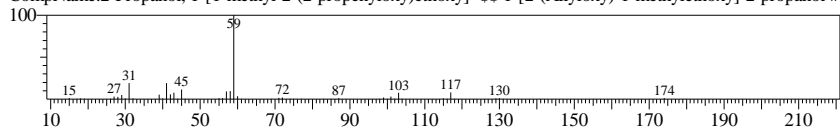
RawMode:Averaged 17.867-17.883(1785-1787) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:90 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

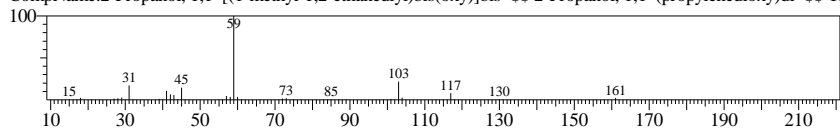
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:37499 Library:NIST08.LIB

SI:87 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

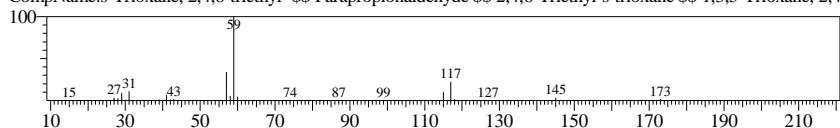
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:3 Entry:27552 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

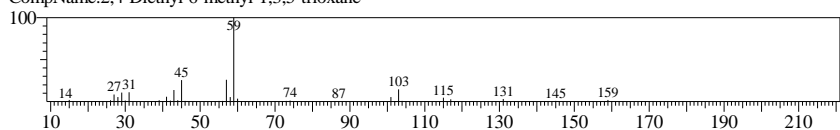
CompName:s-Trioxane, 2,4,6-triethyl- \$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:4 Entry:19968 Library:NIST08.LIB

SI:86 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

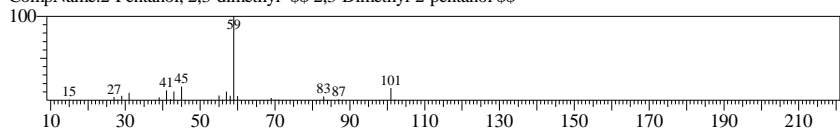
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:5 Entry:4678 Library:NIST08.LIB

SI:85 Formula:C7H16O CAS:4911-70-0 MolWeight:116 RetIndex:745

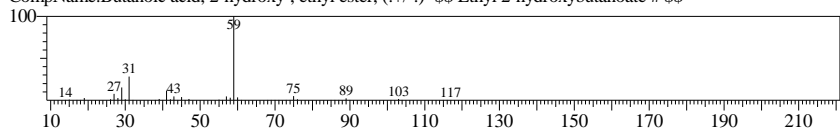
CompName:2-Pentanol, 2,3-dimethyl- \$ 2,3-Dimethyl-2-pentanol \$



Hit#:6 Entry:8423 Library:NIST08.LIB

SI:85 Formula:C6H12O3 CAS:68057-83-0 MolWeight:132 RetIndex:947

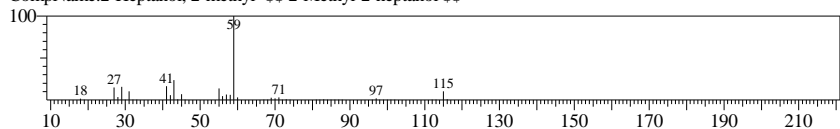
CompName:Butanoic acid, 2-hydroxy-, ethyl ester, (+/-)- \$ Ethyl 2-hydroxybutanoate # \$



Hit#:7 Entry:5482 Library:NIST08.LIB

SI:85 Formula:C8H18O CAS:625-25-2 MolWeight:130 RetIndex:908

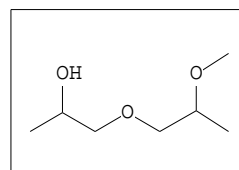
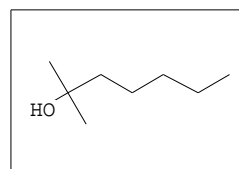
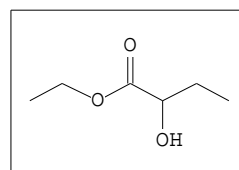
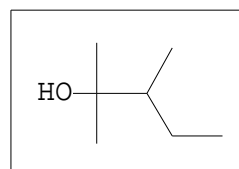
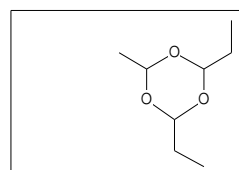
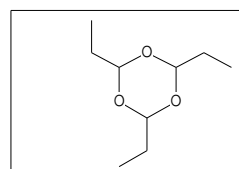
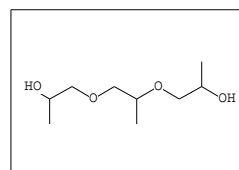
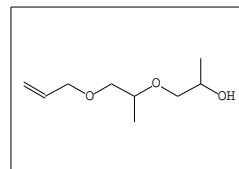
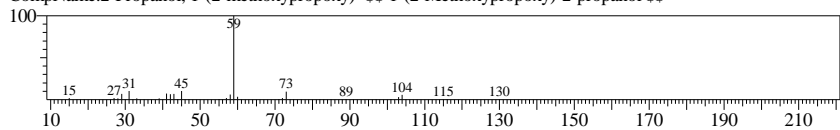
CompName:2-Heptanol, 2-methyl- \$ 2-Methyl-2-heptanol \$



Hit#:8 Entry:14294 Library:NIST08.LIB

SI:85 Formula:C7H16O3 CAS:13429-07-7 MolWeight:148 RetIndex:967

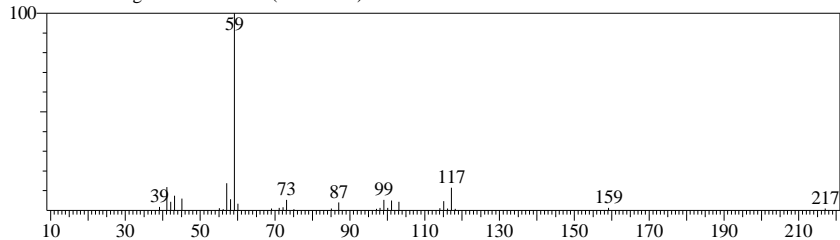
CompName:2-Propanol, 1-(2-methoxypropoxy)- \$ 1-(2-Methoxypropoxy)-2-propanol \$



<< Target >>

Line# 15 R.Time:17.933(Scan#:1793) BasePeak:59.10(488270)

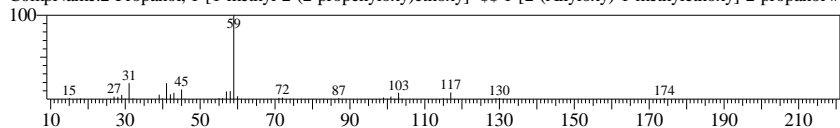
RawMode:Averaged 17.925-17.942(1792-1794) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:90 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

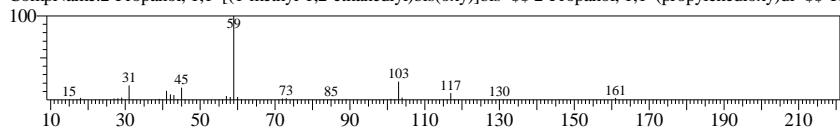
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:37499 Library:NIST08.LIB

SI:87 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

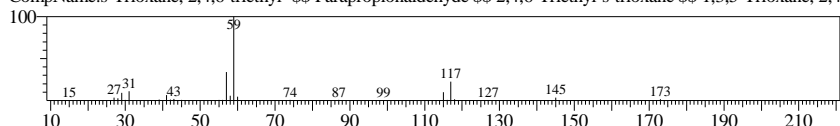
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:3 Entry:27552 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

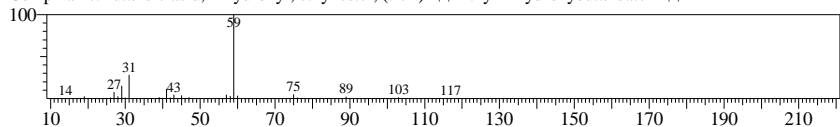
CompName:s-Trioxane, 2,4,6-triethyl- \$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:4 Entry:8423 Library:NIST08.LIB

SI:85 Formula:C6H12O3 CAS:68057-83-0 MolWeight:132 RetIndex:947

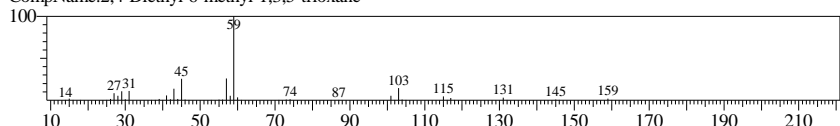
CompName:Butanoic acid, 2-hydroxy-, ethyl ester, (+/-)- \$ Ethyl 2-hydroxybutanoate # \$



Hit#:5 Entry:19968 Library:NIST08.LIB

SI:85 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

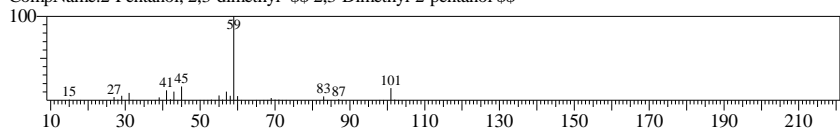
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:6 Entry:4678 Library:NIST08.LIB

SI:84 Formula:C7H16O CAS:4911-70-0 MolWeight:116 RetIndex:745

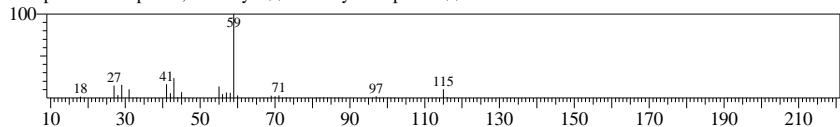
CompName:2-Pentanol, 2,3-dimethyl- \$ 2,3-Dimethyl-2-pentanol \$



Hit#:7 Entry:5482 Library:NIST08s.LIB

SI:84 Formula:C8H18O CAS:625-25-2 MolWeight:130 RetIndex:908

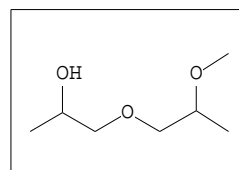
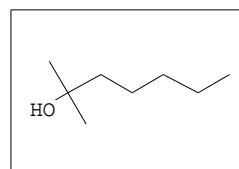
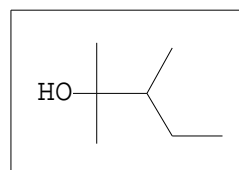
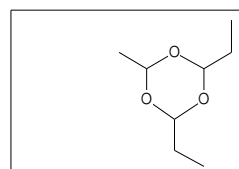
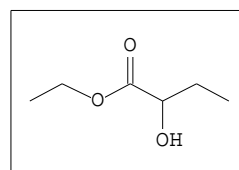
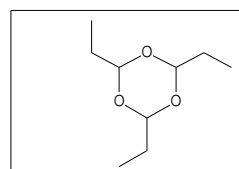
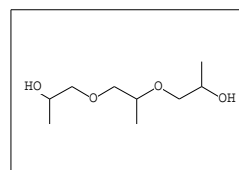
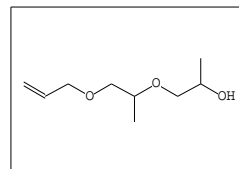
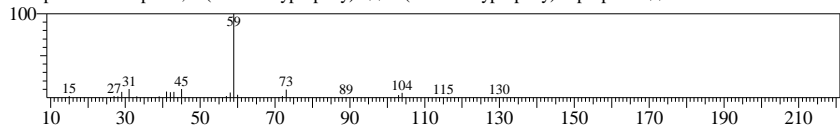
CompName:2-Heptanol, 2-methyl- \$ 2-Methyl-2-heptanol \$



Hit#:8 Entry:14294 Library:NIST08.LIB

SI:84 Formula:C7H16O3 CAS:13429-07-7 MolWeight:148 RetIndex:967

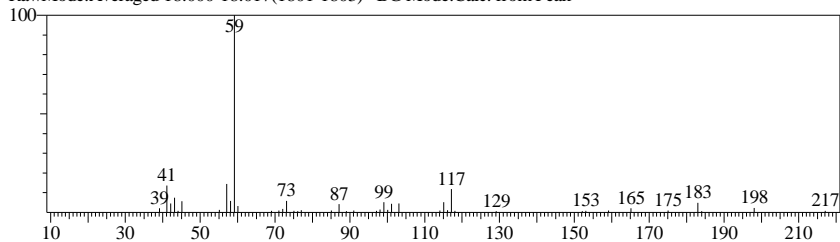
CompName:2-Propanol, 1-(2-methoxypropoxy)- \$ 1-(2-Methoxypropoxy)-2-propanol \$



<< Target >>

Line#:16 R.Time:18.008(Scan#:1802) BasePeak:59.10(259887)

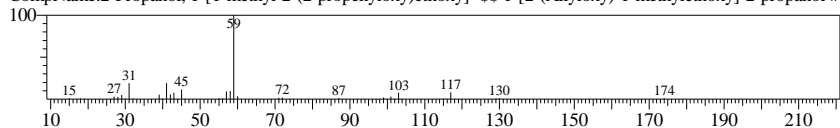
RawMode:Averaged 18.000-18.017(1801-1803) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:87 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

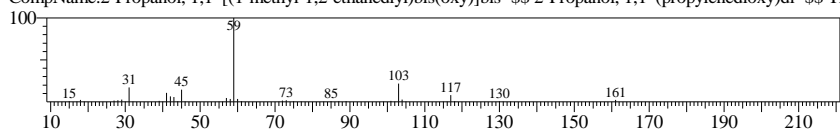
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:37499 Library:NIST08.LIB

SI:84 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

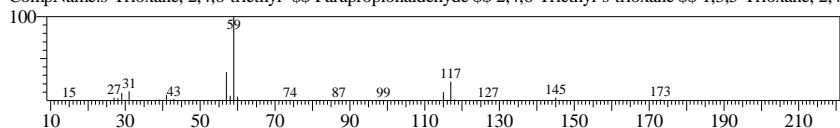
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:3 Entry:27552 Library:NIST08.LIB

SI:83 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

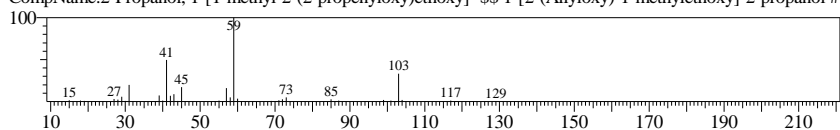
CompName:s-Trioxane, 2,4,6-triethyl- \$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:4 Entry:12792 Library:NIST08s.LIB

SI:83 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

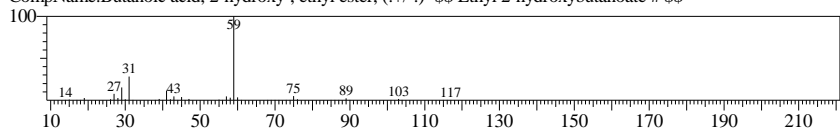
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:5 Entry:8423 Library:NIST08.LIB

SI:83 Formula:C6H12O3 CAS:68057-83-0 MolWeight:132 RetIndex:947

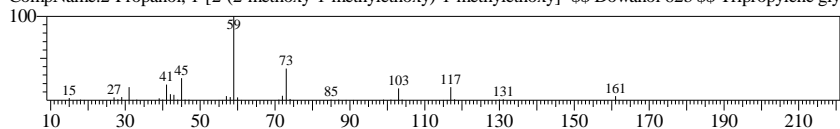
CompName:Butanoic acid, 2-hydroxy-, ethyl ester, (+/-)- \$ Ethyl 2-hydroxybutanoate # \$



Hit#:6 Entry:46317 Library:NIST08.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

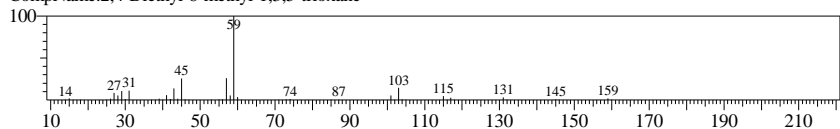
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



Hit#:7 Entry:19968 Library:NIST08.LIB

SI:82 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

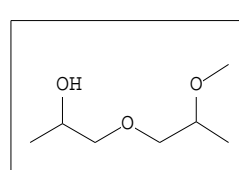
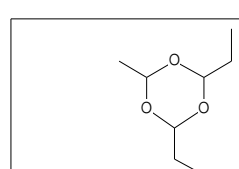
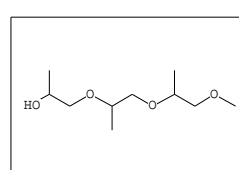
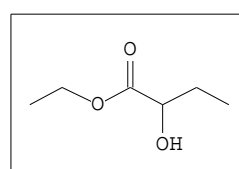
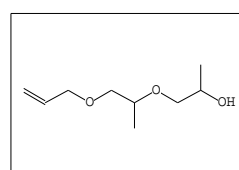
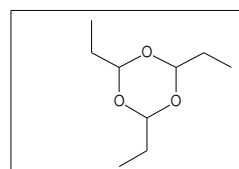
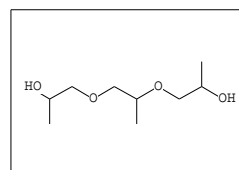
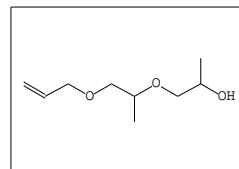
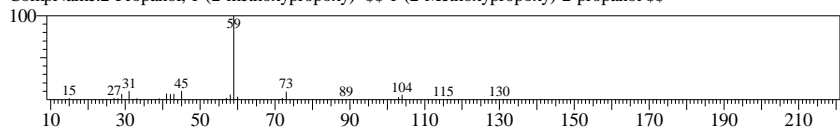
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:8 Entry:14294 Library:NIST08.LIB

SI:82 Formula:C7H16O3 CAS:13429-07-7 MolWeight:148 RetIndex:967

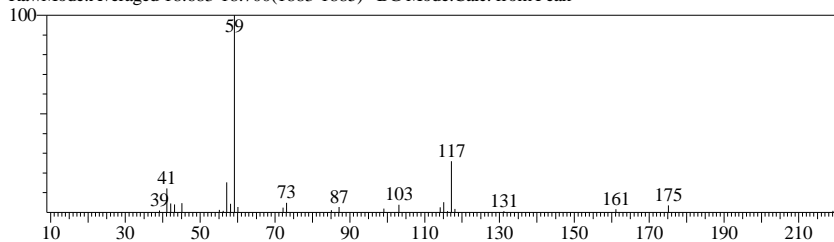
CompName:2-Propanol, 1-(2-methoxypropoxy)- \$ 1-(2-Methoxypropoxy)-2-propanol \$



<< Target >>

Line# 17 R.Time:18.692(Scan#:1884) BasePeak:59.10(172215)

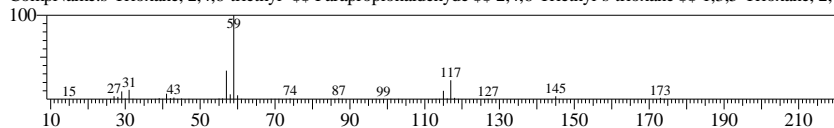
RawMode:Averaged 18.683-18.700(1883-1885) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:88 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

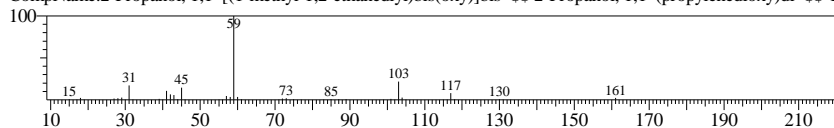
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:37499 Library:NIST08.LIB

SI:86 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

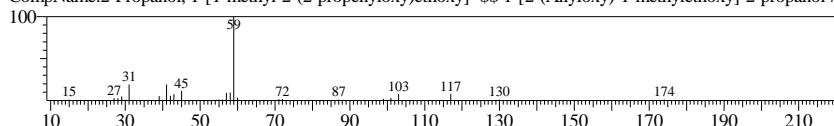
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

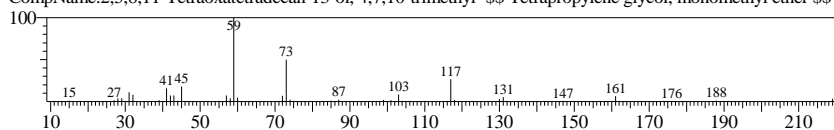
CompName:2-Propanol, 1-[2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:85908 Library:NIST08.LIB

SI:85 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

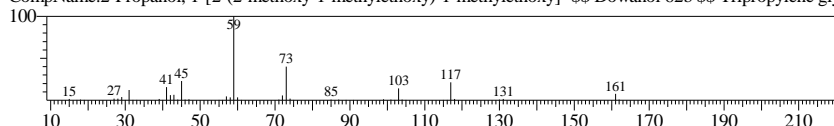
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#:5 Entry:17307 Library:NIST08s.LIB

SI:85 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

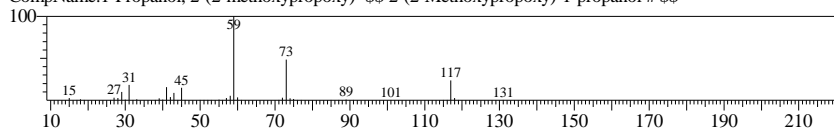
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:14297 Library:NIST08.LIB

SI:85 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

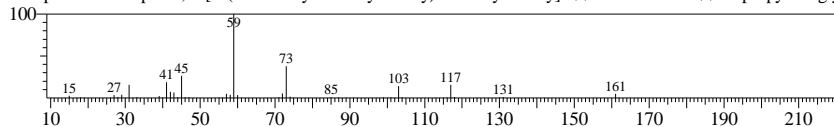
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:84 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

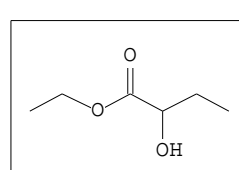
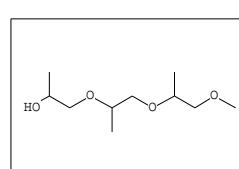
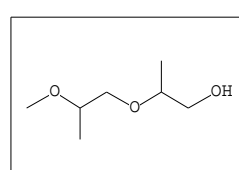
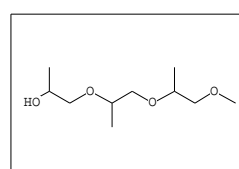
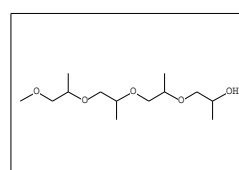
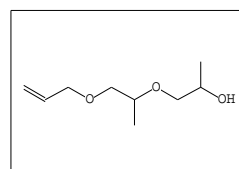
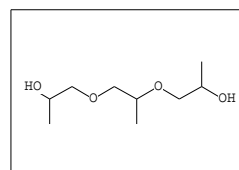
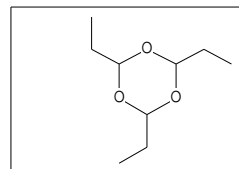
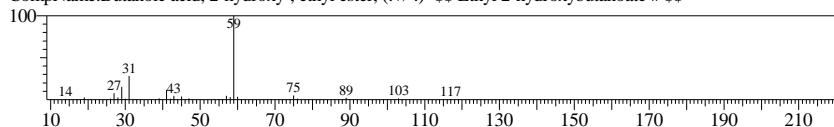
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:8423 Library:NIST08.LIB

SI:84 Formula:C6H12O3 CAS:68057-83-0 MolWeight:132 RetIndex:947

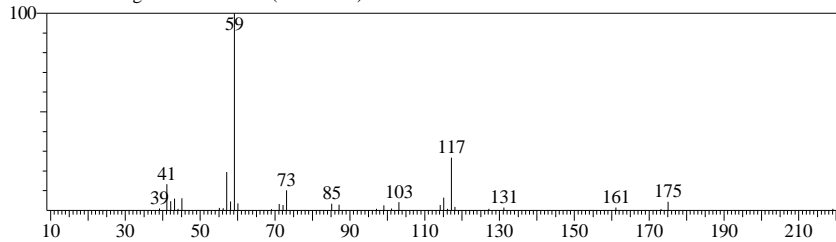
CompName:Butanoic acid, 2-hydroxy-, ethyl ester, (+/-)- \$\$ Ethyl 2-hydroxybutanoate # \$\$



<< Target >>

Line# 18 R.Time:18.742(Scan#:1890) BasePeak:59.10(209577)

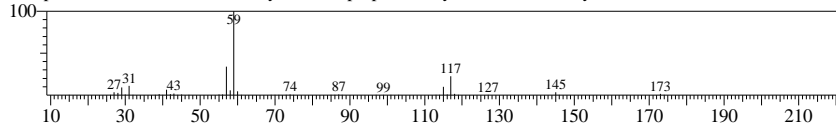
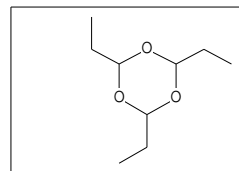
RawMode:Averaged 18.733-18.750(1889-1891) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:88 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

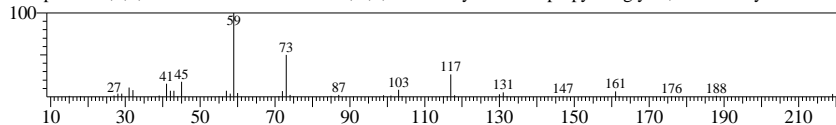
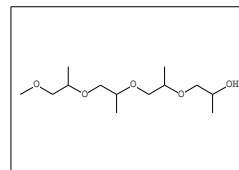
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:86 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

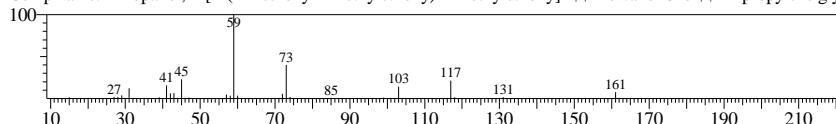
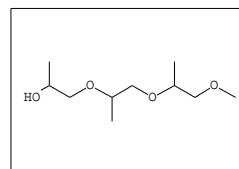
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#3 Entry:17307 Library:NIST08s.LIB

SI:86 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

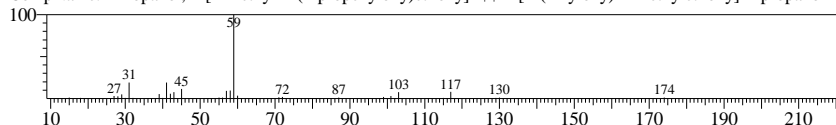
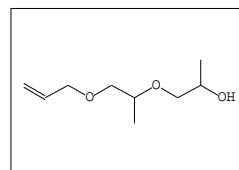
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#4 Entry:27550 Library:NIST08.LIB

SI:85 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

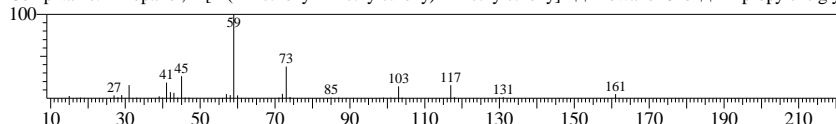
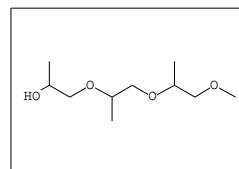
CompName:2-Propanol, 1-[1-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#5 Entry:46317 Library:NIST08.LIB

SI:85 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

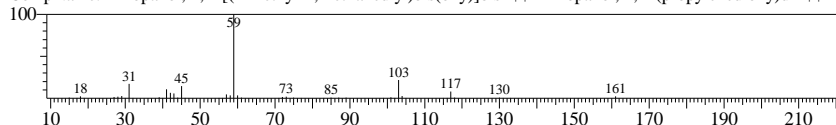
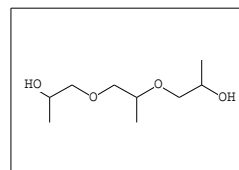
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#6 Entry:37499 Library:NIST08.LIB

SI:85 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

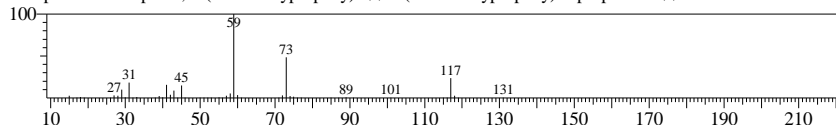
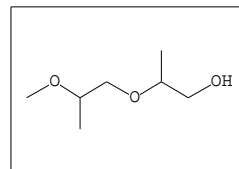
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#7 Entry:14297 Library:NIST08.LIB

SI:85 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

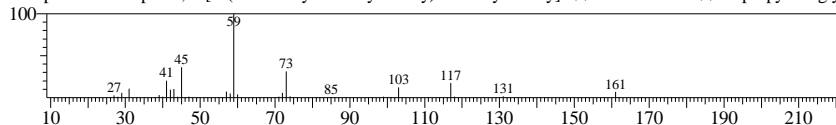
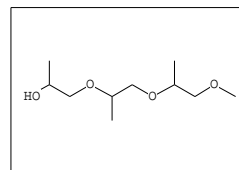
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#8 Entry:17306 Library:NIST08s.LIB

SI:84 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

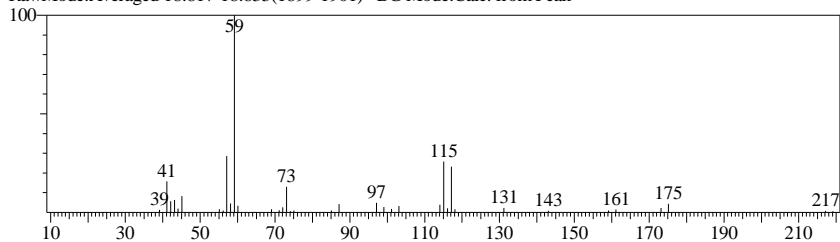
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line#:19 R.Time:18.825(Scan#:1900) BasePeak:59.10(109169)

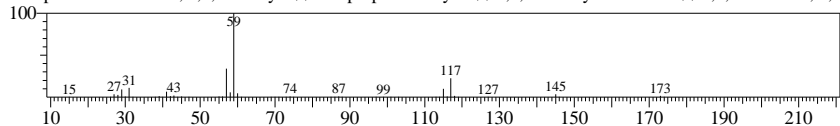
RawMode:Averaged 18.817-18.833(1899-1901) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:87 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

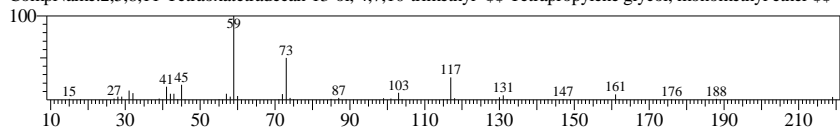
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:85908 Library:NIST08.LIB

SI:84 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

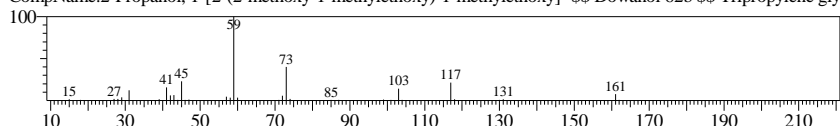
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#:3 Entry:17307 Library:NIST08s.LIB

SI:84 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

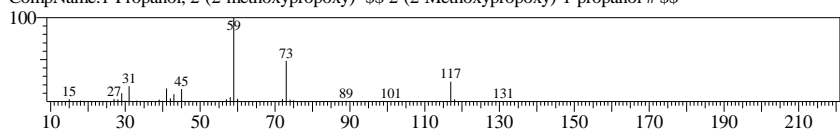
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:4 Entry:14297 Library:NIST08.LIB

SI:83 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

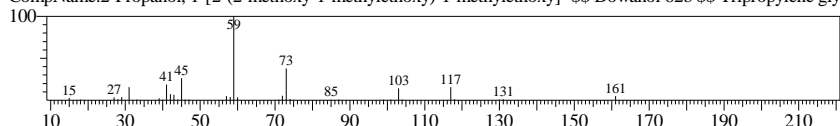
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:5 Entry:46317 Library:NIST08.LIB

SI:83 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

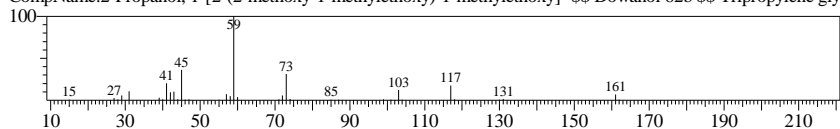
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:17306 Library:NIST08s.LIB

SI:83 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

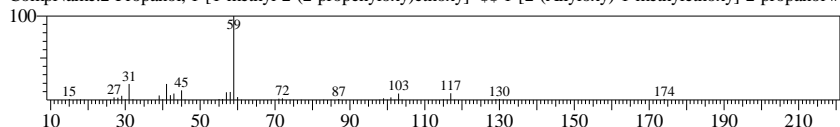
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:7 Entry:27550 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

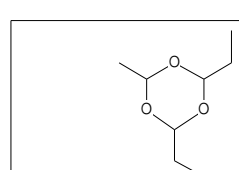
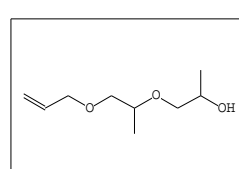
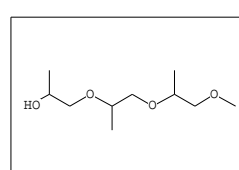
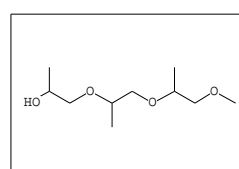
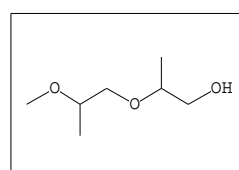
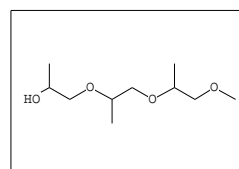
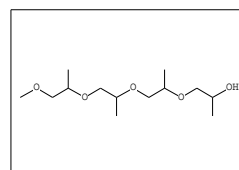
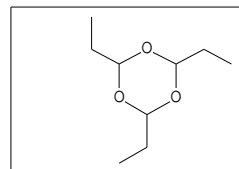
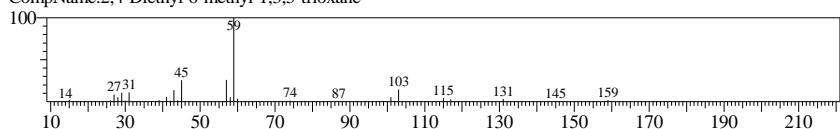
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:8 Entry:19968 Library:NIST08.LIB

SI:82 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

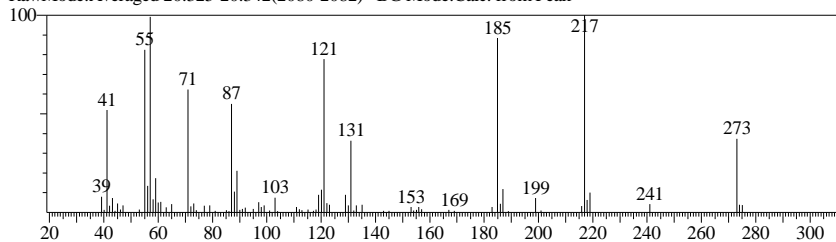
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line# 20 R.Time:20.333(Scan#:2081) BasePeak:216.95(747109)

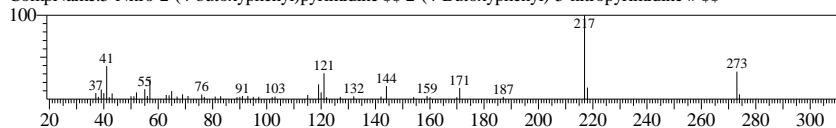
RawMode:Averaged 20.325-20.342(2080-2082) BG Mode:Calc. from Peak



Hit#:1 Entry:92637 Library:NIST08.LIB

SI:57 Formula:C14H15N3O3 CAS:84610-06-0 MolWeight:273 RetIndex:2238

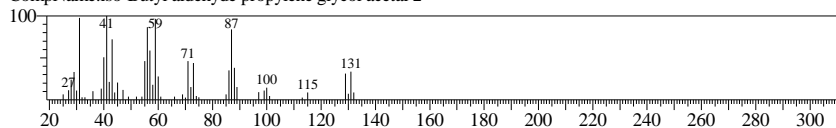
CompName:5-Nitro-2-(4-butoxyphenyl)pyrimidine \$ 2-(4-Butoxyphenyl)-5-nitropyrimidine # \$ \$



Hit#:2 Entry:7864 Library:NIST08.LIB

SI:56 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

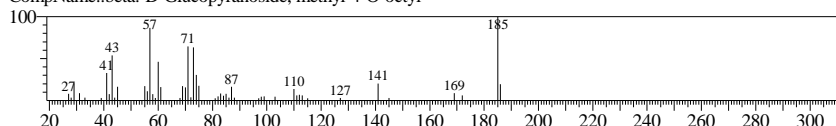
CompName:iso-Butyl aldehyde propylene glycol acetal 2



Hit#:3 Entry:116065 Library:NIST08.LIB

SI:55 Formula:C15H30O6 CAS:0-00-0 MolWeight:306 RetIndex:2358

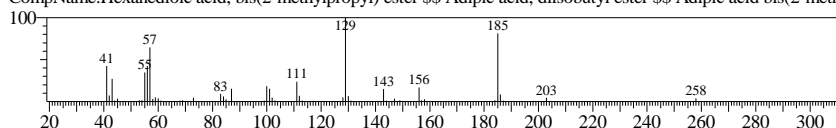
CompName:beta.-D-Glucopyranoside, methyl-4-O-octyl-



Hit#:4 Entry:21972 Library:NIST08s.LIB

SI:55 Formula:C14H26O4 CAS:141-04-8 MolWeight:258 RetIndex:1619

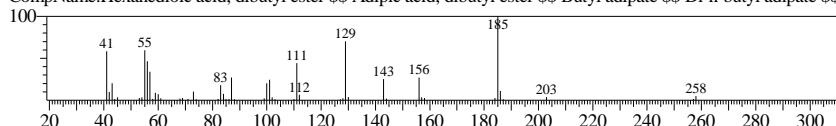
CompName:Hexanedioic acid, bis(2-methylpropyl) ester \$ Adipic acid, diisobutyl ester \$ Adipic acid bis(2-meth



Hit#:5 Entry:81926 Library:NIST08.LIB

SI:55 Formula:C14H26O4 CAS:105-99-7 MolWeight:258 RetIndex:1747

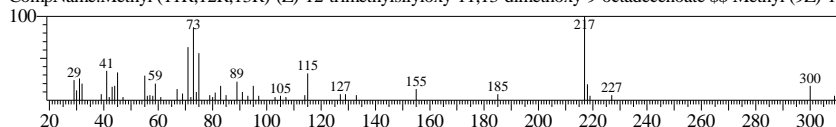
CompName:Hexanedioic acid, dibutyl ester \$ Adipic acid, dibutyl ester \$ Butyl adipate \$ Di-n-butyl adipate \$



Hit#:6 Entry:178021 Library:NIST08.LIB

SI:54 Formula:C24H48O5Si CAS:83303-85-9 MolWeight:444 RetIndex:2528

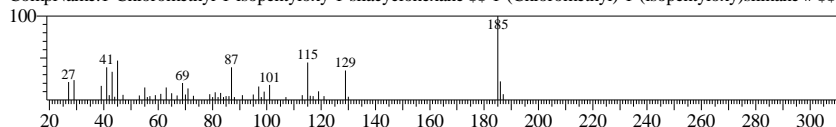
CompName:Methyl (11R,12R,13R)-(Z)-12-trimethylsilyloxy-11,13-dimethoxy-9-octadecenoate \$ Methyl (9Z)-1-



Hit#:7 Entry:65040 Library:NIST08.LIB

SI:54 Formula:C11H23ClOSi CAS:232270-70-1 MolWeight:234 RetIndex:1332

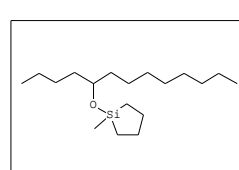
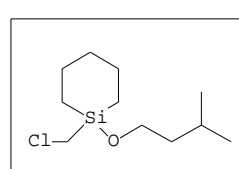
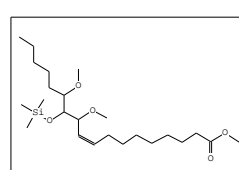
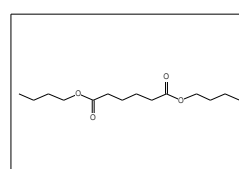
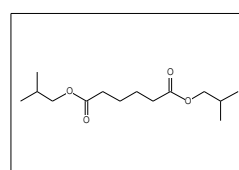
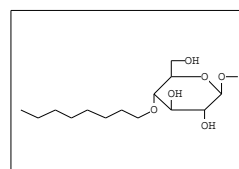
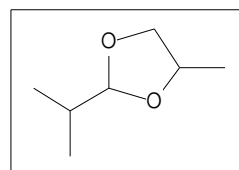
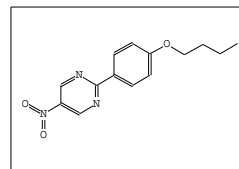
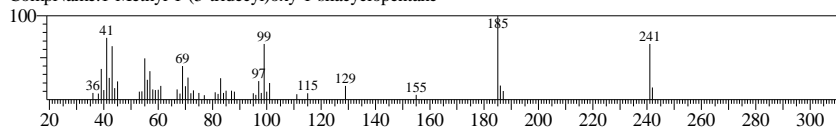
CompName:1-Chloromethyl-1-isopentyloxy-1-silacyclohexane \$ 1-(Chloromethyl)-1-(isopentyloxy)silane # \$



Hit#:8 Entry:110749 Library:NIST08.LIB

SI:53 Formula:C18H38OSi CAS:0-00-0 MolWeight:298 RetIndex:1782

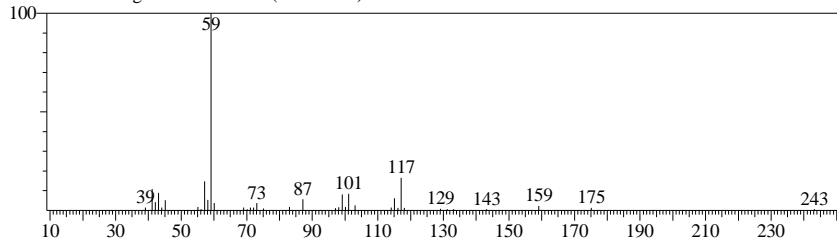
CompName:1-Methyl-1-(5-tridecyl)oxy-1-silacyclopentane



<< Target >>

Line# 21 R.Time:20.558(Scan#:2108) BasePeak:59.10(83057)

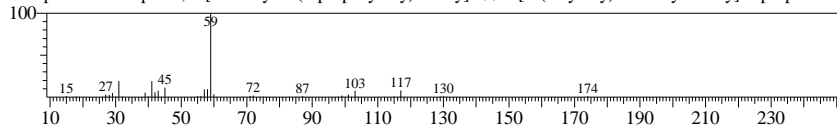
RawMode:Averaged 20.550-20.567(2107-2109) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:88 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

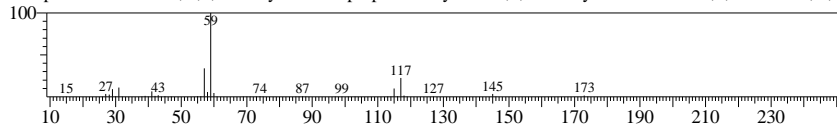
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:27552 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

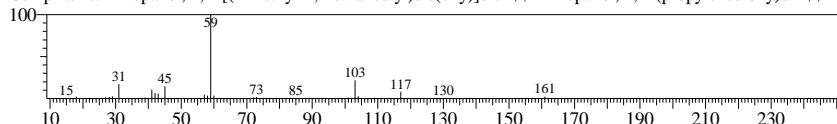
CompName:s-Trioxane, 2,4,6-triethyl- \$- Parapropionaldehyde \$- 2,4,6-Triethyl-s-trioxane \$- 1,3,5-Trioxane, 2,4



Hit#:3 Entry:37499 Library:NIST08.LIB

SI:84 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

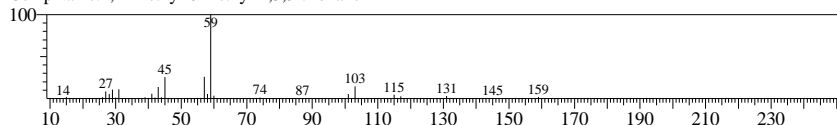
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$- 2-Propanol, 1,1'-(propylenedioxy)di- \$- T



Hit#:4 Entry:19968 Library:NIST08.LIB

SI:84 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

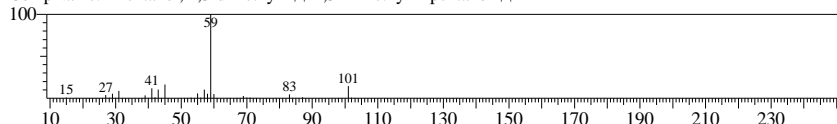
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:5 Entry:4678 Library:NIST08.LIB

SI:83 Formula:C7H16O CAS:4911-70-0 MolWeight:116 RetIndex:745

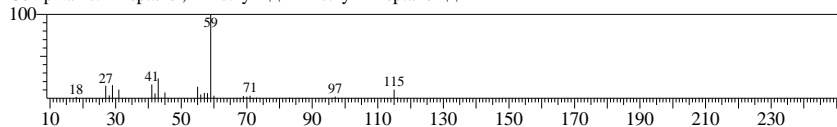
CompName:2-Pentanol, 2,3-dimethyl- \$- 2,3-Dimethyl-2-pentanol \$-



Hit#:6 Entry:5482 Library:NIST08.LIB

SI:83 Formula:C8H18O CAS:625-25-2 MolWeight:130 RetIndex:908

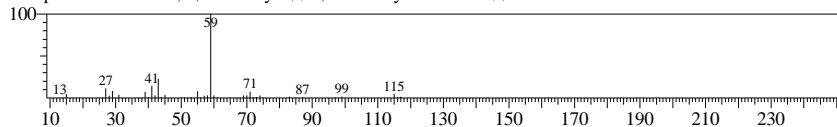
CompName:2-Heptanol, 2-methyl- \$- 2-Methyl-2-heptanol \$-



Hit#:7 Entry:8021 Library:NIST08.LIB

SI:83 Formula:C8H18O CAS:19550-03-9 MolWeight:130 RetIndex:844

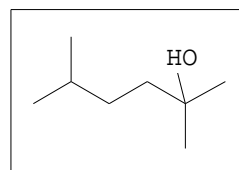
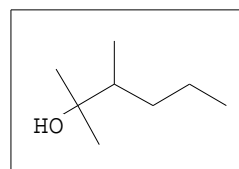
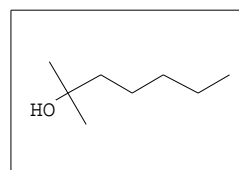
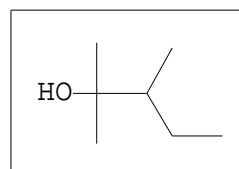
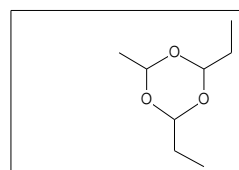
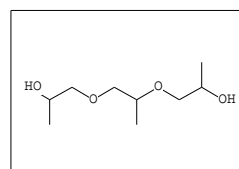
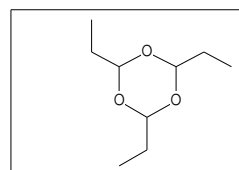
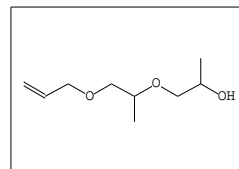
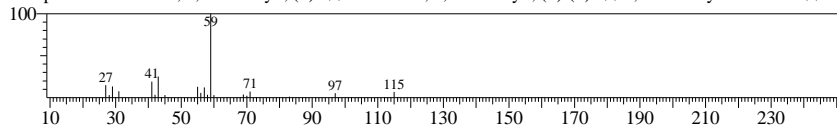
CompName:2-Hexanol, 2,3-dimethyl- \$- 2,3-Dimethyl-2-hexanol \$-



Hit#:8 Entry:5483 Library:NIST08.LIB

SI:82 Formula:C8H18O CAS:3730-60-7 MolWeight:130 RetIndex:844

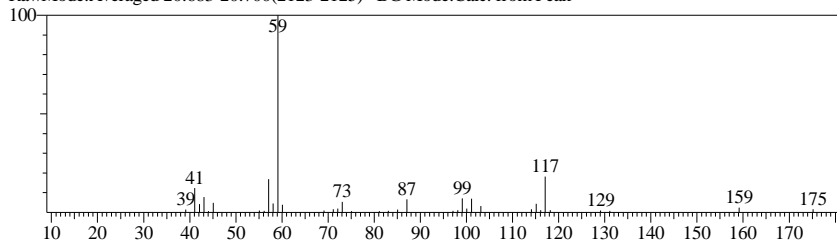
CompName:2-Hexanol, 2,5-dimethyl-, (S)- \$- 2-Hexanol, 2,5-dimethyl-, (S)-(+)- \$- 2,5-Dimethyl-2-hexanol \$- 2-



<< Target >>

Line#:22 R.Time:20.692(Scan#:2124) BasePeak:59.10(209707)

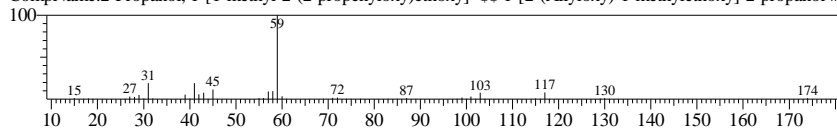
RawMode:Averaged 20.683-20.700(2123-2125) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:88 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

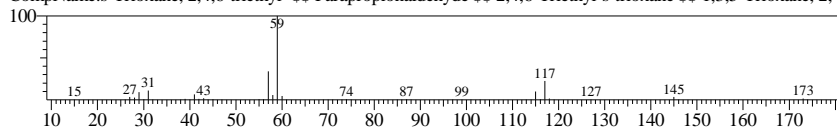
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:27552 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

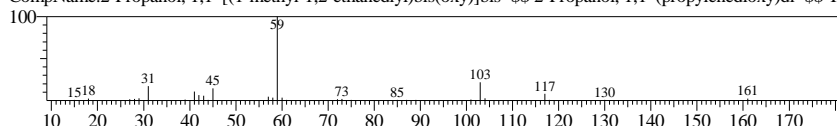
CompName:s-Trioxane, 2,4,6-triethyl- \$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:3 Entry:37499 Library:NIST08.LIB

SI:85 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

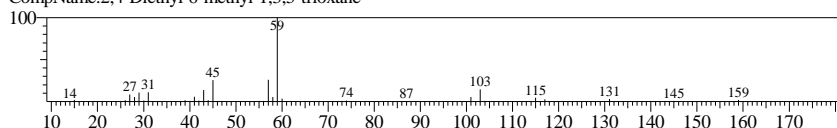
CompName:2-Propanol, 1,1'-[(1-methyl-2-ethanediy)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:4 Entry:19968 Library:NIST08.LIB

SI:84 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

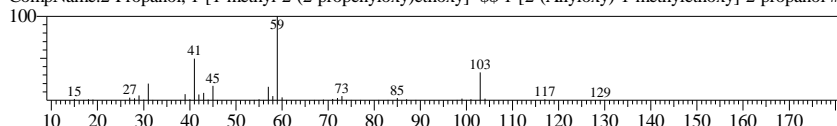
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:5 Entry:12792 Library:NIST08s.LIB

SI:83 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

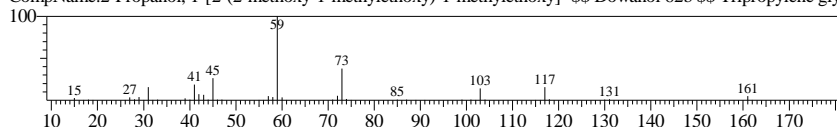
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:6 Entry:46317 Library:NIST08.LIB

SI:83 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

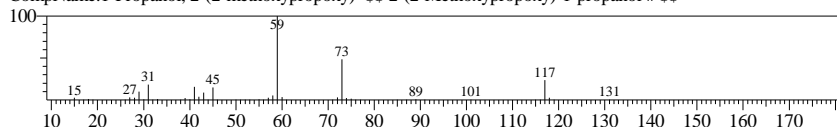
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



Hit#:7 Entry:14297 Library:NIST08.LIB

SI:83 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

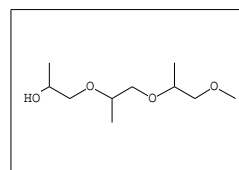
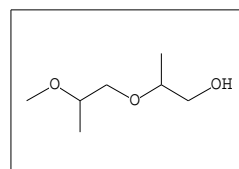
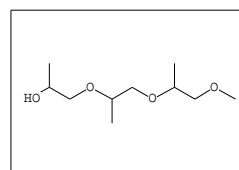
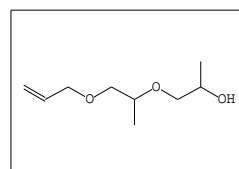
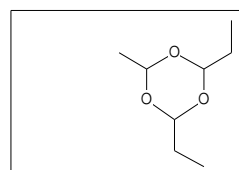
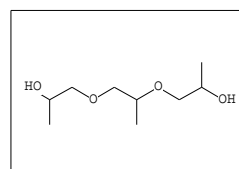
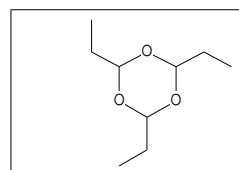
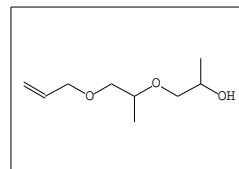
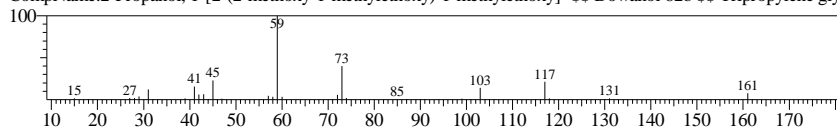
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$ 2-(2-Methoxypropoxy)-1-propanol # \$ \$



Hit#:8 Entry:17307 Library:NIST08s.LIB

SI:83 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

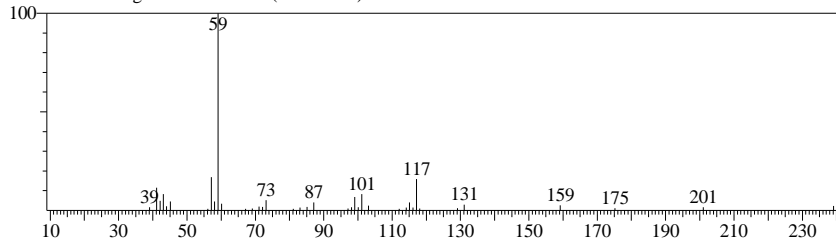
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



<< Target >>

Line# 23 R.Time:20.783(Scan#:2135) BasePeak:59.10(97322)

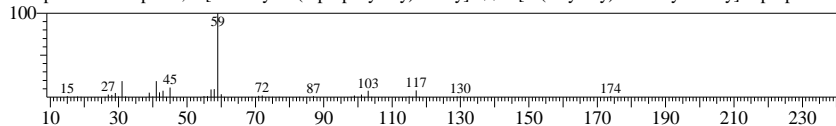
RawMode:Averaged 20.775-20.792(2134-2136) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

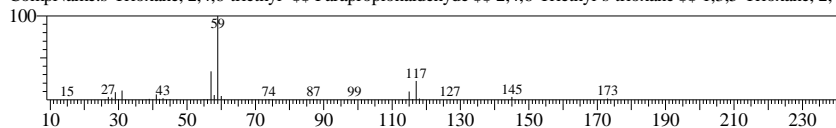
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:27552 Library:NIST08.LIB

SI:84 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

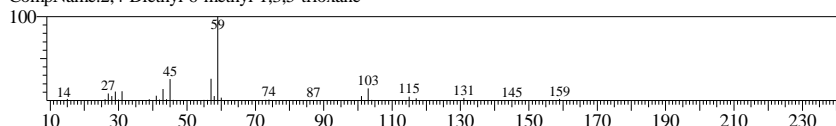
CompName:s-Trioxane, 2,4,6-triethyl- \$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:3 Entry:19968 Library:NIST08.LIB

SI:84 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

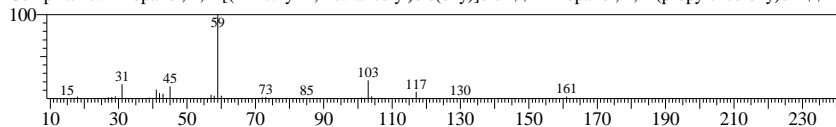
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:4 Entry:37499 Library:NIST08.LIB

SI:83 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

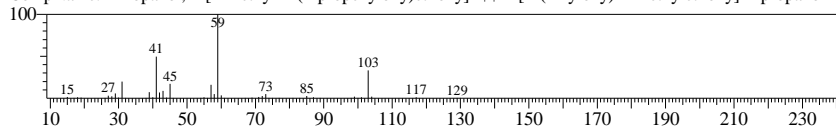
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:5 Entry:12792 Library:NIST08s.LIB

SI:82 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

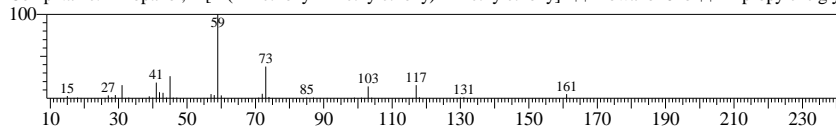
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:6 Entry:46317 Library:NIST08.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

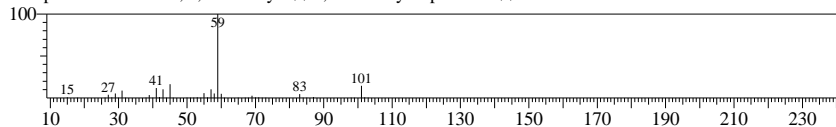
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



Hit#:7 Entry:4678 Library:NIST08.LIB

SI:81 Formula:C7H16O CAS:4911-70-0 MolWeight:116 RetIndex:745

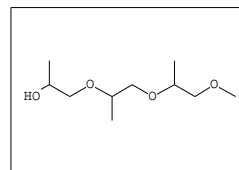
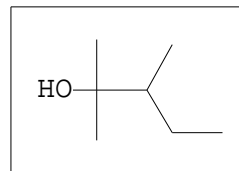
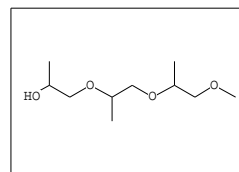
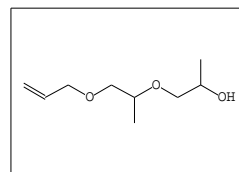
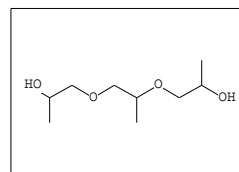
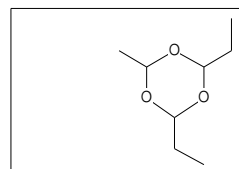
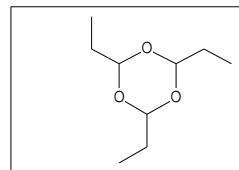
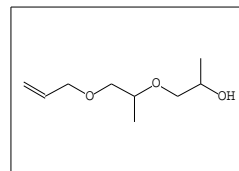
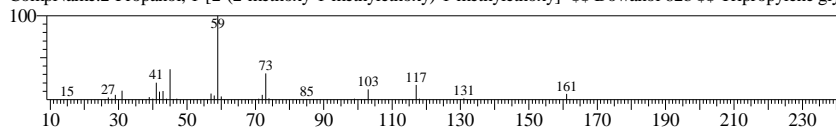
CompName:2-Pentanol, 2,3-dimethyl- \$ 2,3-Dimethyl-2-pentanol \$



Hit#:8 Entry:17306 Library:NIST08s.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

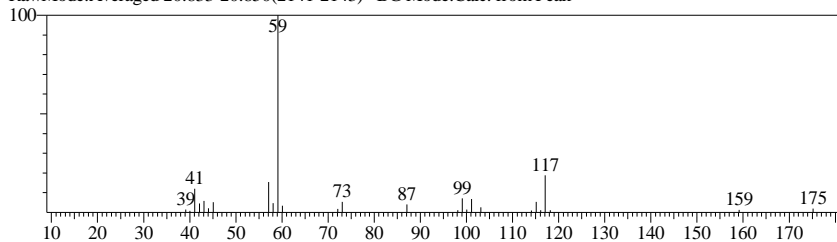
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



<< Target >>

Line# 24 R.Time:20.842(Scan#:2142) BasePeak:59.10(112700)

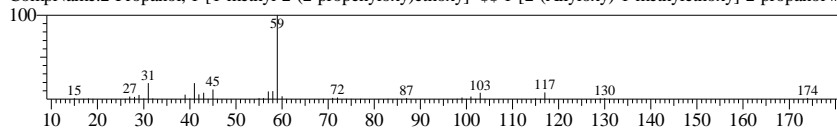
RawMode:Averaged 20.833-20.850(2141-2143) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:87 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

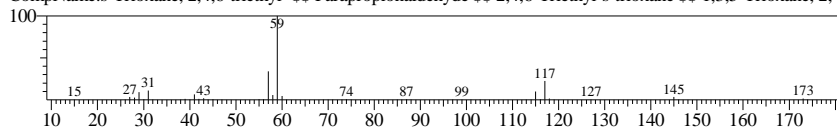
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:27552 Library:NIST08.LIB

SI:87 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

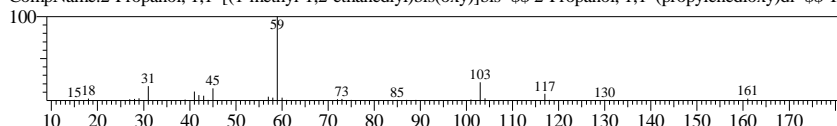
CompName:s-Trioxane, 2,4,6-triethyl- \$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:3 Entry:37499 Library:NIST08.LIB

SI:86 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

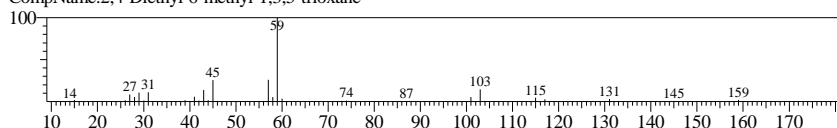
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanedyl)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:4 Entry:19968 Library:NIST08.LIB

SI:84 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

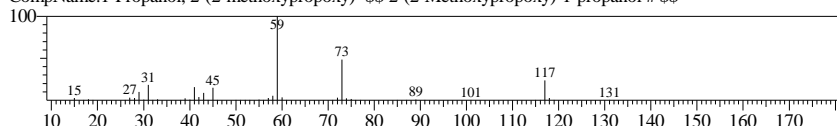
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:5 Entry:14297 Library:NIST08.LIB

SI:84 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

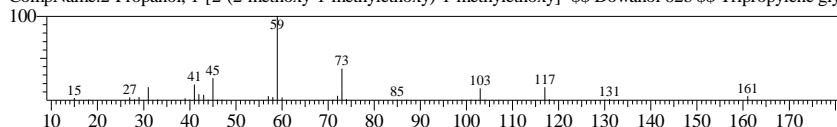
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$ 2-(2-Methoxypropoxy)-1-propanol # \$ \$



Hit#:6 Entry:46317 Library:NIST08.LIB

SI:84 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

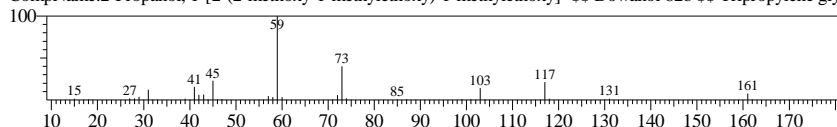
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



Hit#:7 Entry:17307 Library:NIST08s.LIB

SI:83 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

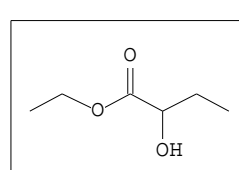
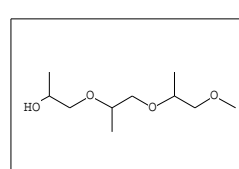
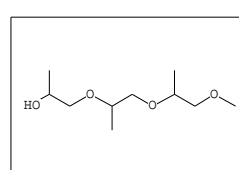
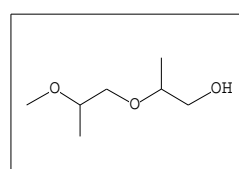
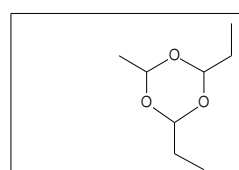
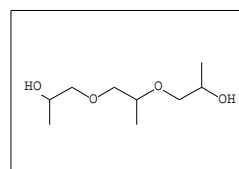
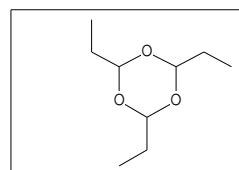
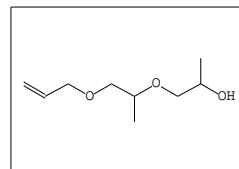
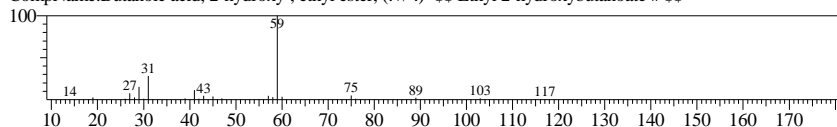
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



Hit#:8 Entry:8423 Library:NIST08.LIB

SI:83 Formula:C6H12O3 CAS:68057-83-0 MolWeight:132 RetIndex:947

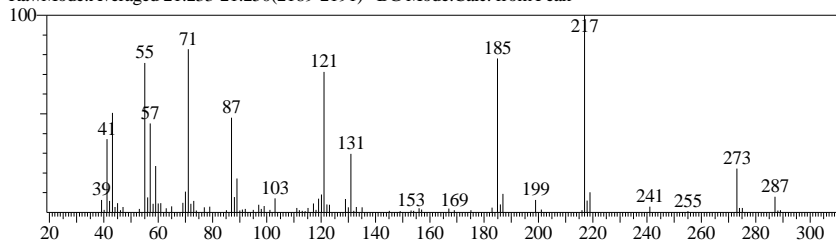
CompName:Butanoic acid, 2-hydroxy-, ethyl ester, (+/-)- \$ Ethyl 2-hydroxybutanoate # \$ \$



<< Target >>

Line# 25 R.Time:21.242(Scan#:2190) BasePeak:216.95(349054)

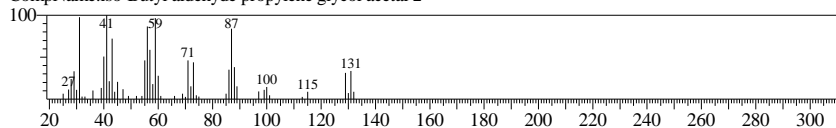
RawMode:Averaged 21.233-21.250(2189-2191) BG Mode:Calc. from Peak



Hit#:1 Entry:7864 Library:NIST08.LIB

SI:58 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

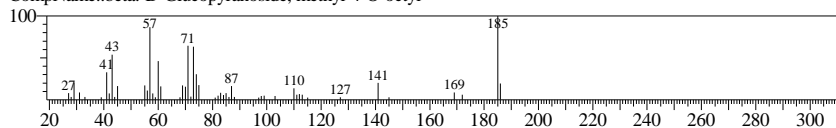
CompName:iso-Butyl aldehyde propylene glycol acetal 2



Hit#:2 Entry:116065 Library:NIST08.LIB

SI:58 Formula:C15H30O6 CAS:0-00-0 MolWeight:306 RetIndex:2358

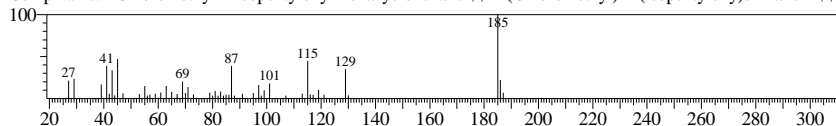
CompName:beta-D-Glucopyranoside, methyl-4-O-octyl-



Hit#:3 Entry:65040 Library:NIST08.LIB

SI:58 Formula:C11H23ClOSi CAS:232270-70-1 MolWeight:234 RetIndex:1332

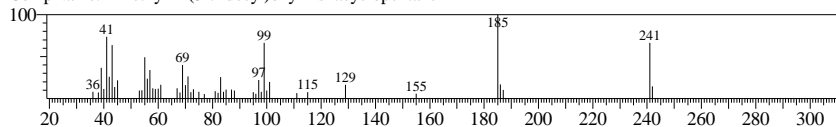
CompName:1-Chloromethyl-1-isopentyloxy-1-silacyclohexane \$\$ 1-(Chloromethyl)-1-(isopentyloxy)silane # \$\$



Hit#:4 Entry:110749 Library:NIST08.LIB

SI:58 Formula:C18H38OSi CAS:0-00-0 MolWeight:298 RetIndex:1782

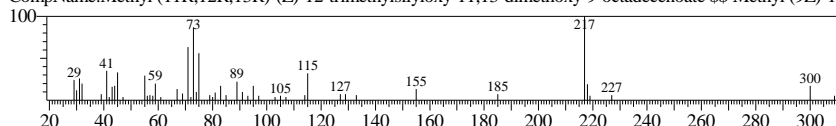
CompName:1-Methyl-1-(5-tridecyl)oxy-1-silacyclopentane



Hit#:5 Entry:178021 Library:NIST08.LIB

SI:57 Formula:C24H48O5Si CAS:83303-85-9 MolWeight:444 RetIndex:2528

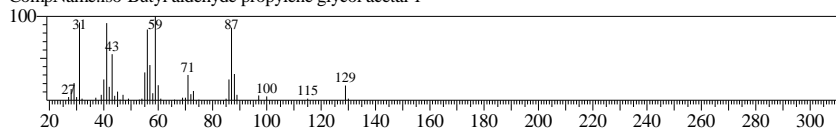
CompName:Methyl (11R,12R,13R)-(Z)-12-trimethylsilyloxy-11,13-dimethoxy-9-octadecenoate \$\$ Methyl (9Z)-1-



Hit#:6 Entry:7906 Library:NIST08.LIB

SI:57 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

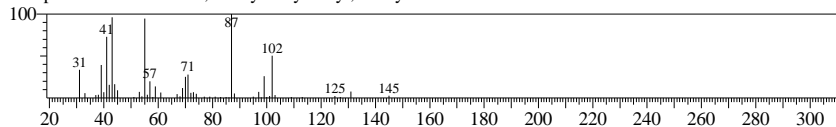
CompName:iso-Butyl aldehyde propylene glycol acetal 1



Hit#:7 Entry:27563 Library:NIST08.LIB

SI:57 Formula:C9H18O3 CAS:0-00-0 MolWeight:174 RetIndex:1181

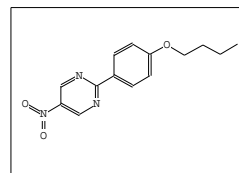
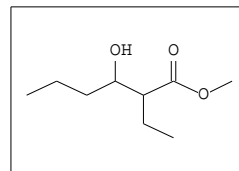
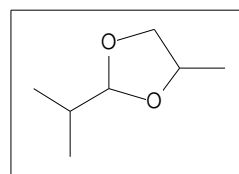
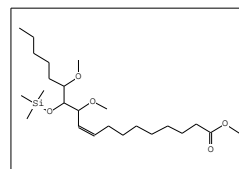
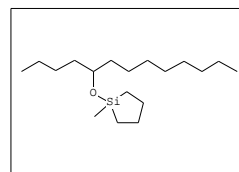
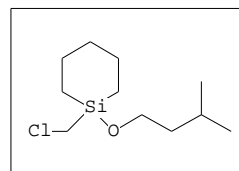
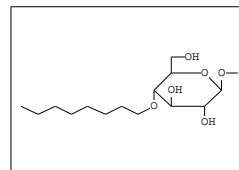
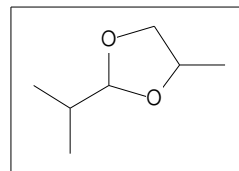
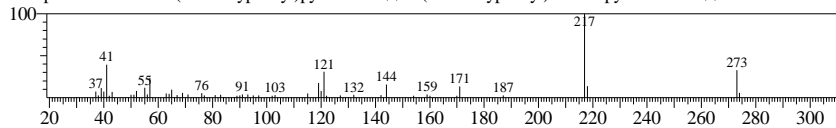
CompName:Hexanoic acid, 2-ethyl-3-hydroxy-, methyl ester



Hit#:8 Entry:92637 Library:NIST08.LIB

SI:57 Formula:C14H15N3O3 CAS:84610-06-0 MolWeight:273 RetIndex:2238

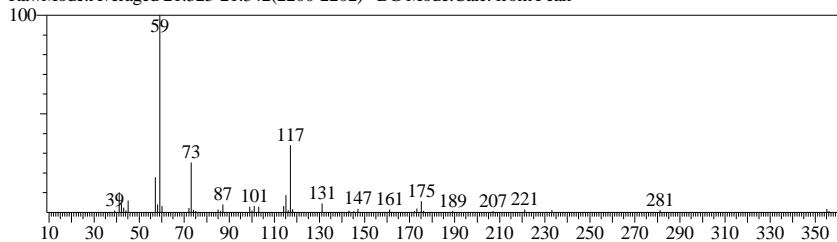
CompName:5-Nitro-2-(4-butoxyphenyl)pyrimidine \$\$ 2-(4-Butoxyphenyl)-5-nitropyrimidine # \$\$



<< Target >>

Line# 26 R.Time: 21.333 (Scan#: 2201) BasePeak: 59.10 (95761)

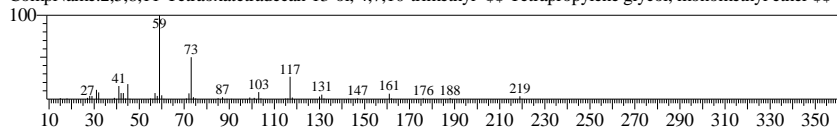
RawMode: Averaged 21.325-21.342 (2200-2202) BG Mode: Calc. from Peak



Hit#: 1 Entry: 85908 Library: NIST08.LIB

SI: 85 Formula: C₁₃H₂₈O₅ CAS: 20324-34-9 MolWeight: 264 RetIndex: 1587

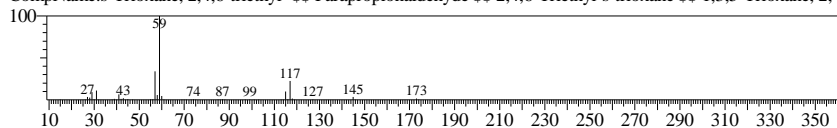
CompName: 2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#: 2 Entry: 27552 Library: NIST08.LIB

SI: 84 Formula: C₉H₁₈O₃ CAS: 2396-42-1 MolWeight: 174 RetIndex: 1168

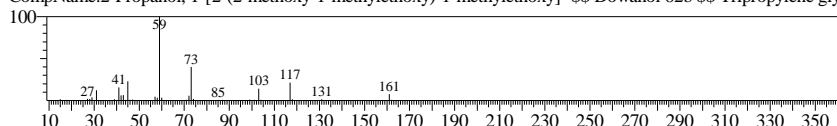
CompName: s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#: 3 Entry: 17307 Library: NIST08s.LIB

SI: 84 Formula: C₁₀H₂₂O₄ CAS: 20324-33-8 MolWeight: 206 RetIndex: 1277

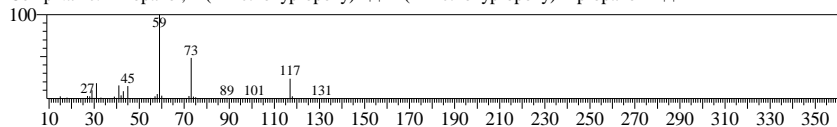
CompName: 2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#: 4 Entry: 14297 Library: NIST08.LIB

SI: 83 Formula: C₇H₁₆O₃ CAS: 13588-28-8 MolWeight: 148 RetIndex: 983

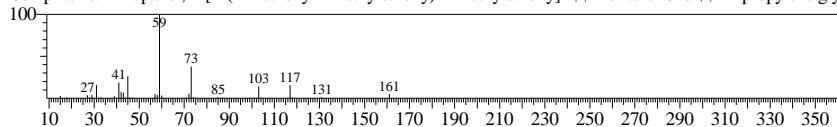
CompName: 1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#: 5 Entry: 46317 Library: NIST08.LIB

SI: 83 Formula: C₁₀H₂₂O₄ CAS: 20324-33-8 MolWeight: 206 RetIndex: 1277

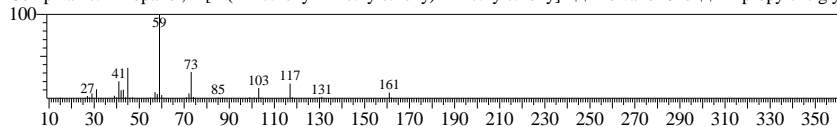
CompName: 2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#: 6 Entry: 17306 Library: NIST08s.LIB

SI: 82 Formula: C₁₀H₂₂O₄ CAS: 20324-33-8 MolWeight: 206 RetIndex: 1277

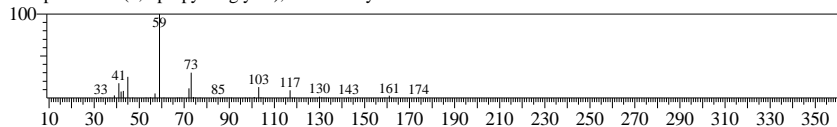
CompName: 2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#: 7 Entry: 46316 Library: NIST08.LIB

SI: 81 Formula: C₁₀H₂₂O₄ CAS: 0-00-0 MolWeight: 206 RetIndex: 1277

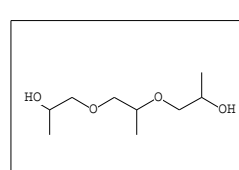
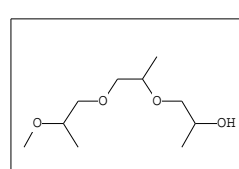
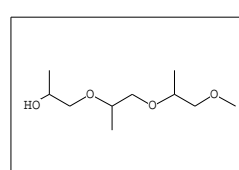
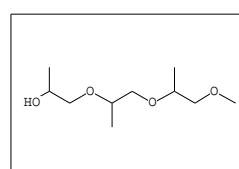
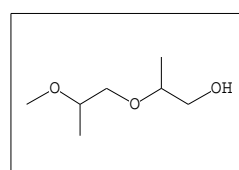
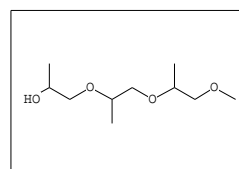
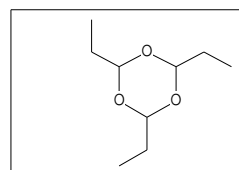
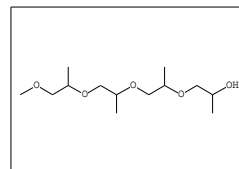
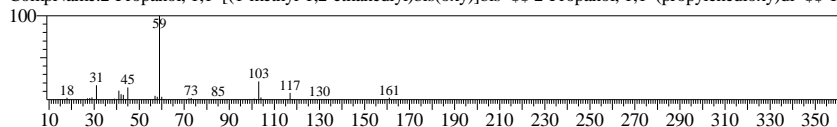
CompName: Tri(1,2-propyleneglycol), monomethyl ether



Hit#: 8 Entry: 37499 Library: NIST08.LIB

SI: 80 Formula: C₉H₂₀O₄ CAS: 1638-16-0 MolWeight: 192 RetIndex: 1328

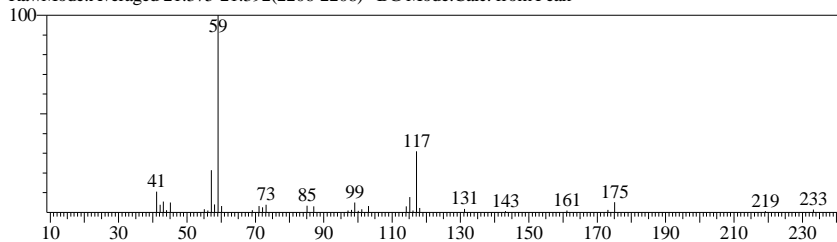
CompName: 2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



<< Target >>

Line#:27 R.Time:21.383(Scan#:2207) BasePeak:59.10(162232)

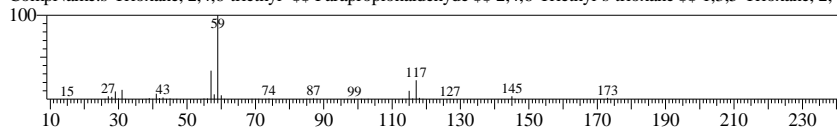
RawMode:Averaged 21.375-21.392(2206-2208) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:87 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

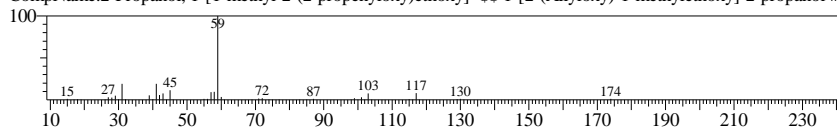
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:85 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

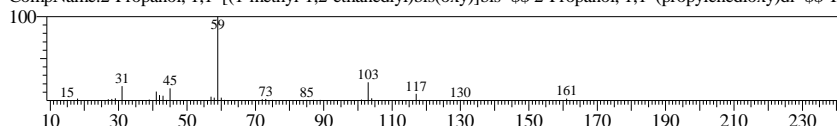
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:37499 Library:NIST08.LIB

SI:83 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

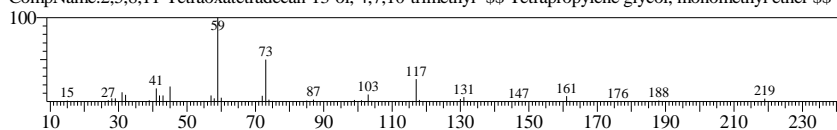
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:4 Entry:85908 Library:NIST08.LIB

SI:83 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

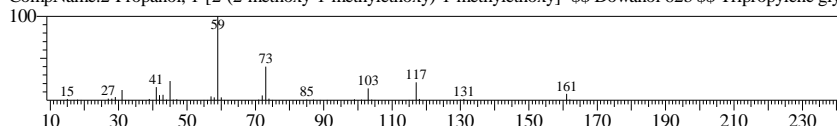
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#:5 Entry:17307 Library:NIST08s.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

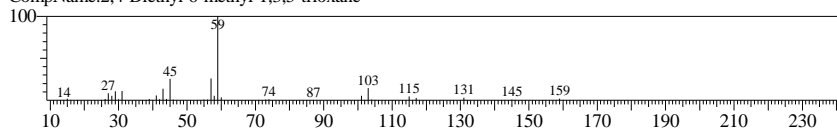
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:19968 Library:NIST08.LIB

SI:82 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

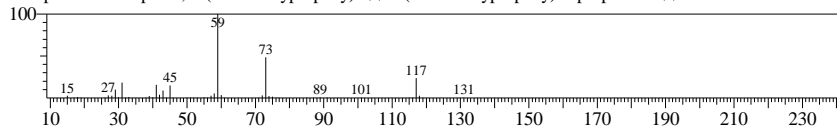
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:7 Entry:14297 Library:NIST08.LIB

SI:82 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

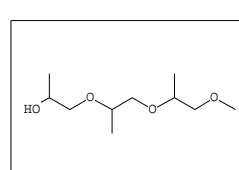
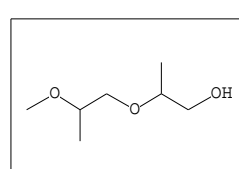
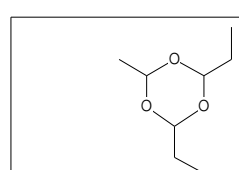
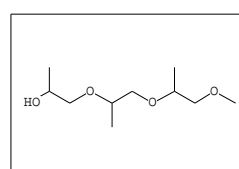
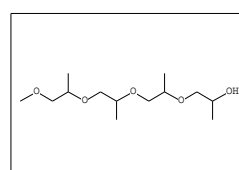
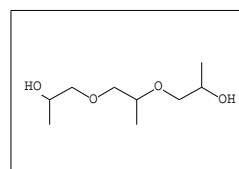
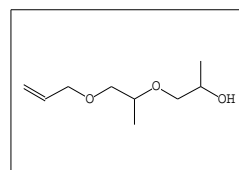
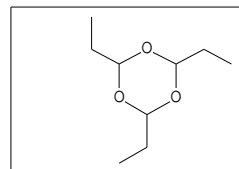
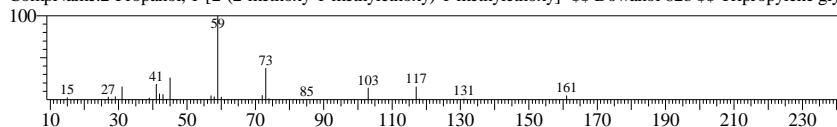
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:8 Entry:46317 Library:NIST08.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

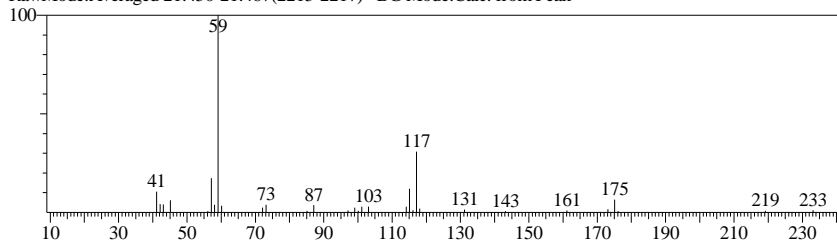
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line# 28 R.Time:21.458(Scan#:2216) BasePeak:59.10(317120)

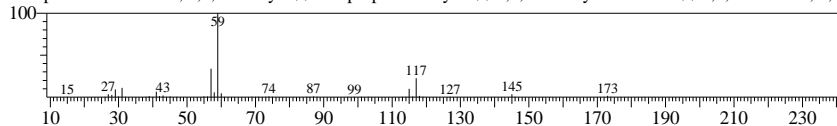
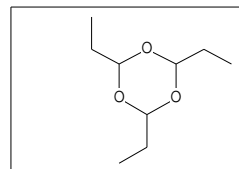
RawMode:Averaged 21.450-21.467(2215-2217) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:87 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

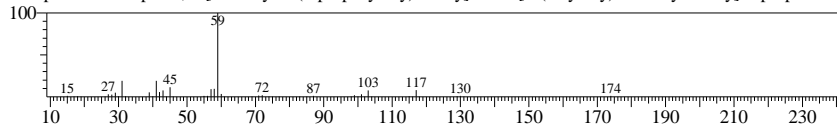
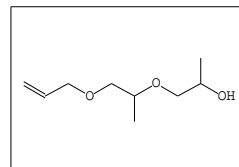
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:84 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

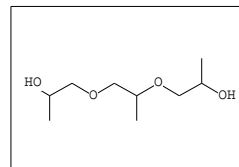
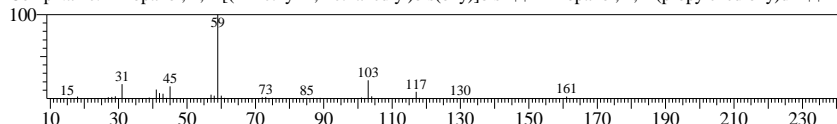
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:37499 Library:NIST08.LIB

SI:83 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

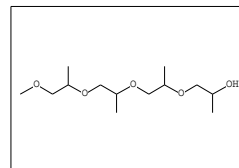
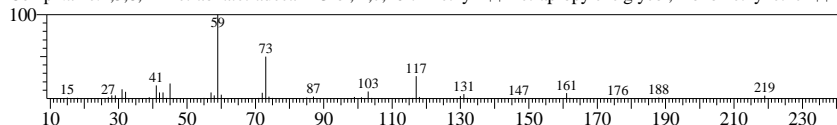
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:4 Entry:85908 Library:NIST08.LIB

SI:83 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

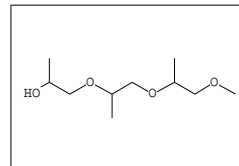
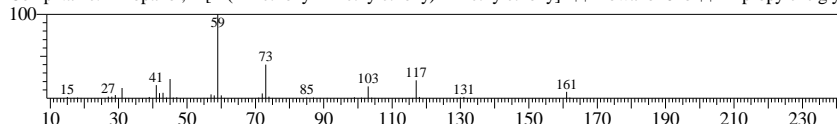
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#:5 Entry:17307 Library:NIST08s.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

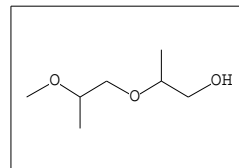
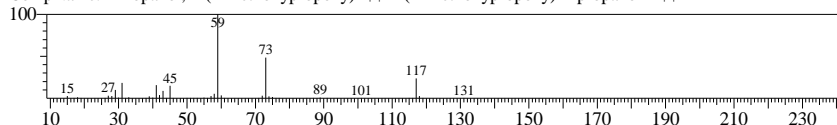
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:14297 Library:NIST08.LIB

SI:82 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

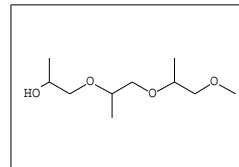
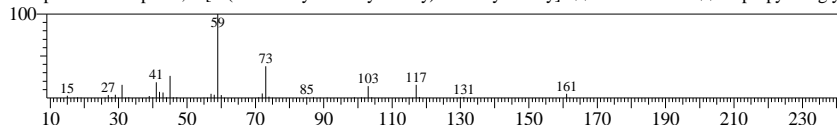
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

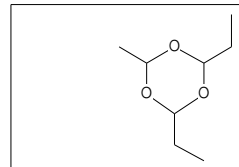
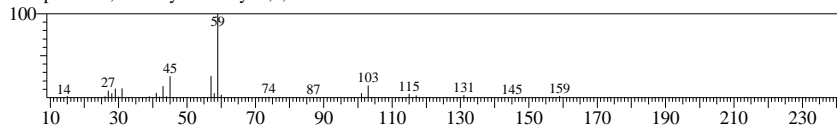
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:19968 Library:NIST08.LIB

SI:81 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

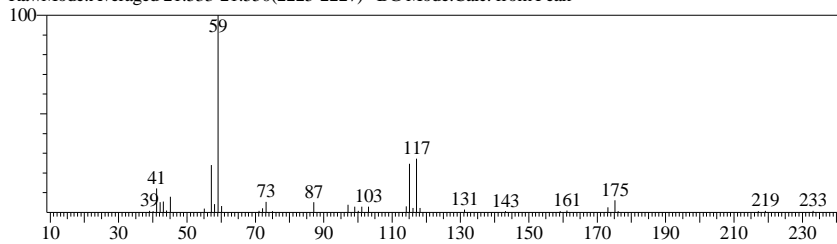
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line# 29 R.Time:21.542(Scan#:2226) BasePeak:59.10(146683)

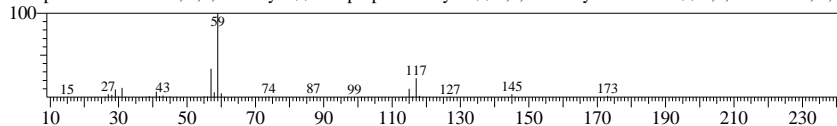
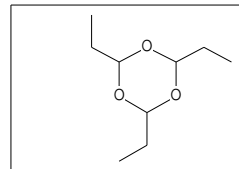
RawMode:Averaged 21.533-21.550(2225-2227) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

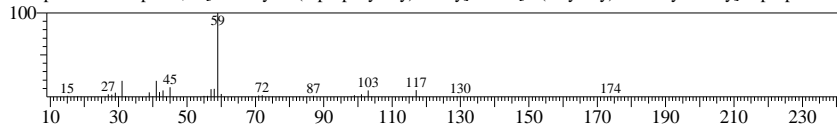
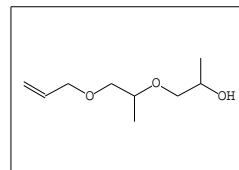
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:83 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

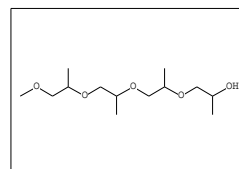
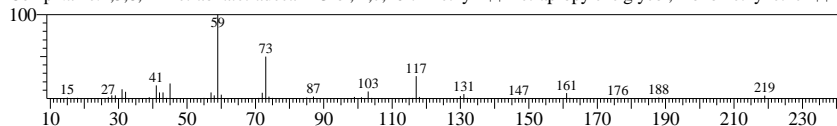
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:85908 Library:NIST08.LIB

SI:82 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

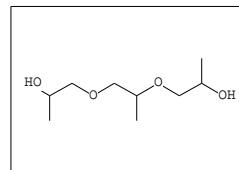
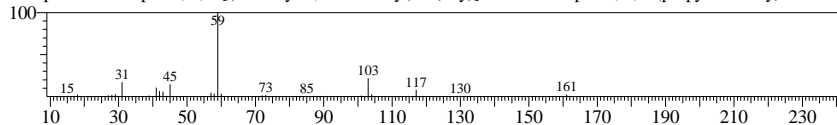
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#:4 Entry:37499 Library:NIST08.LIB

SI:82 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

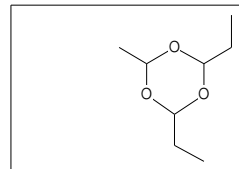
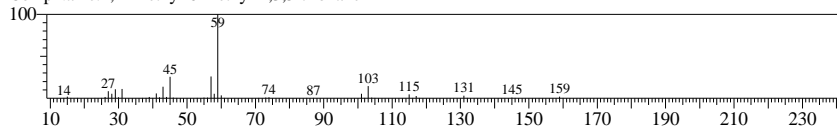
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:5 Entry:19968 Library:NIST08.LIB

SI:82 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

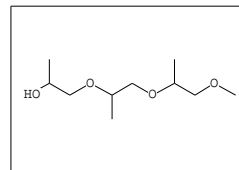
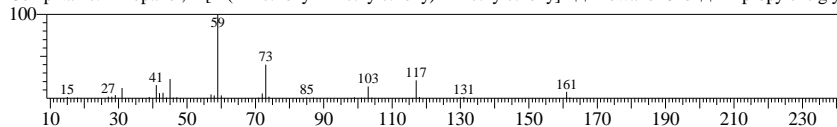
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:6 Entry:17307 Library:NIST08s.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

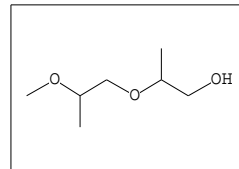
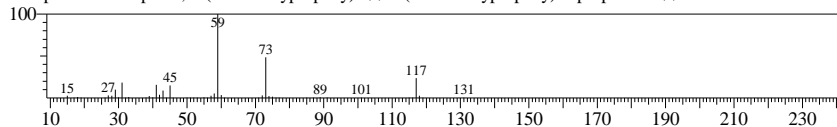
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:7 Entry:14297 Library:NIST08.LIB

SI:81 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

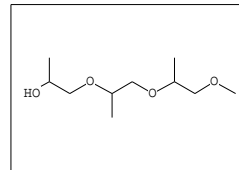
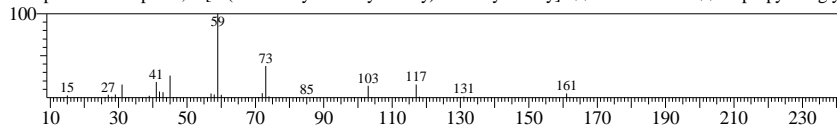
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:8 Entry:46317 Library:NIST08.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

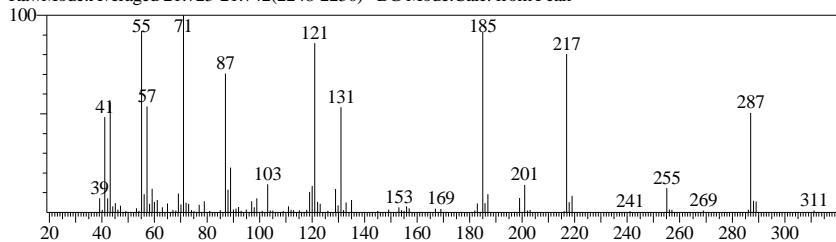
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line#:30 R.Time:21.733(Scan#:2249) BasePeak:71.05(512559)

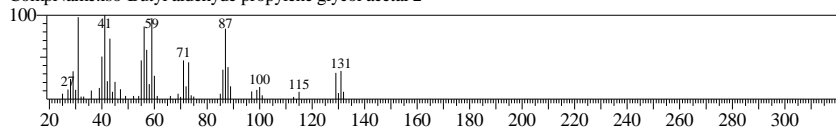
RawMode:Averaged 21.725-21.742(2248-2250) BG Mode:Calc. from Peak



Hit#:1 Entry:7864 Library:NIST08.LIB

SI:58 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

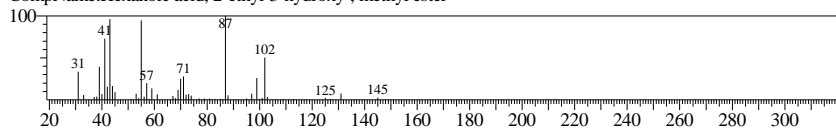
CompName:iso-Butyl aldehyde propylene glycol acetal 2



Hit#:2 Entry:27563 Library:NIST08.LIB

SI:56 Formula:C9H18O3 CAS:0-00-0 MolWeight:174 RetIndex:1181

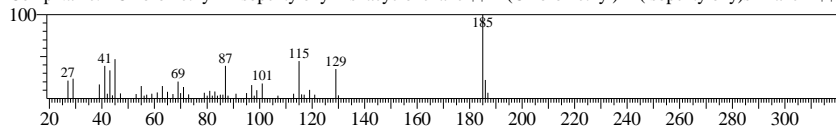
CompName:Hexanoic acid, 2-ethyl-3-hydroxy-, methyl ester



Hit#:3 Entry:65040 Library:NIST08.LIB

SI:56 Formula:C11H23ClOSi CAS:232270-70-1 MolWeight:234 RetIndex:1332

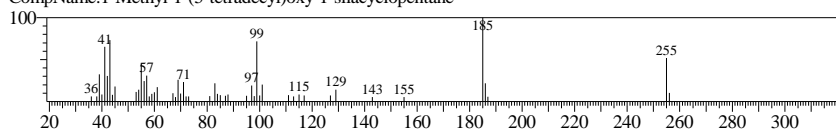
CompName:1-Chloromethyl-1-isopentyloxy-1-silacyclohexane \$\$ 1-(Chloromethyl)-1-(isopentyloxy)silane # \$\$



Hit#:4 Entry:120665 Library:NIST08.LIB

SI:56 Formula:C19H40OSi CAS:0-00-0 MolWeight:312 RetIndex:1881

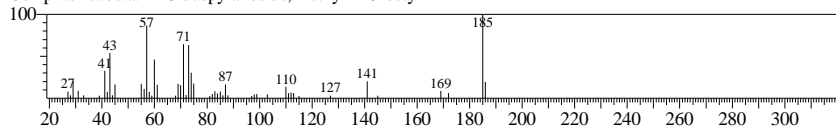
CompName:1-Methyl-1-(5-tetradecyloxy)-1-silacyclopentane



Hit#:5 Entry:116065 Library:NIST08.LIB

SI:56 Formula:C15H30O6 CAS:0-00-0 MolWeight:306 RetIndex:2358

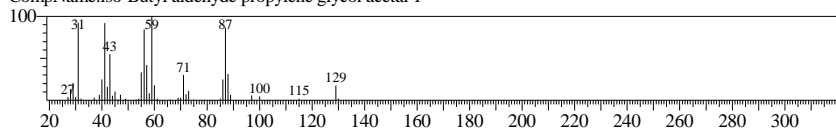
CompName:beta-D-Glucopyranoside, methyl-4-O-octyl-



Hit#:6 Entry:7906 Library:NIST08.LIB

SI:55 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

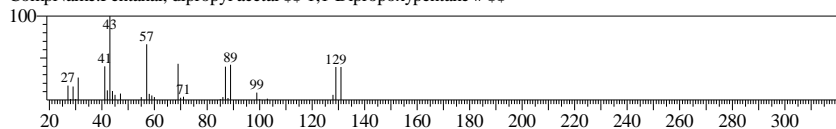
CompName:iso-Butyl aldehyde propylene glycol acetal 1



Hit#:7 Entry:35579 Library:NIST08.LIB

SI:55 Formula:C11H24O2 CAS:13112-64-6 MolWeight:188 RetIndex:1202

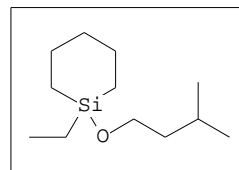
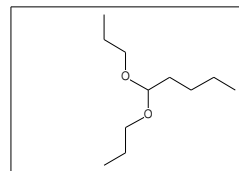
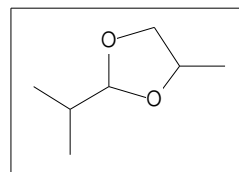
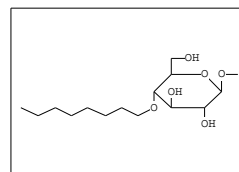
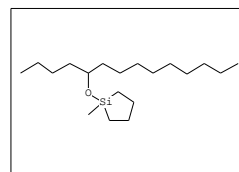
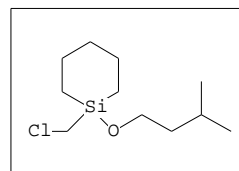
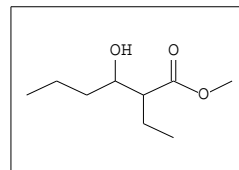
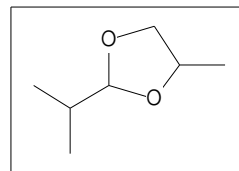
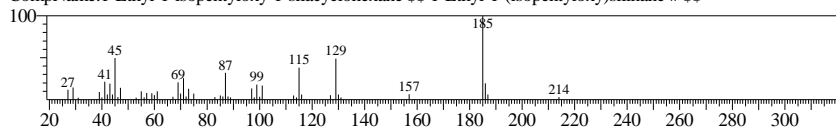
CompName:Pentanal, dipropyl acetal \$\$ 1,1-Dipropoxypentane # \$\$



Hit#:8 Entry:51928 Library:NIST08.LIB

SI:55 Formula:C12H26OSi CAS:232270-69-8 MolWeight:214 RetIndex:1206

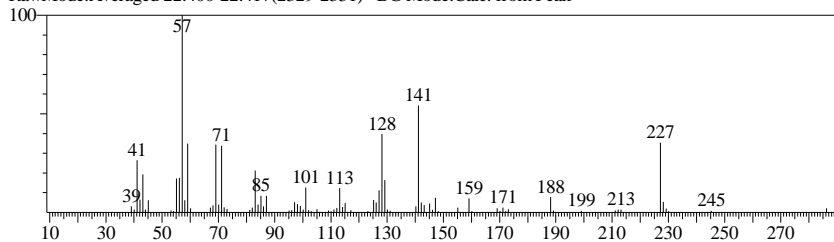
CompName:1-Ethyl-1-isopentyloxy-1-silacyclohexane \$\$ 1-Ethyl-1-(isopentyloxy)silane # \$\$



<< Target >>

Line#:31 R.Time:22.408(Scan#:2330) BasePeak:57.10(97608)

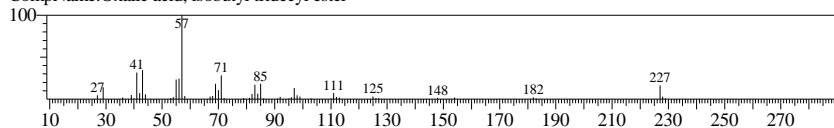
RawMode:Averaged 22.400-22.417(2329-2331) BG Mode:Calc. from Peak



Hit#:1 Entry:131301 Library:NIST08.LIB

SI:68 Formula:C19H36O4 CAS:0-00-0 MolWeight:328 RetIndex:2180

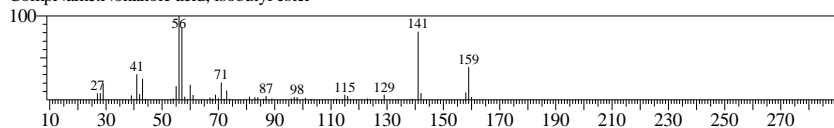
CompName:Oxalic acid, isobutyl tridecyl ester



Hit#:2 Entry:52004 Library:NIST08.LIB

SI:67 Formula:C13H26O2 CAS:0-00-0 MolWeight:214 RetIndex:1417

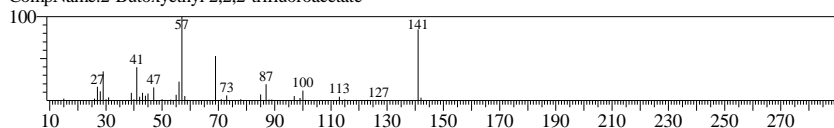
CompName:Nonanoic acid, isobutyl ester



Hit#:3 Entry:51556 Library:NIST08.LIB

SI:67 Formula:C8H13F3O3 CAS:0-00-0 MolWeight:214 RetIndex:894

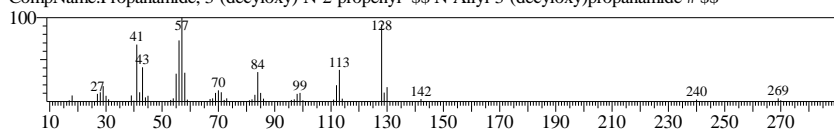
CompName:2-Butoxyethyl 2,2,2-trifluoroacetate



Hit#:4 Entry:89752 Library:NIST08.LIB

SI:67 Formula:C16H31NO2 CAS:54889-73-5 MolWeight:269 RetIndex:2078

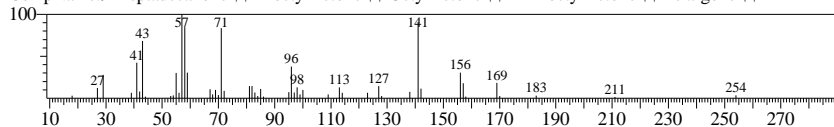
CompName:Propanamide, 3-(decyloxy)-N-2-propenyl- N-Allyl-3-(decyloxy)propanamide # \$ \$



Hit#:5 Entry:21706 Library:NIST08s.LIB

SI:66 Formula:C17H34O CAS:540-08-9 MolWeight:254 RetIndex:1847

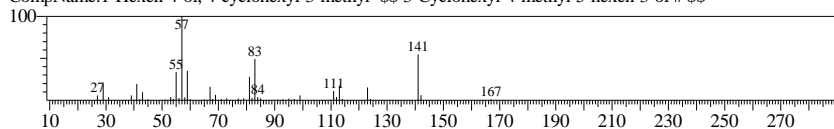
CompName:9-Heptadecanone \$ \$ Dioctyl ketone \$ \$ Octyl ketone \$ \$ Di-n-octyl ketone \$ \$ Pelargone \$ \$



Hit#:6 Entry:40635 Library:NIST08.LIB

SI:66 Formula:C13H24O CAS:344308-85-6 MolWeight:196 RetIndex:1395

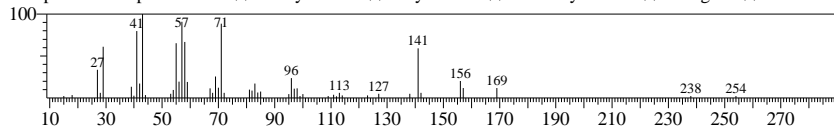
CompName:1-Hexen-4-ol, 4-cyclohexyl-3-methyl- 3-Cyclohexyl-4-methyl-5-hexen-3-ol # \$ \$



Hit#:7 Entry:21704 Library:NIST08s.LIB

SI:66 Formula:C17H34O CAS:540-08-9 MolWeight:254 RetIndex:1847

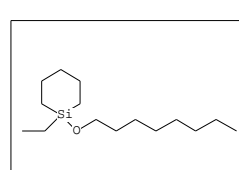
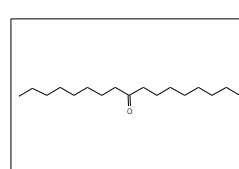
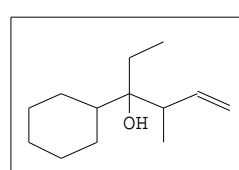
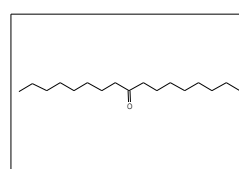
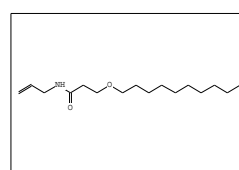
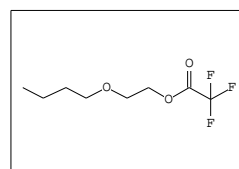
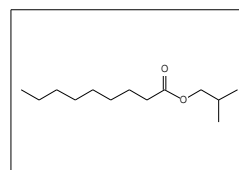
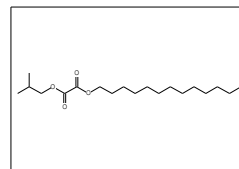
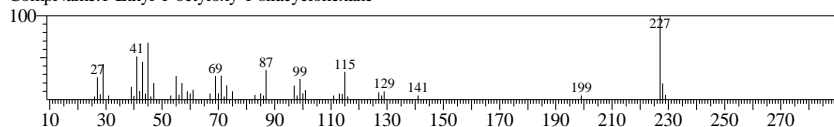
CompName:9-Heptadecanone \$ \$ Dioctyl ketone \$ \$ Octyl ketone \$ \$ Di-n-octyl ketone \$ \$ Pelargone \$ \$



Hit#:8 Entry:80626 Library:NIST08.LIB

SI:66 Formula:C15H32OSi CAS:0-00-0 MolWeight:256 RetIndex:1568

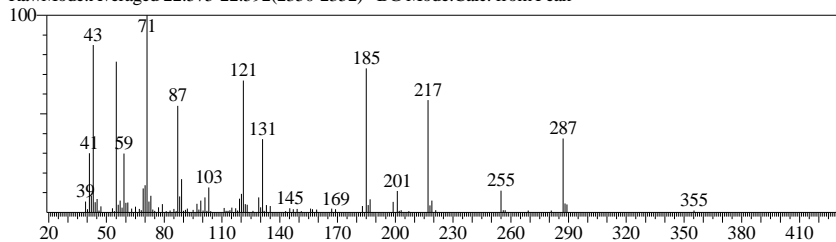
CompName:1-Ethyl-1-octyloxy-1-silacyclohexane



<< Target >>

Line#:32 R.Time:22.583(Scan#:2351) BasePeak:71.05(128290)

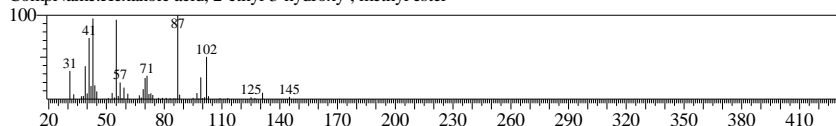
RawMode:Averaged 22.575-22.592(2350-2352) BG Mode:Calc. from Peak



Hit#:1 Entry:27563 Library:NIST08.LIB

SI:61 Formula:C9H18O3 CAS:0-00-0 MolWeight:174 RetIndex:1181

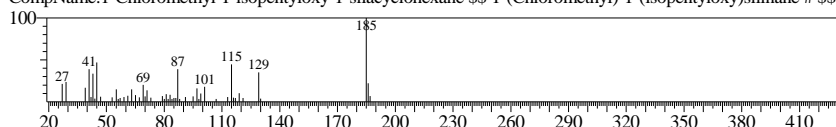
CompName:Hexanoic acid, 2-ethyl-3-hydroxy-, methyl ester



Hit#:2 Entry:65040 Library:NIST08.LIB

SI:59 Formula:C11H23ClOSi CAS:232270-70-1 MolWeight:234 RetIndex:1332

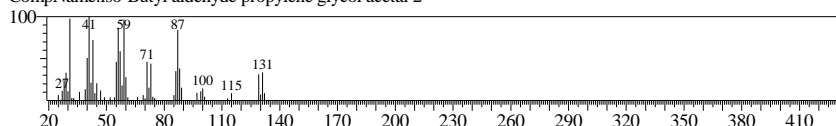
CompName:1-Chloromethyl-1-isopentyloxy-1-silacyclohexane \$\$ 1-(Chloromethyl)-1-(isopentyloxy)silane # \$\$



Hit#:3 Entry:7864 Library:NIST08.LIB

SI:58 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

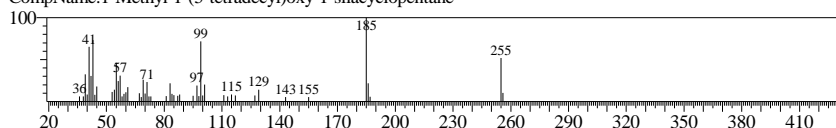
CompName:iso-Butyl aldehyde propylene glycol acetal 2



Hit#:4 Entry:120665 Library:NIST08.LIB

SI:58 Formula:C19H40OSi CAS:0-00-0 MolWeight:312 RetIndex:1881

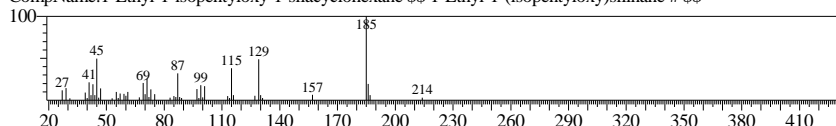
CompName:1-Methyl-1-(5-tetradecyloxy)-1-silacyclopentane



Hit#:5 Entry:51928 Library:NIST08.LIB

SI:58 Formula:C12H26OSi CAS:232270-69-8 MolWeight:214 RetIndex:1206

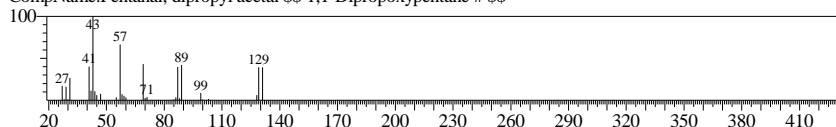
CompName:1-Ethyl-1-isopentyloxy-1-silacyclohexane \$\$ 1-Ethyl-1-(isopentyloxy)silane # \$\$



Hit#:6 Entry:35579 Library:NIST08.LIB

SI:57 Formula:C11H24O2 CAS:13112-64-6 MolWeight:188 RetIndex:1202

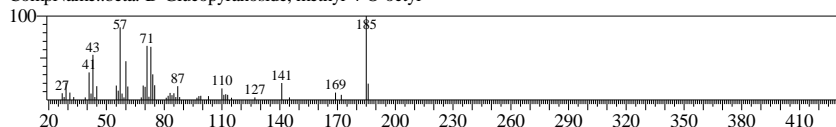
CompName:Pentanal, dipropyl acetal \$\$ 1,1-Dipropoxypentane # \$\$



Hit#:7 Entry:116065 Library:NIST08.LIB

SI:56 Formula:C15H30O6 CAS:0-00-0 MolWeight:306 RetIndex:2358

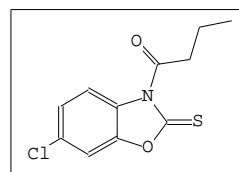
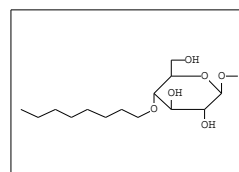
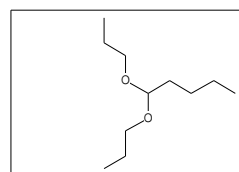
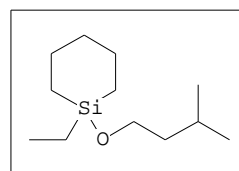
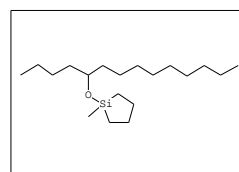
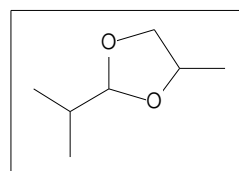
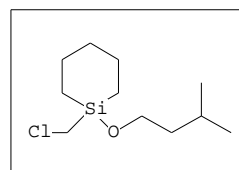
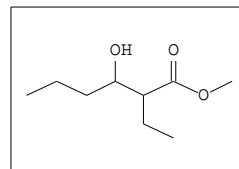
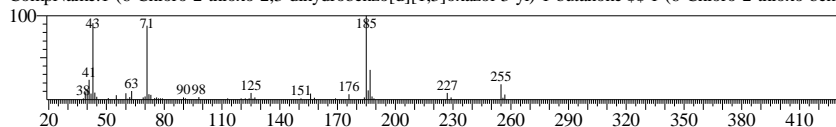
CompName:.beta.-D-Glucopyranoside, methyl-4-O-octyl-



Hit#:8 Entry:21742 Library:NIST08s.LIB

SI:56 Formula:C11H10ClNO2S CAS:31315-64-7 MolWeight:255 RetIndex:2071

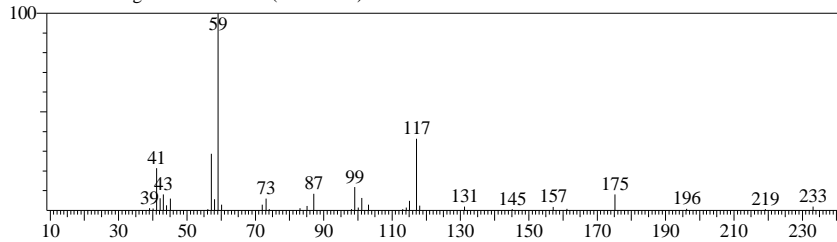
CompName:1-(6-Chloro-2-thioxo-2,3-dihydrobenzo[d][1,3]oxazol-3-yl)-1-butanone \$\$ 1-(6-Chloro-2-thioxo-benz



<< Target >>

Line#33 R.Time:22.958(Scan#:2396) BasePeak:59.10(65985)

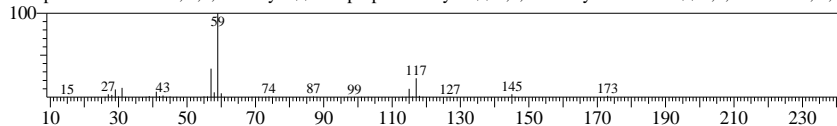
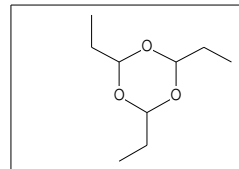
RawMode:Averaged 22.950-22.967(2395-2397) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:84 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

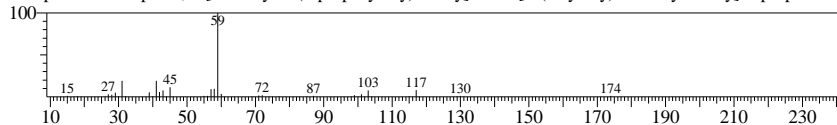
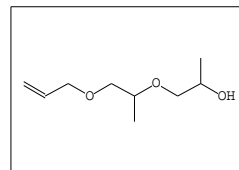
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:83 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

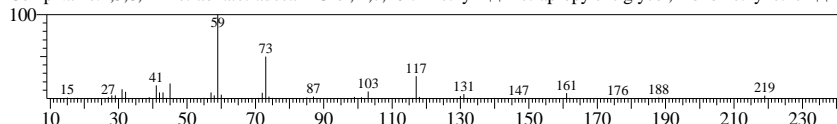
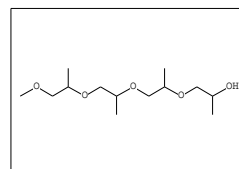
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:85908 Library:NIST08.LIB

SI:82 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

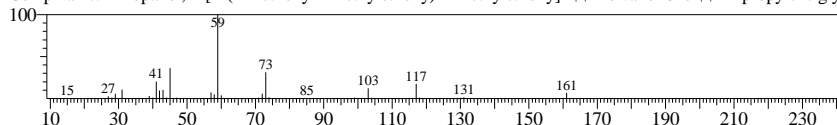
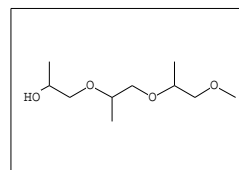
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#:4 Entry:17306 Library:NIST08s.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

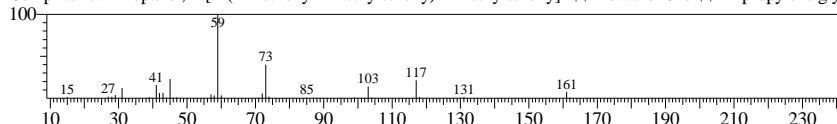
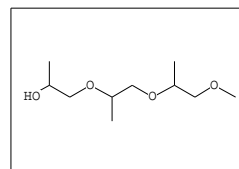
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:5 Entry:17307 Library:NIST08s.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

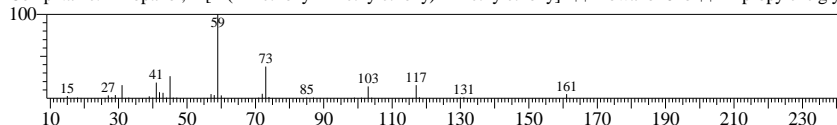
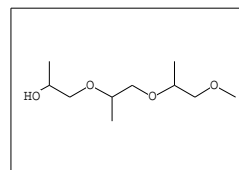
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:46317 Library:NIST08.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

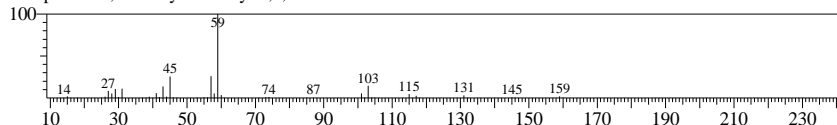
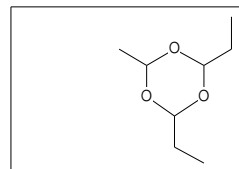
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:7 Entry:19968 Library:NIST08.LIB

SI:81 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

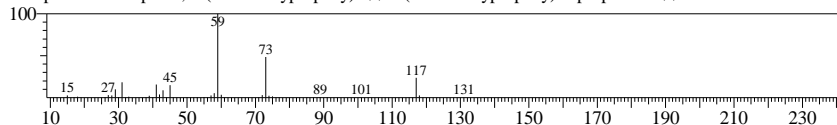
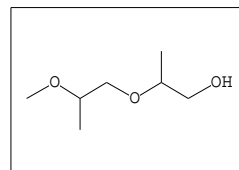
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:8 Entry:14297 Library:NIST08.LIB

SI:81 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

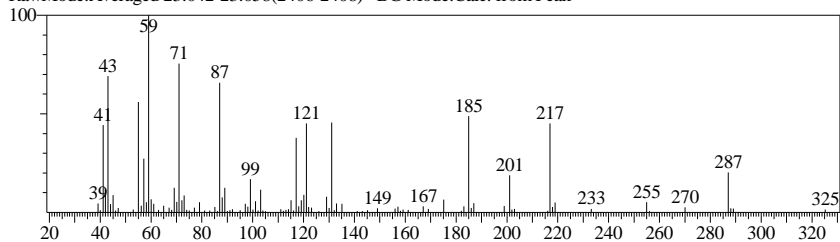
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



<< Target >>

Line#:34 R.Time:23.050(Scan#:2407) BasePeak:59.10(160299)

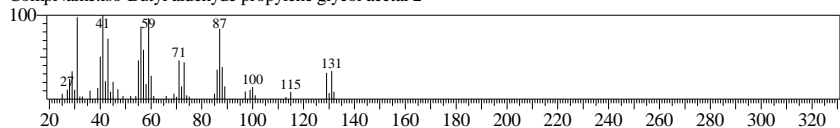
RawMode:Averaged 23.042-23.058(2406-2408) BG Mode:Calc. from Peak



Hit#:1 Entry:7864 Library:NIST08.LIB

SI:66 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

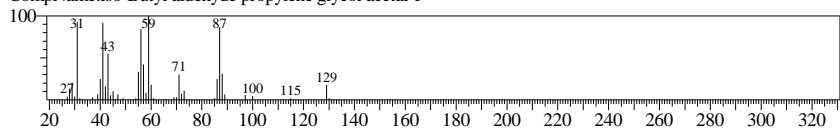
CompName:iso-Butyl aldehyde propylene glycol acetal 2



Hit#:2 Entry:7906 Library:NIST08.LIB

SI:64 Formula:C7H14O2 CAS:0-00-0 MolWeight:130 RetIndex:834

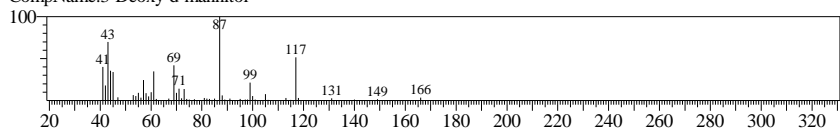
CompName:iso-Butyl aldehyde propylene glycol acetal 1



Hit#:3 Entry:22754 Library:NIST08.LIB

SI:63 Formula:C6H14O5 CAS:0-00-0 MolWeight:166 RetIndex:1590

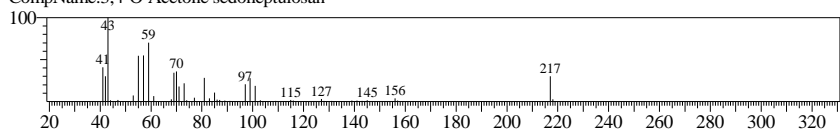
CompName:3-Deoxy-d-mannitol



Hit#:4 Entry:63643 Library:NIST08.LIB

SI:63 Formula:C10H16O6 CAS:0-00-0 MolWeight:232 RetIndex:1718

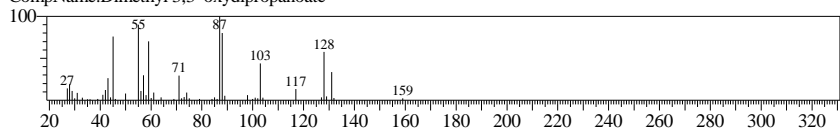
CompName:3,4-O-Acetone sedoheptulosan



Hit#:5 Entry:36354 Library:NIST08.LIB

SI:63 Formula:C8H14O5 CAS:94102-60-0 MolWeight:190 RetIndex:1227

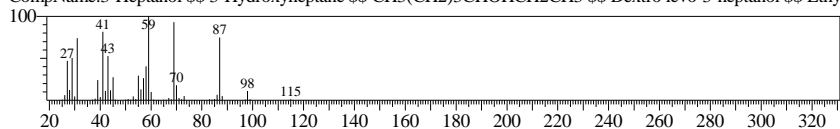
CompName:Dimethyl 3,3'-oxydipropionate



Hit#:6 Entry:3601 Library:NIST08.LIB

SI:63 Formula:C7H16O CAS:589-82-2 MolWeight:116 RetIndex:879

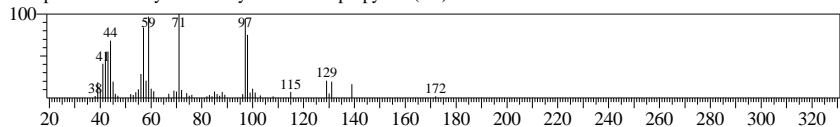
CompName:3-Heptanol \$ 3-Hydroxyheptane \$ CH3(CH2)3CHOHCH2CH3 \$ Dextro levo-3-heptanol \$ Ethy



Hit#:7 Entry:26422 Library:NIST08.LIB

SI:63 Formula:C8H16N2S CAS:0-00-0 MolWeight:172 RetIndex:1138

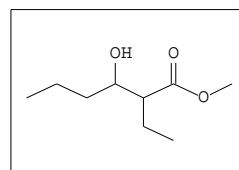
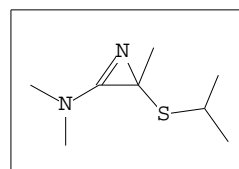
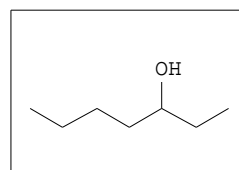
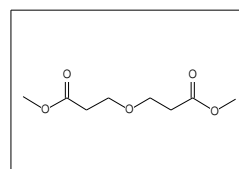
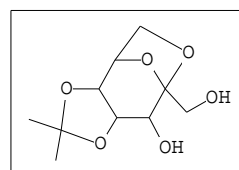
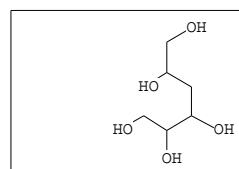
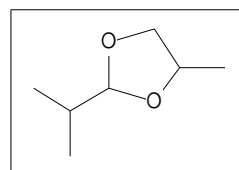
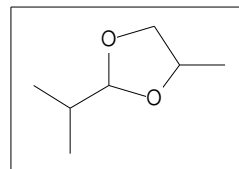
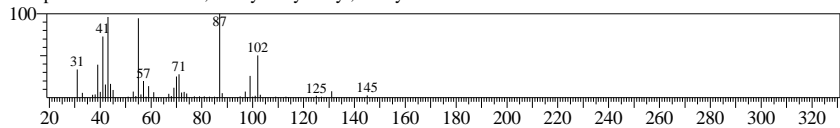
CompName:2-Methyl-3-dimethylamino-2-isopropylthio(2H)azirine



Hit#:8 Entry:27563 Library:NIST08.LIB

SI:62 Formula:C9H18O3 CAS:0-00-0 MolWeight:174 RetIndex:1181

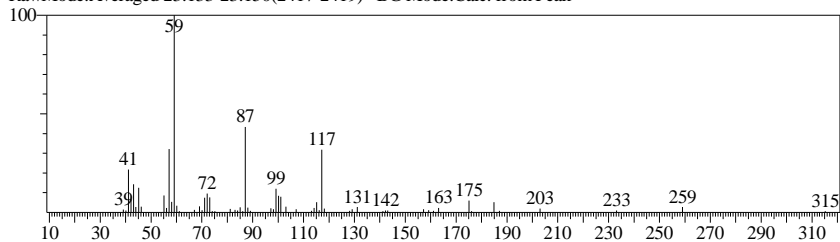
CompName:Hexanoic acid, 2-ethyl-3-hydroxy-, methyl ester



<< Target >>

Line#:35 R.Time:23.142(Scan#:2418) BasePeak:59.10(130174)

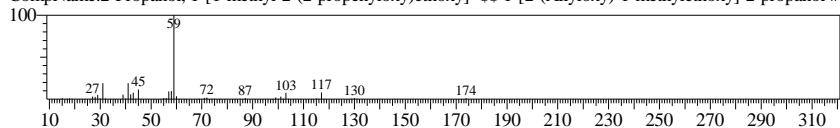
RawMode:Averaged 23.133-23.150(2417-2419) BG Mode:Calc. from Peak



Hit#:1 Entry:27550 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

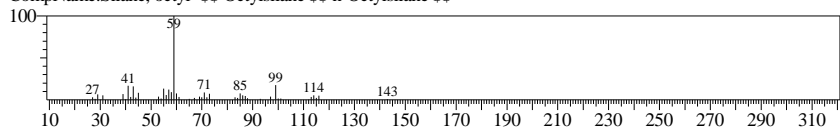
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:2 Entry:12984 Library:NIST08.LIB

SI:77 Formula:C8H20Si CAS:871-92-1 MolWeight:144 RetIndex:0

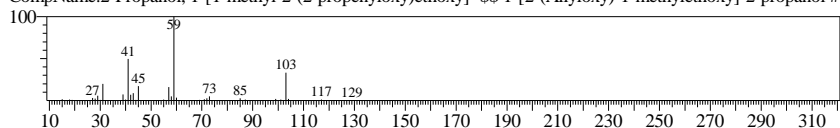
CompName:Silane, octyl- \$- Octylsilane \$- n-Octylsilane \$-



Hit#:3 Entry:12792 Library:NIST08s.LIB

SI:76 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

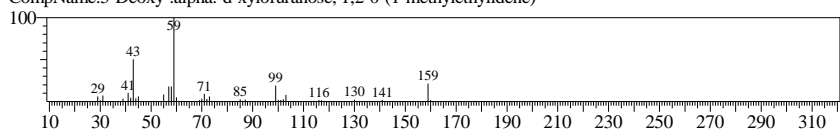
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:27412 Library:NIST08.LIB

SI:76 Formula:C8H14O4 CAS:0-00-0 MolWeight:174 RetIndex:1212

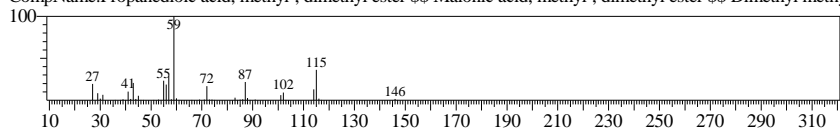
CompName:5-Deoxy-.alpha.-d-xylofuranose, 1,2-O-(1-methylethylidene)-



Hit#:5 Entry:13447 Library:NIST08.LIB

SI:76 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

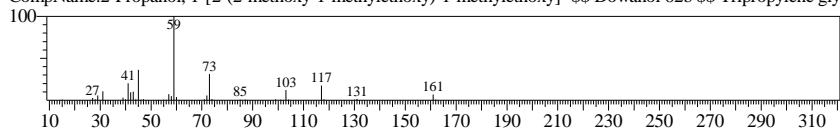
CompName:Propanedioic acid, methyl-, dimethyl ester \$- Malonic acid, methyl-, dimethyl ester \$- Dimethyl meth



Hit#:6 Entry:17306 Library:NIST08s.LIB

SI:76 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

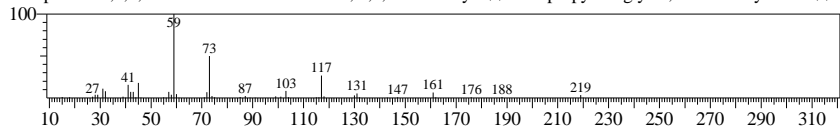
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#:7 Entry:85908 Library:NIST08.LIB

SI:76 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

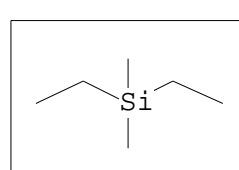
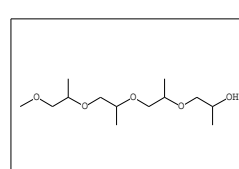
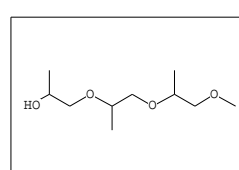
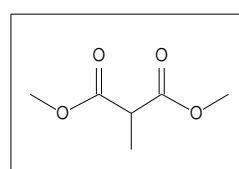
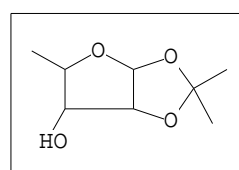
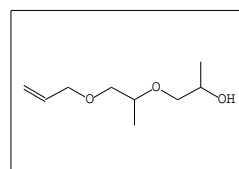
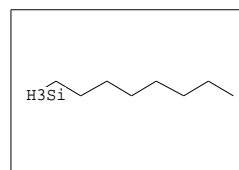
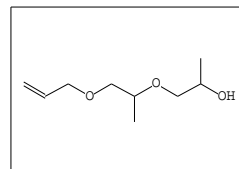
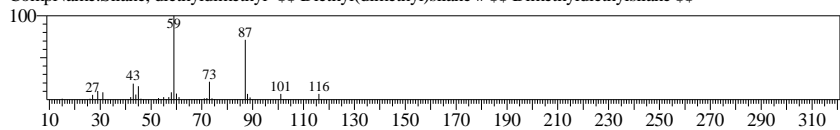
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$- Tetrapropylene glycol, monomethyl ether \$- 4



Hit#:8 Entry:4645 Library:NIST08.LIB

SI:76 Formula:C6H16Si CAS:756-81-0 MolWeight:116 RetIndex:528

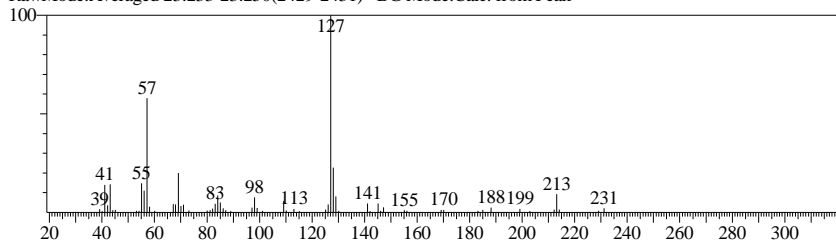
CompName:Silane, diethyldimethyl- \$- Diethyl(dimethyl)silane # \$- Dimethyldiethylsilane \$-



<< Target >>

Line#:36 R.Time:23.242(Scan#:2430) BasePeak:127.15(453308)

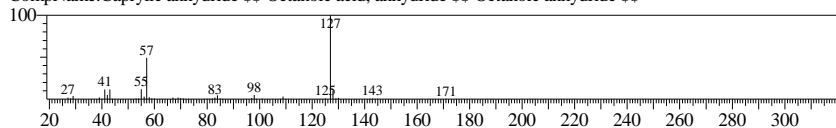
RawMode:Averaged 23.233-23.250(2429-2431) BG Mode:Calc. from Peak



Hit#:1 Entry:90583 Library:NIST08.LIB

SI:84 Formula:C16H30O3 CAS:623-66-5 MolWeight:270 RetIndex:1915

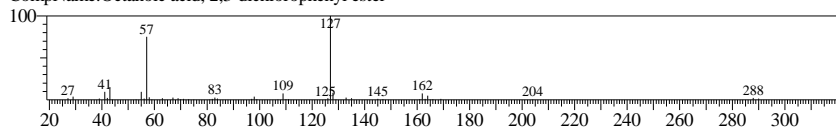
CompName:Caprylic anhydride \$ Octanoic acid, anhydride \$ Octanoic anhydride \$



Hit#:2 Entry:103265 Library:NIST08.LIB

SI:81 Formula:C14H18Cl2O2 CAS:0-00-0 MolWeight:288 RetIndex:2017

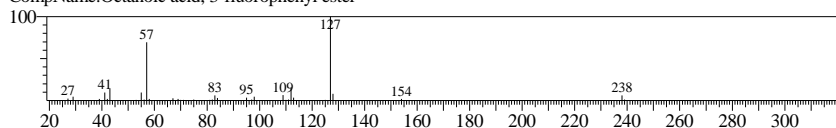
CompName:Octanoic acid, 2,3-dichlorophenyl ester



Hit#:3 Entry:67925 Library:NIST08.LIB

SI:80 Formula:C14H19FO2 CAS:0-00-0 MolWeight:238 RetIndex:1632

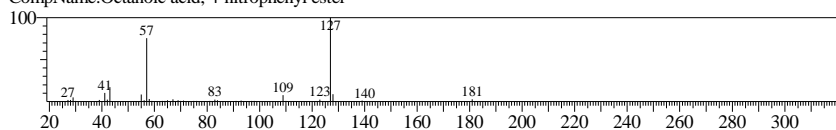
CompName:Octanoic acid, 3-fluorophenyl ester



Hit#:4 Entry:86636 Library:NIST08.LIB

SI:80 Formula:C14H19NO4 CAS:0-00-0 MolWeight:265 RetIndex:2052

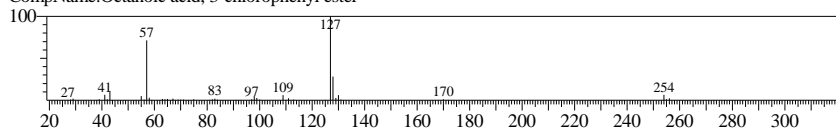
CompName:Octanoic acid, 4-nitrophenyl ester



Hit#:5 Entry:78856 Library:NIST08.LIB

SI:80 Formula:C14H19ClO2 CAS:0-00-0 MolWeight:254 RetIndex:1837

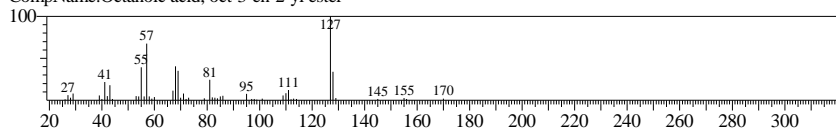
CompName:Octanoic acid, 3-chlorophenyl ester



Hit#:6 Entry:79219 Library:NIST08.LIB

SI:80 Formula:C16H30O2 CAS:0-00-0 MolWeight:254 RetIndex:1723

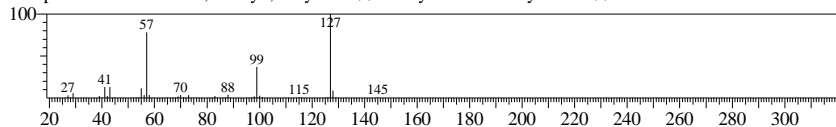
CompName:Octanoic acid, oct-3-en-2-yl ester



Hit#:7 Entry:90584 Library:NIST08.LIB

SI:80 Formula:C16H30O3 CAS:36765-89-6 MolWeight:270 RetIndex:1787

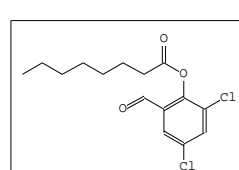
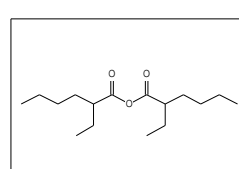
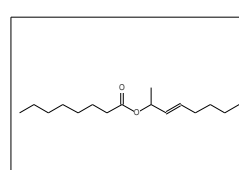
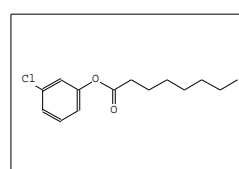
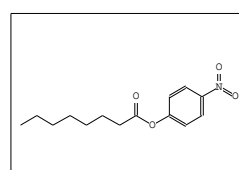
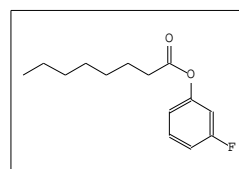
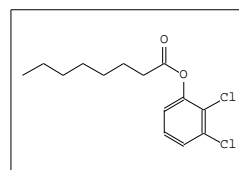
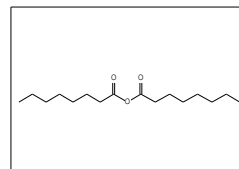
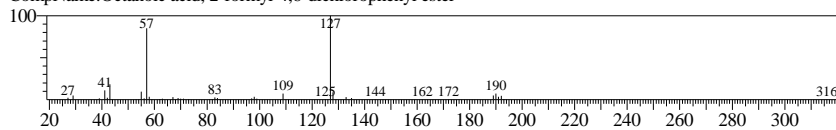
CompName:Hexanoic acid, 2-ethyl-, anhydride \$ 1-Ethylhexanoic anhydride # \$



Hit#:8 Entry:123073 Library:NIST08.LIB

SI:80 Formula:C15H18Cl2O3 CAS:0-00-0 MolWeight:316 RetIndex:2318

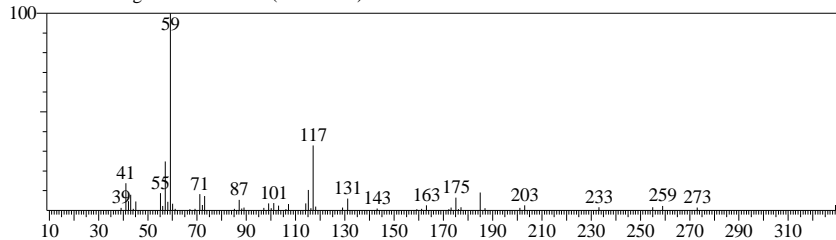
CompName:Octanoic acid, 2-formyl-4,6-dichlorophenyl ester



<< Target >>

Line#:37 R.Time:23.575(Scan#:2470) BasePeak:59.10(127384)

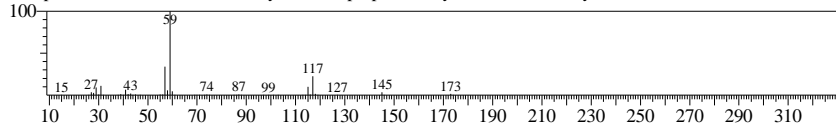
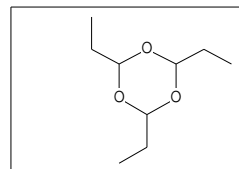
RawMode:Averaged 23.567-23.583(2469-2471) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:80 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

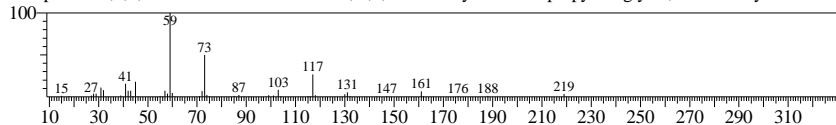
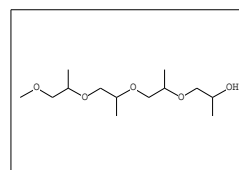
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:85908 Library:NIST08.LIB

SI:78 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

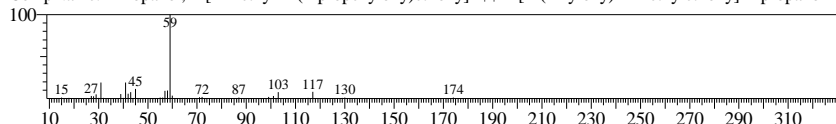
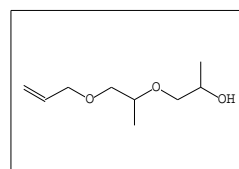
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

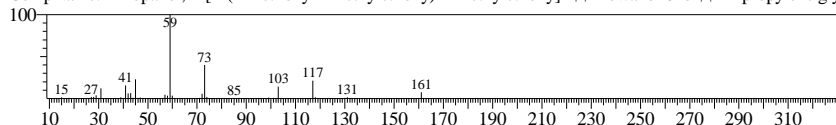
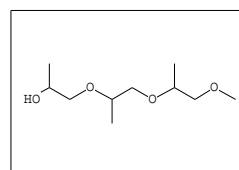
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:17307 Library:NIST08s.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

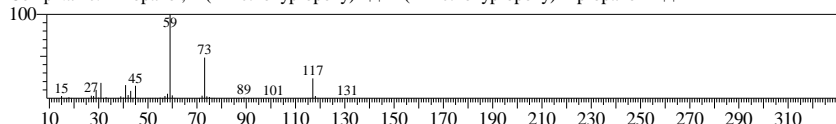
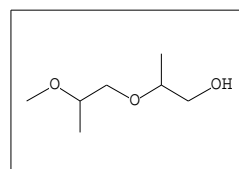
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:5 Entry:14297 Library:NIST08.LIB

SI:77 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

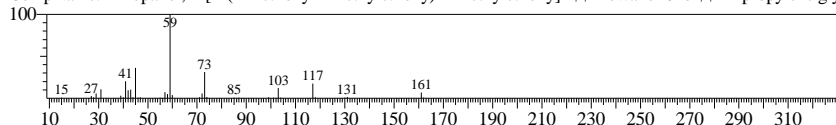
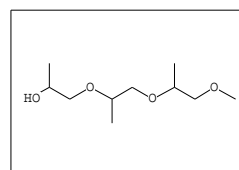
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:6 Entry:17306 Library:NIST08s.LIB

SI:76 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

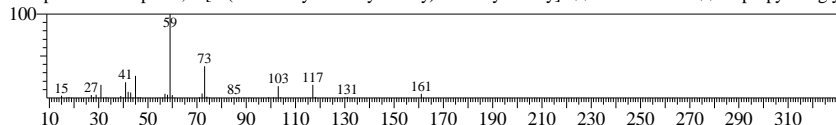
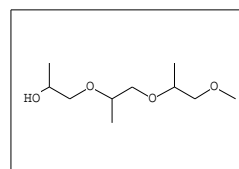
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:76 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

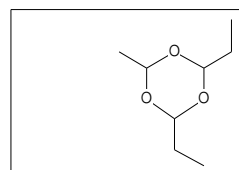
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:19968 Library:NIST08.LIB

SI:76 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

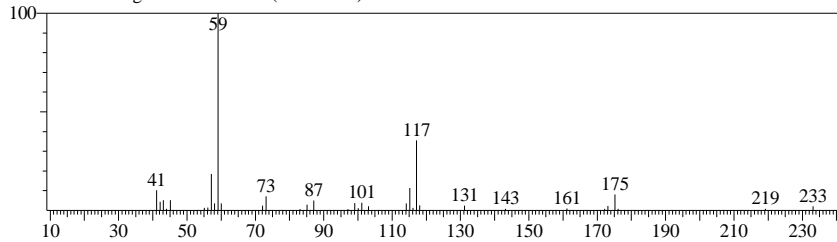
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line#:38 R.Time:23.667(Scan#:2481) BasePeak:59.10(532051)

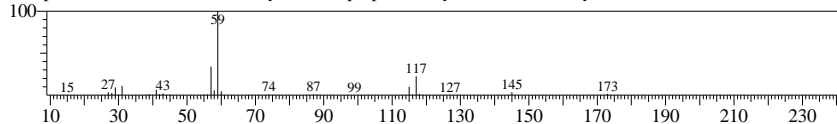
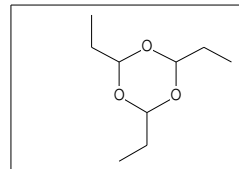
RawMode:Averaged 23.658-23.675(2480-2482) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:85 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

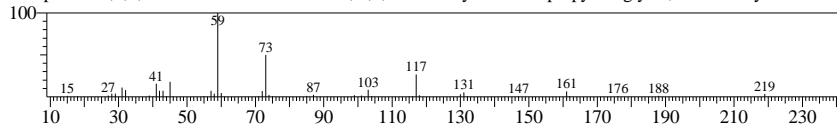
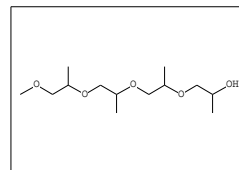
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:85908 Library:NIST08.LIB

SI:82 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

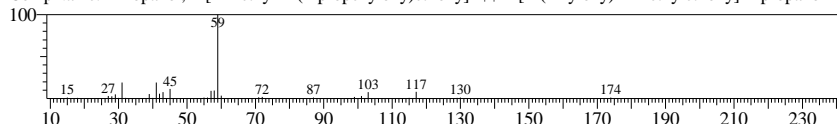
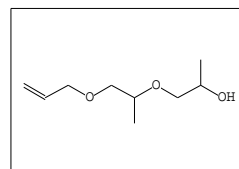
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl-



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

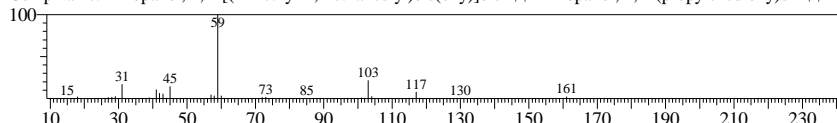
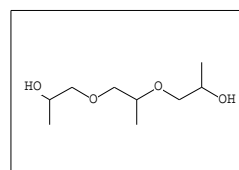
CompName:2-Propanol, 1-[2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:37499 Library:NIST08.LIB

SI:82 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

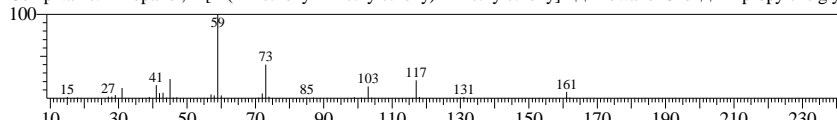
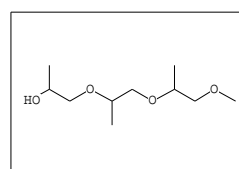
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:5 Entry:17307 Library:NIST08.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

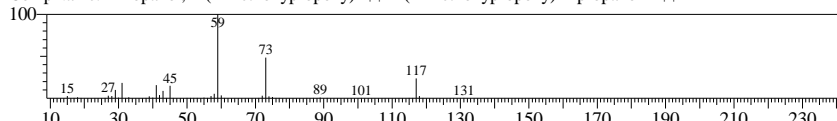
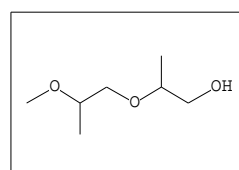
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:14297 Library:NIST08.LIB

SI:81 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

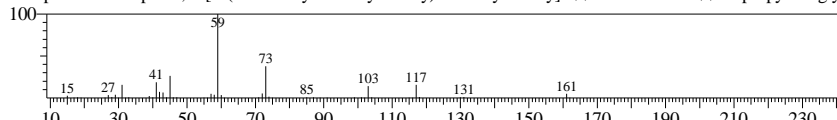
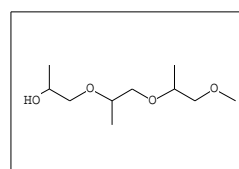
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

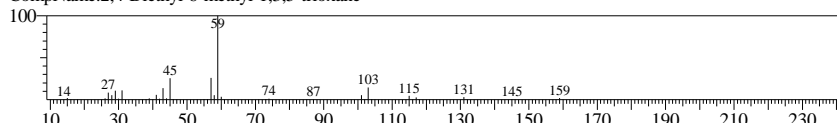
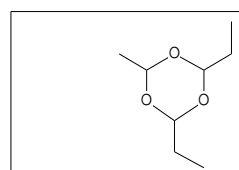
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:19968 Library:NIST08.LIB

SI:80 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

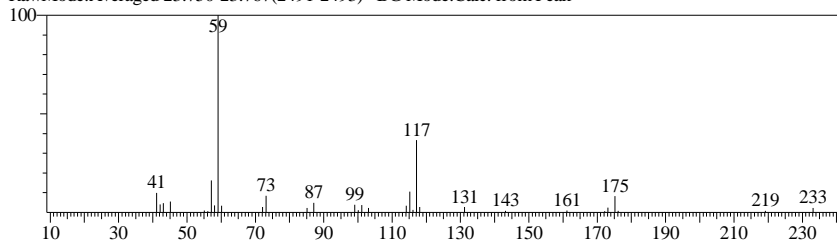
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line#:39 R.Time:23.758(Scan#:2492) BasePeak:59.10(656002)

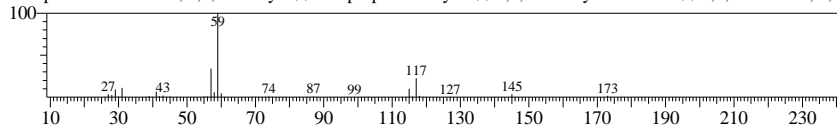
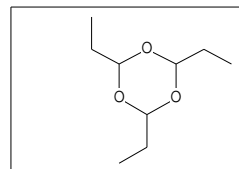
RawMode:Averaged 23.750-23.767(2491-2493) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:85 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

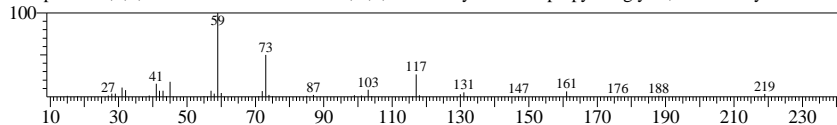
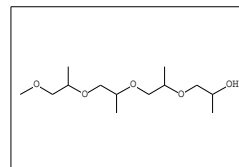
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:85908 Library:NIST08.LIB

SI:83 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

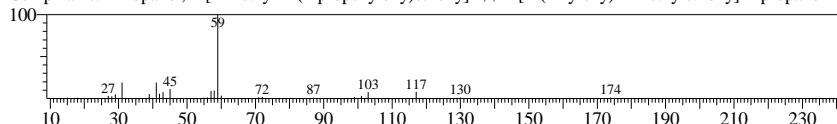
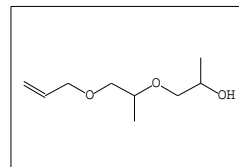
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

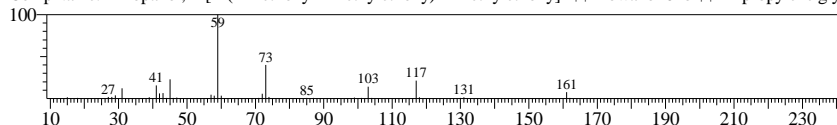
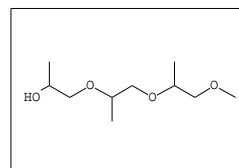
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:17307 Library:NIST08s.LIB

SI:82 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

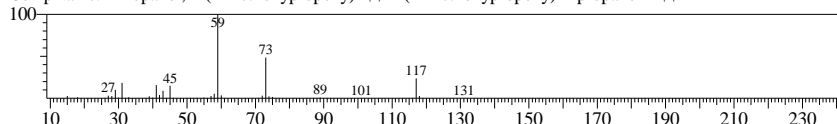
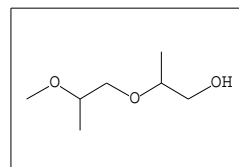
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:5 Entry:14297 Library:NIST08.LIB

SI:82 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

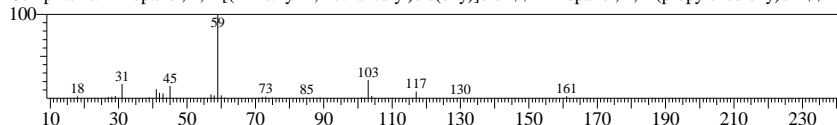
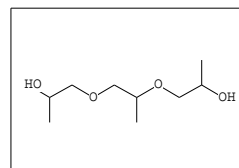
CompName:1-Propanol, 1-[1-(1-methoxypropoxy)-2-(2-methoxypropoxy)-1-propanol #



Hit#:6 Entry:37499 Library:NIST08.LIB

SI:81 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

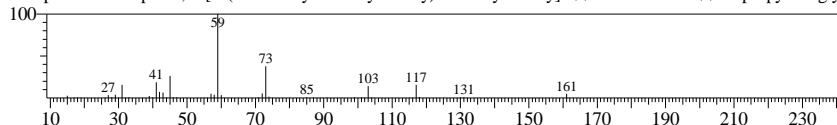
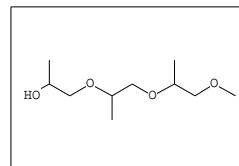
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

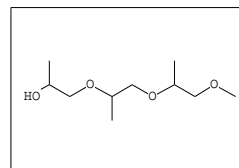
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:17306 Library:NIST08s.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

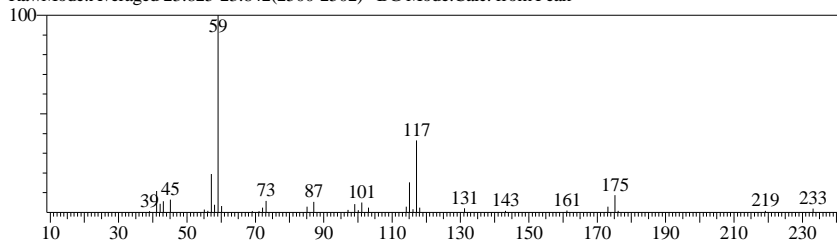
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line#:40 R.Time:23.833(Scan#:2501) BasePeak:59.10(458829)

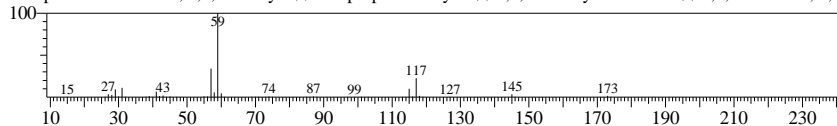
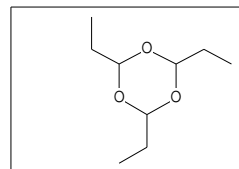
RawMode:Averaged 23.825-23.842(2500-2502) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:85 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

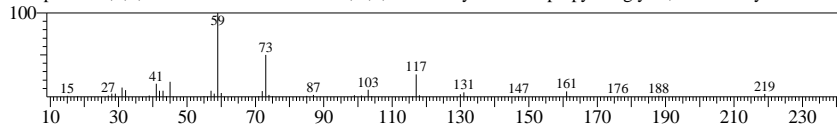
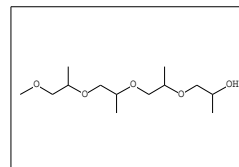
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:85908 Library:NIST08.LIB

SI:82 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

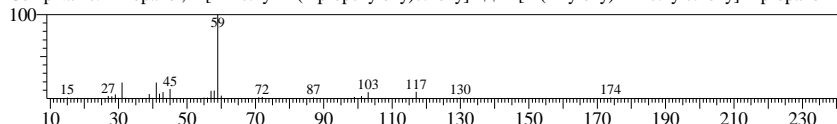
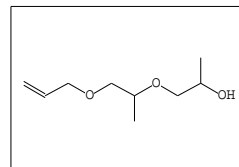
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl-



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

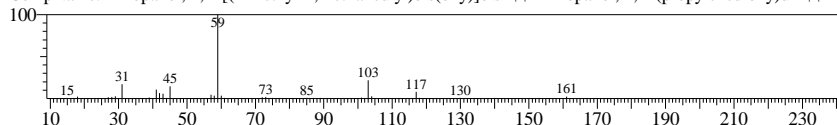
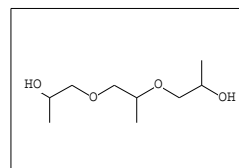
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:37499 Library:NIST08.LIB

SI:81 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

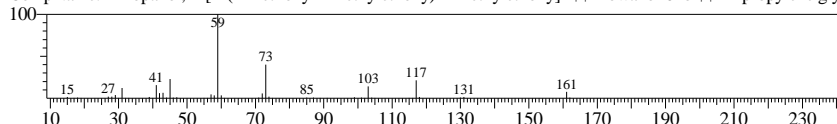
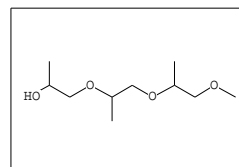
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ 2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis-



Hit#:5 Entry:17307 Library:NIST08.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

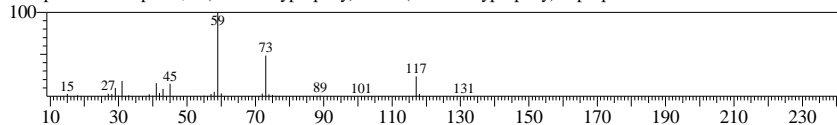
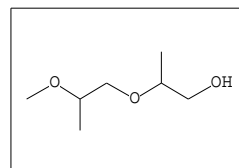
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:6 Entry:14297 Library:NIST08.LIB

SI:81 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

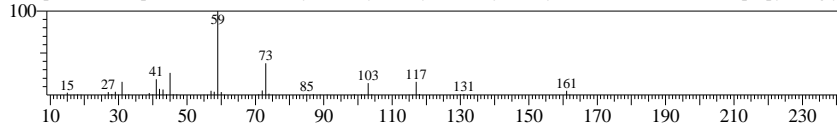
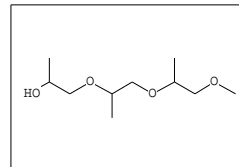
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

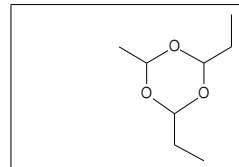
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:19968 Library:NIST08.LIB

SI:80 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

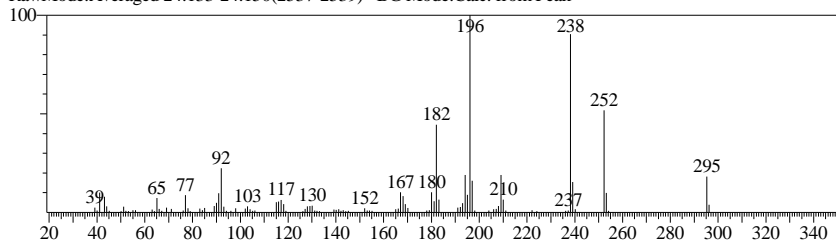
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line#:41 R.Time:24.142(Scan#:2538) BasePeak:196.10(66211)

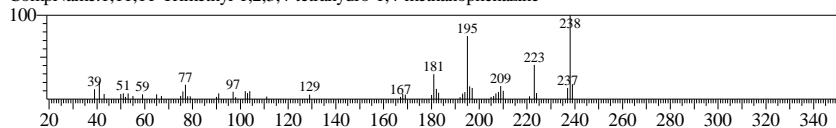
RawMode:Averaged 24.133-24.150(2537-2539) BG Mode:Calc. from Peak



Hit#:1 Entry:68171 Library:NIST08.LIB

SI:61 Formula:C16H18N2 CAS:0-00-0 MolWeight:238 RetIndex:1812

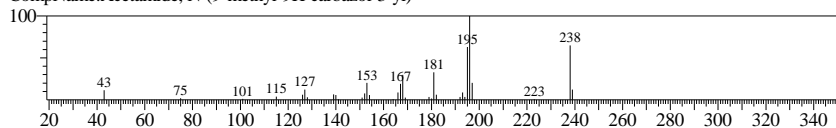
CompName:1,11,11-Trimethyl-1,2,3,4-tetrahydro-1,4-methanophenazine



Hit#:2 Entry:68009 Library:NIST08.LIB

SI:61 Formula:C15H14N2O CAS:0-00-0 MolWeight:238 RetIndex:2219

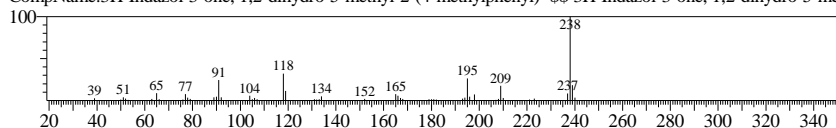
CompName:Acetamide, N-(9-methyl-9H-carbazol-3-yl)-



Hit#:3 Entry:68017 Library:NIST08.LIB

SI:59 Formula:C15H14N2O CAS:17049-55-7 MolWeight:238 RetIndex:2313

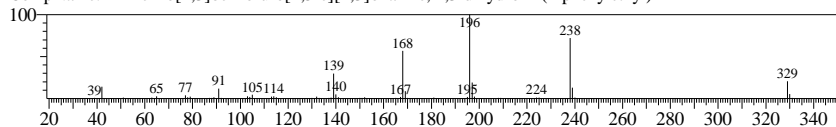
CompName:3H-Indazol-3-one, 1,2-dihydro-5-methyl-2-(4-methylphenyl)-



Hit#:4 Entry:132150 Library:NIST08.LIB

SI:58 Formula:C22H19NO2 CAS:0-00-0 MolWeight:329 RetIndex:2720

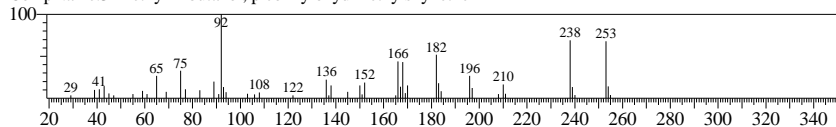
CompName:1H-Benzofuro[2,3]benzofuro[4,5-e][1,3]oxazine, 2,3-dihydro-2-(2-phenylethyl)-



Hit#:5 Entry:78200 Library:NIST08.LIB

SI:57 Formula:C13H23NO2Si CAS:0-00-0 MolWeight:253 RetIndex:1481

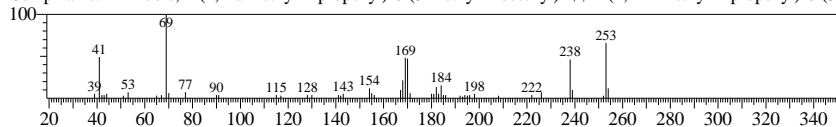
CompName:3-Methyl-1-butanol, picolinoyloxymethyl ether



Hit#:6 Entry:78406 Library:NIST08.LIB

SI:57 Formula:C18H23N CAS:55045-00-6 MolWeight:253 RetIndex:2063

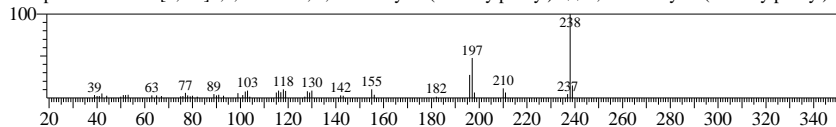
CompName:1H-Indole, 2-(1,1-dimethyl-2-propenyl)-6-(3-methyl-2-butenyl)-



Hit#:7 Entry:67916 Library:NIST08.LIB

SI:56 Formula:C14H14N4 CAS:61802-70-8 MolWeight:238 RetIndex:2060

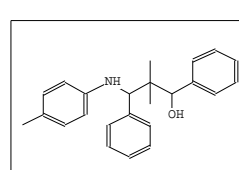
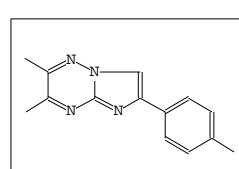
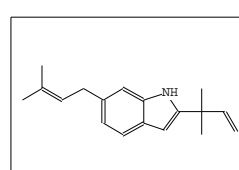
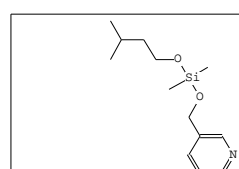
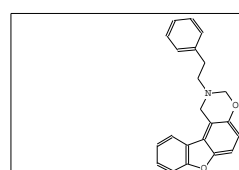
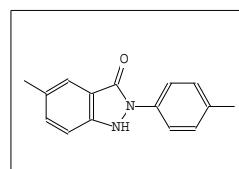
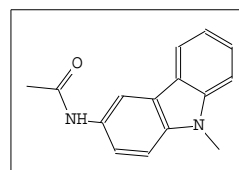
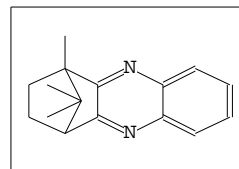
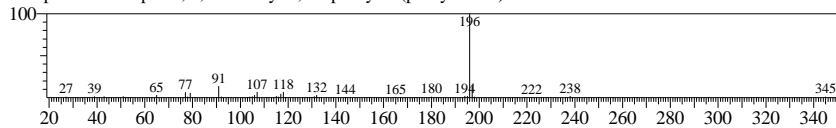
CompName:Imidazo[1,2-b]1,2,4-triazine, 2,3-dimethyl-6-(4-methylphenyl)-



Hit#:8 Entry:142282 Library:NIST08.LIB

SI:56 Formula:C24H27NO CAS:0-00-0 MolWeight:345 RetIndex:2863

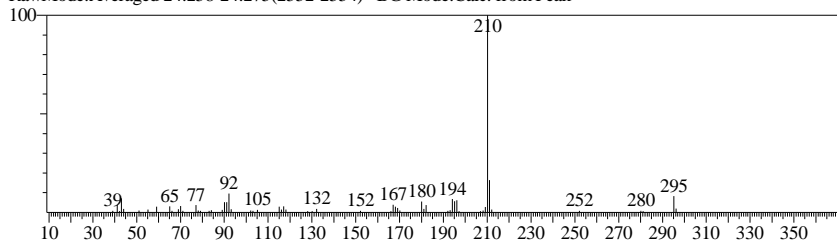
CompName:1-Propanol, 2,2-dimethyl-1,3-diphenyl-3-(p-tolylamino)-



<< Target >>

Line#:42 R.Time:24.267(Scan#:2553) BasePeak:210.10(96850)

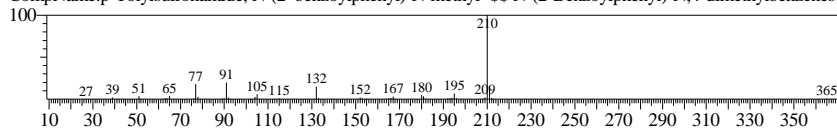
RawMode:Averaged 24.258-24.275(2552-2554) BG Mode:Calc. from Peak



Hit#:1 Entry:152984 Library:NIST08.LIB

SI:78 Formula:C21H19NO3S CAS:1859-73-0 MolWeight:365 RetIndex:3058

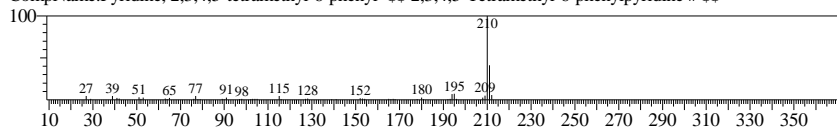
CompName:p-Tolylsulfonamide, N-(2'-benzoylphenyl)-N-methyl- \$\$ N-(2-Benzoylphenyl)-N,4-dimethylbenzenes



Hit#:2 Entry:50086 Library:NIST08.LIB

SI:76 Formula:C15H17N CAS:80206-51-5 MolWeight:211 RetIndex:1814

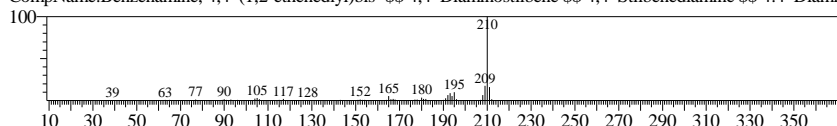
CompName:Pyridine, 2,3,4,5-tetramethyl-6-phenyl- \$\$ 2,3,4,5-Tetramethyl-6-phenylpyridine # \$\$



Hit#:3 Entry:49526 Library:NIST08.LIB

SI:76 Formula:C14H14N2 CAS:621-96-5 MolWeight:210 RetIndex:2197

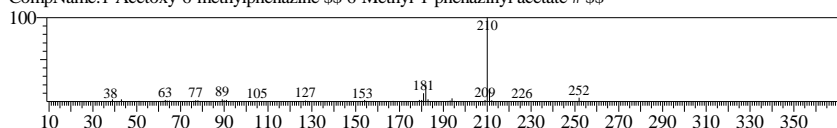
CompName:Benzenamine, 4,4'-(1,2-ethenediyl)bis- \$\$ 4,4'-Diaminostilbene \$\$ 4,4'-Diami



Hit#:4 Entry:77607 Library:NIST08.LIB

SI:74 Formula:C15H12N2O2 CAS:14031-09-5 MolWeight:252 RetIndex:2263

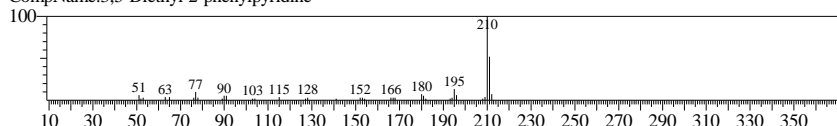
CompName:1-Acetoxy-6-methylphenazine \$\$ 6-Methyl-1-phenazinyl acetate # \$\$



Hit#:5 Entry:50085 Library:NIST08.LIB

SI:74 Formula:C15H17N CAS:73669-43-9 MolWeight:211 RetIndex:1786

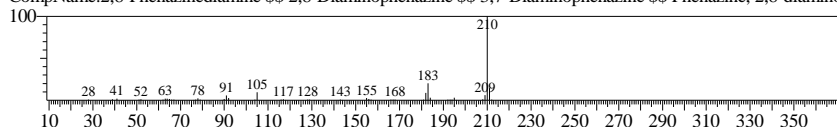
CompName:3,5-Diethyl-2-phenylpyridine



Hit#:6 Entry:49240 Library:NIST08.LIB

SI:74 Formula:C12H10N4 CAS:7704-40-7 MolWeight:210 RetIndex:2393

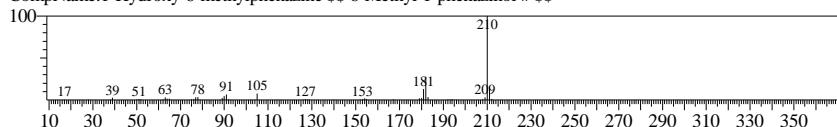
CompName:2,8-Phenazinediamine \$\$ 2,8-Diaminophenazine \$\$ 3,7-Diaminophenazine \$\$ Phenazine, 2,8-diaminc



Hit#:7 Entry:49380 Library:NIST08.LIB

SI:74 Formula:C13H10N2O CAS:14031-08-4 MolWeight:210 RetIndex:2104

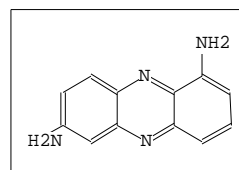
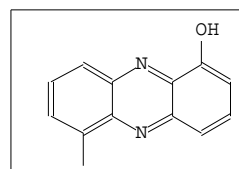
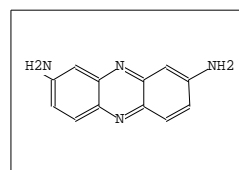
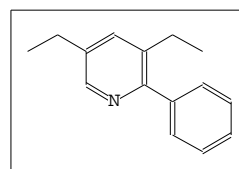
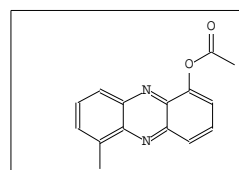
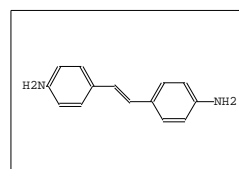
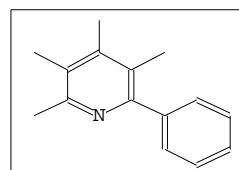
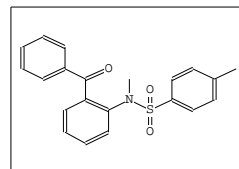
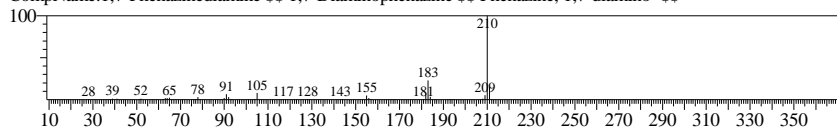
CompName:1-Hydroxy-6-methylphenazine \$\$ 6-Methyl-1-phenazinol # \$\$



Hit#:8 Entry:49241 Library:NIST08.LIB

SI:73 Formula:C12H10N4 CAS:28124-29-0 MolWeight:210 RetIndex:2393

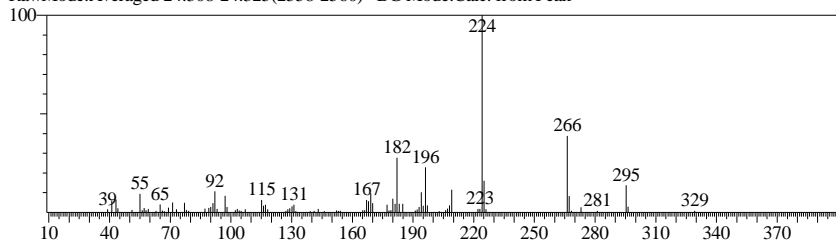
CompName:1,7-Phenazinediamine \$\$ 1,7-Diaminophenazine \$\$ Phenazine, 1,7-diamino- \$\$



<< Target >>

Line#:43 R.Time:24.317(Scan#:2559) BasePeak:224.10(85774)

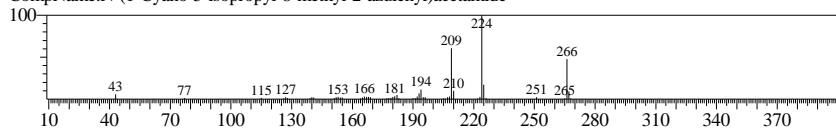
RawMode:Averaged 24.308-24.325(2558-2560) BG Mode:Calc. from Peak



Hit#:1 Entry:87607 Library:NIST08.LIB

SI:69 Formula:C17H18N2O CAS:93648-58-9 MolWeight:266 RetIndex:2307

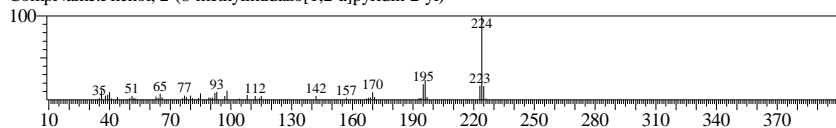
CompName:N-(1-Cyano-5-isopropyl-8-methyl-2-azulenyl)acetamide



Hit#:2 Entry:58742 Library:NIST08.LIB

SI:61 Formula:C14H12N2O CAS:0-00-0 MolWeight:224 RetIndex:2002

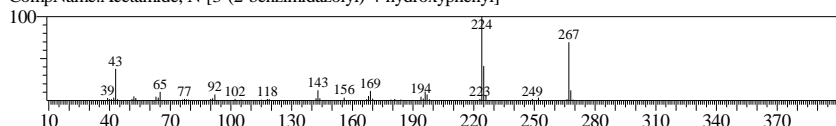
CompName:Phenol, 2-(8-methylimidazo[1,2-a]pyridin-2-yl)-



Hit#:3 Entry:88153 Library:NIST08.LIB

SI:60 Formula:C15H13N3O2 CAS:0-00-0 MolWeight:267 RetIndex:2768

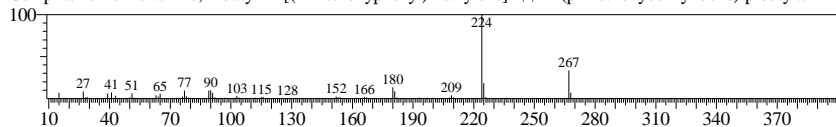
CompName:Acetamide, N-[3-(2-benzimidazolyl)-4-hydroxyphenyl]-



Hit#:4 Entry:88336 Library:NIST08.LIB

SI:60 Formula:C18H21NO CAS:26227-73-6 MolWeight:267 RetIndex:2283

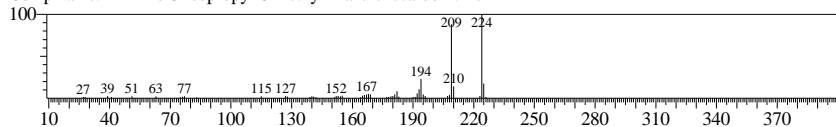
CompName:Benzenamine, 4-butyl-N-[(4-methoxyphenyl)methylene]- N-(p-Methoxybenzylidene)-p-butylaniline



Hit#:5 Entry:58863 Library:NIST08.LIB

SI:60 Formula:C15H16N2 CAS:93946-48-6 MolWeight:224 RetIndex:1951

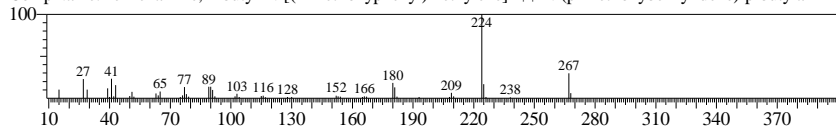
CompName:2-Amino-5-isopropyl-8-methyl-1-azulenecarbonitrile



Hit#:6 Entry:22530 Library:NIST08s.LIB

SI:59 Formula:C18H21NO CAS:26227-73-6 MolWeight:267 RetIndex:2283

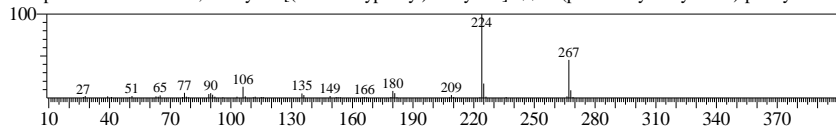
CompName:Benzenamine, 4-butyl-N-[(4-methoxyphenyl)methylene]- N-(p-Methoxybenzylidene)-p-butylaniline



Hit#:7 Entry:22531 Library:NIST08s.LIB

SI:59 Formula:C18H21NO CAS:26227-73-6 MolWeight:267 RetIndex:2283

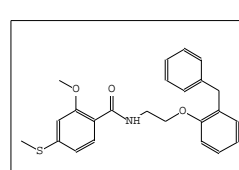
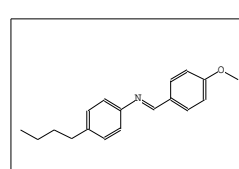
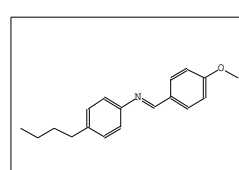
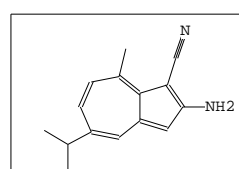
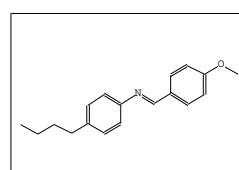
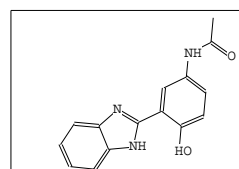
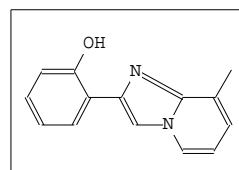
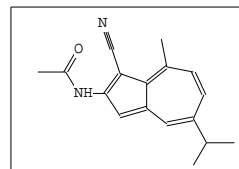
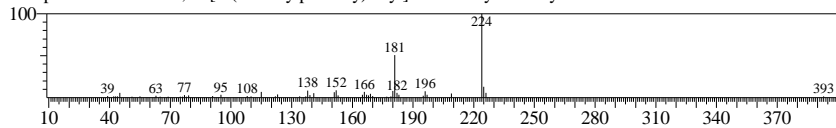
CompName:Benzenamine, 4-butyl-N-[(4-methoxyphenyl)methylene]- N-(p-Methoxybenzylidene)-p-butylaniline



Hit#:8 Entry:169852 Library:NIST08.LIB

SI:59 Formula:C24H25NO3S CAS:0-00-0 MolWeight:407 RetIndex:3481

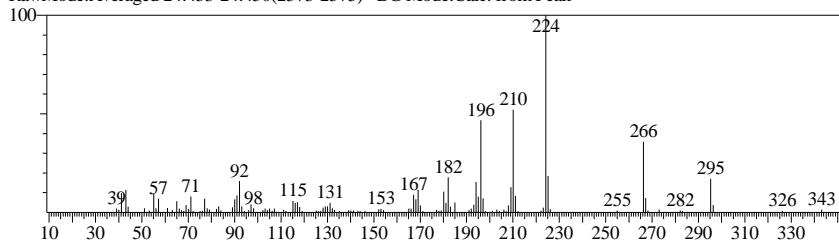
CompName:Benamide, N-[2-(2-benzylphenoxy)ethyl]-2-methoxy-4-methylthio-



<< Target >>

Line#:44 R.Time:24.442(Scan#:2574) BasePeak:224.10(64529)

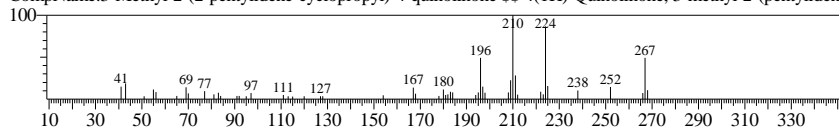
RawMode:Averaged 24.433-24.450(2573-2575) BG Mode:Calc. from Peak



Hit#:1 Entry:88335 Library:NIST08.LIB

SI:71 Formula:C18H21NO CAS:68978-12-1 MolWeight:267 RetIndex:2305

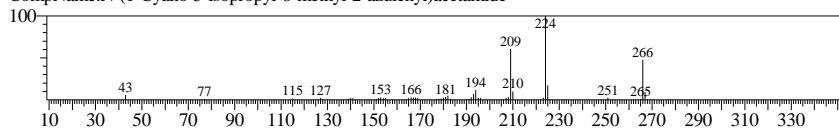
CompName:3-Methyl-2-(2-pentylidene-cyclopropyl)-4-quinolinone



Hit#:2 Entry:87607 Library:NIST08.LIB

SI:65 Formula:C17H18N2O CAS:93648-58-9 MolWeight:266 RetIndex:2307

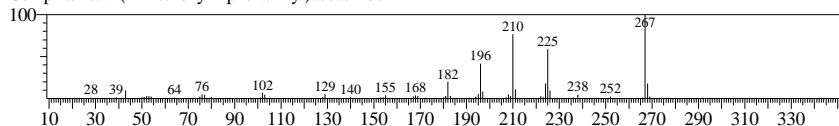
CompName:N-(1-Cyano-5-isopropyl-8-methyl-2-azulenyl)acetamide



Hit#:3 Entry:88162 Library:NIST08.LIB

SI:60 Formula:C15H13N3O2 CAS:28103-04-0 MolWeight:267 RetIndex:2572

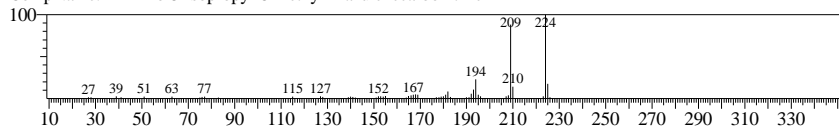
CompName:N-(4-Methoxy-1-phenazinyl)acetamide



Hit#:4 Entry:58863 Library:NIST08.LIB

SI:60 Formula:C15H16N2 CAS:93946-48-6 MolWeight:224 RetIndex:1951

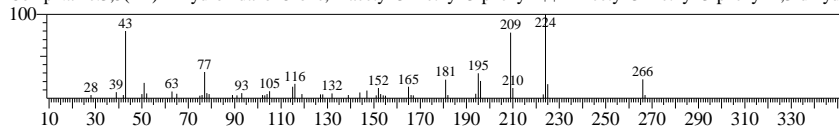
CompName:2-Amino-5-isopropyl-8-methyl-1-azulenecarbonitrile



Hit#:5 Entry:87462 Library:NIST08.LIB

SI:58 Formula:C16H14N2O2 CAS:59341-21-8 MolWeight:266 RetIndex:2251

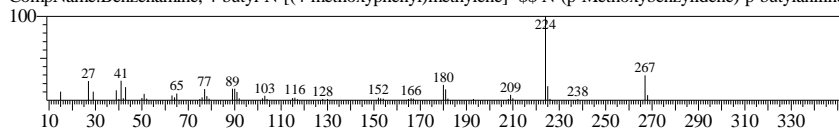
CompName:3,5(2H)-Dihydroindazol-5-one, 2-acetyl-3-methyl-3-phenyl-2,3-dihyd



Hit#:6 Entry:22530 Library:NIST08s.LIB

SI:58 Formula:C18H21NO CAS:26227-73-6 MolWeight:267 RetIndex:2283

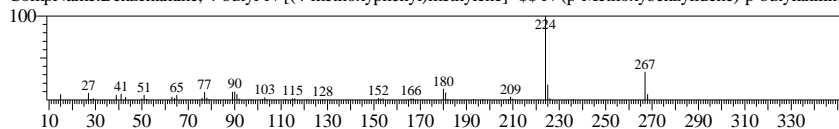
CompName:Benzenamine, 4-butyl-N-[(4-methoxyphenyl)methylene]-



Hit#:7 Entry:88336 Library:NIST08.LIB

SI:58 Formula:C18H21NO CAS:26227-73-6 MolWeight:267 RetIndex:2283

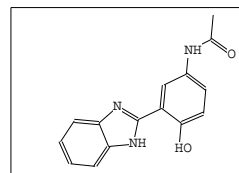
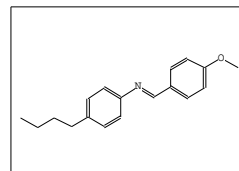
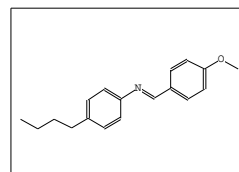
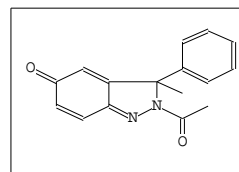
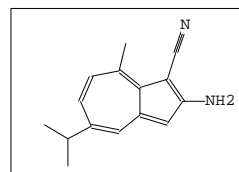
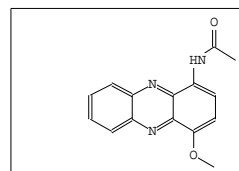
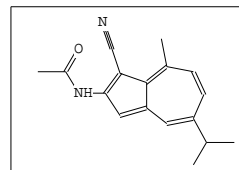
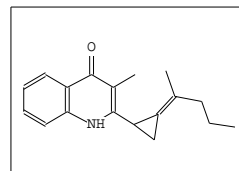
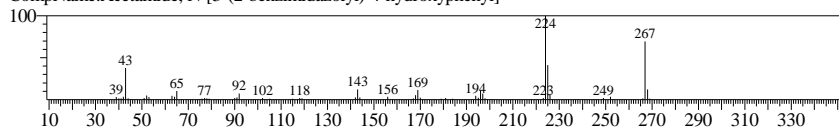
CompName:Benzenamine, 4-butyl-N-[(4-methoxyphenyl)methylene]-



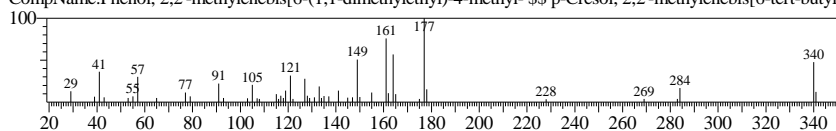
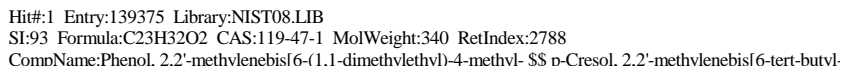
Hit#:8 Entry:88153 Library:NIST08.LIB

SI:57 Formula:C15H13N3O2 CAS:0-00-0 MolWeight:267 RetIndex:2768

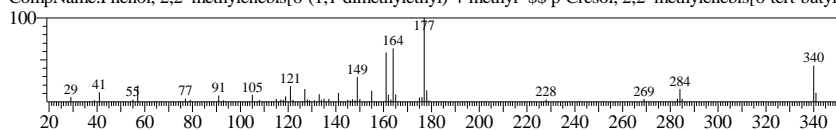
CompName:Acetamide, N-[3-(2-benzimidazolyl)-4-hydroxyphenyl]-



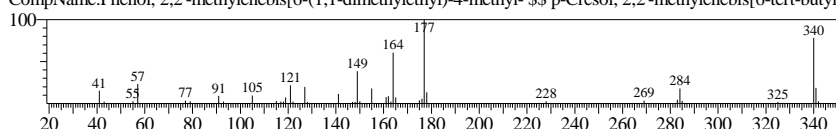
Line#:45 R.Time:24.508(Scan#:2582) BasePeak:177.10(132928)
RawMode:Averaged 24.500-24.517(2581-2583) BG Mode:Calc. from Peak



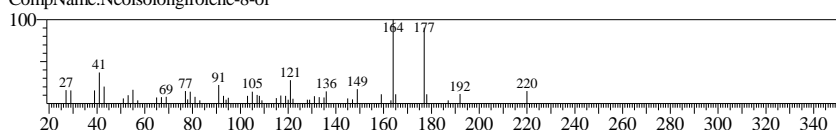
Hit#:2 Entry:25940 Library:NIST08s.LIB
SI:90 Formula:C23H32O2 CAS:119-47-1 MolWeight:340 RetIndex:2788
CompName:Phenol, 2,2'-methylenebis[6-(1,1-dimethylethyl)-4-methyl-5-*n*-Cresol, 2,2'-methylenebis[6-tert-butyl-



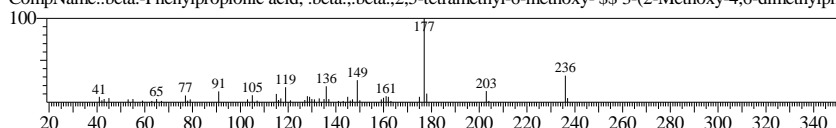
Hit#3 Entry:25941 Library:NIST08s.LIB
 SI:84 Formula:C23H32O2 CAS:119-47-1 MolWeight:340 RefIndex:2788
 ComName:Phenol 2,2'-methylenebis[6-(1,1-dimethylethyl)-4-methyl-5-n-Cresol 2,2'-methylenebis[6-tert-butyl-



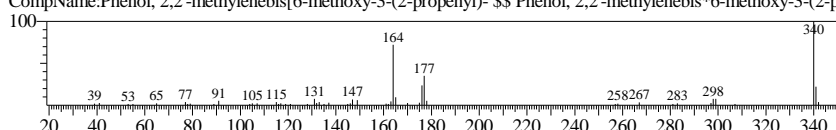
Hit#4 Entry:56010 Library:NIST08.LIB
SI:61 Formula:C15H24O CAS:0-00-0 MolWeight:220 RetIndex:1604
CompName:Neoisolongifolene-8-ol



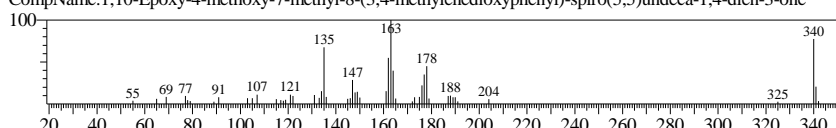
Hit#5 Entry:66607 Library:NIST08.LIB
 SI:61 Formula:C14H20O3 CAS:186317-88-4 MolWeight:236 RefIndex:1879
 CompName: beta -Phenylpropionic acid, beta, beta, 2,5-tetramethyl-6-methoxy-, \$S\$ 3-(2-Methoxy-4,6-dimethylph



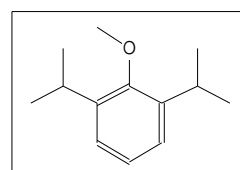
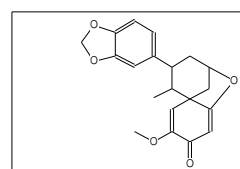
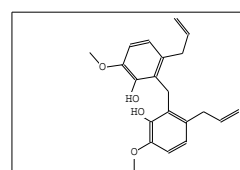
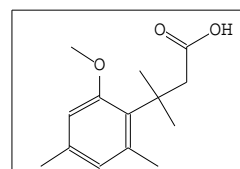
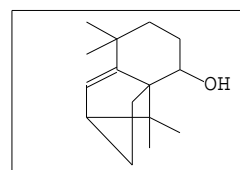
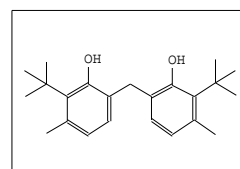
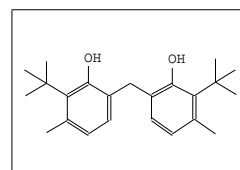
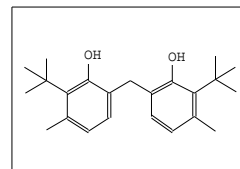
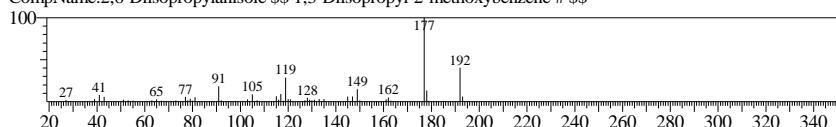
Hit#:6 Entry:139219 Library:NIST08.LIB
 SI:61 Formula:C21H24O4 CAS:55334-55-9 MolWeight:340 RetIndex:2890
 CompName:Phenol 2,2'-methylenebis[6-methoxy-3-(2-propenyl)-] \$\$_{2}\$\$ Phenol 2,2'-methylenebis[6-methoxy-3-(2-propenyl)-]



Hit#:7 Entry:139005 Library:NIST08.LIB
 SI:60 Formula:C20H20O5 CAS:20156-97-2 MolWeight:340 RetIndex:2614
 CompName:1,10-Epoxy-4-methoxy-7-methyl-8-(3,4-methylenedioxyphenyl)-spiro[5.5]undeca-1,4-dien-3-one



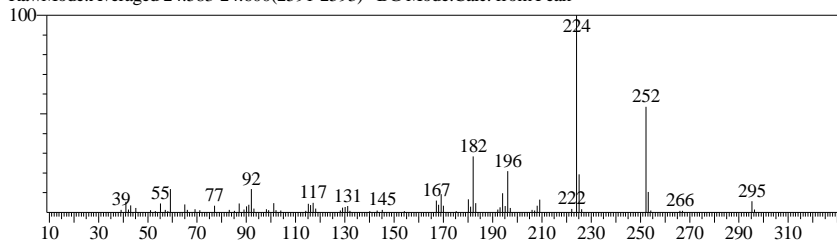
Hit#:8 Entry:37992 Library:NIST08.LIB
 SI:58 Formula:C13H20O CAS:2944-52-7 MolWeight:192 RetIndex:1365
 CompName:2,6-Diisopropylanisole \$\$ 1,3-Diisopropyl-2-methoxybenzene # \$\$



<< Target >>

Line#:46 R.Time:24.592(Scan#:2592) BasePeak:224.10(74119)

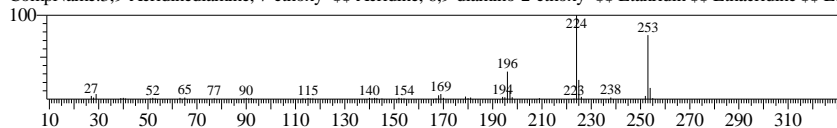
RawMode:Averaged 24.583-24.600(2591-2593) BG Mode:Calc. from Peak



Hit#:1 Entry:78306 Library:NIST08.LIB

SI:67 Formula:C15H15N3O CAS:442-16-0 MolWeight:253 RetIndex:2688

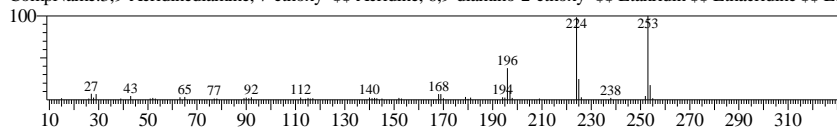
CompName:3,9-Acridinediamine, 7-ethoxy- \$\$ Acridine, 6,9-diamino-2-ethoxy- \$\$ Etakridin \$\$ Ethacridine \$\$ Et



Hit#:2 Entry:21604 Library:NIST08s.LIB

SI:66 Formula:C15H15N3O CAS:442-16-0 MolWeight:253 RetIndex:2688

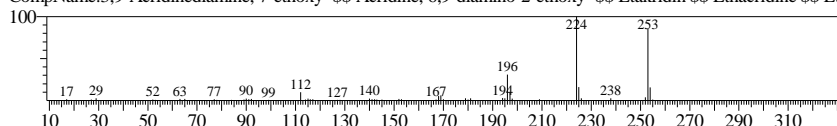
CompName:3,9-Acridinediamine, 7-ethoxy- \$\$ Acridine, 6,9-diamino-2-ethoxy- \$\$ Etakridin \$\$ Ethacridine \$\$ Et



Hit#:3 Entry:21603 Library:NIST08s.LIB

SI:65 Formula:C15H15N3O CAS:442-16-0 MolWeight:253 RetIndex:2688

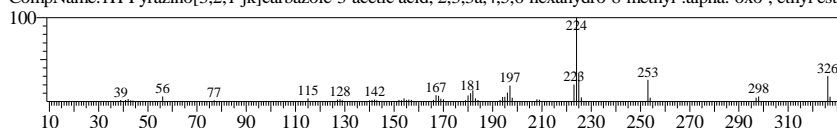
CompName:3,9-Acridinediamine, 7-ethoxy- \$\$ Acridine, 6,9-diamino-2-ethoxy- \$\$ Etakridin \$\$ Ethacridine \$\$ Et



Hit#:4 Entry:129874 Library:NIST08.LIB

SI:61 Formula:C19H22N2O3 CAS:0-00-0 MolWeight:326 RetIndex:2486

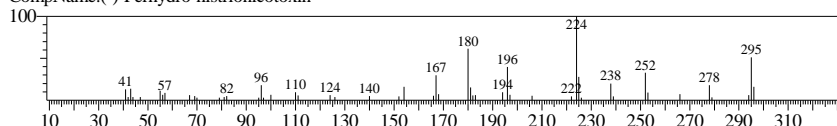
CompName:1H-Pyrazino[3,2,1-jk]carbazole-3-acetic acid, 2,3,3a,4,5,6-hexahydro-8-methyl-.alpha.-oxo-, ethyl est



Hit#:5 Entry:108468 Library:NIST08.LIB

SI:60 Formula:C19H37NO CAS:0-00-0 MolWeight:295 RetIndex:2385

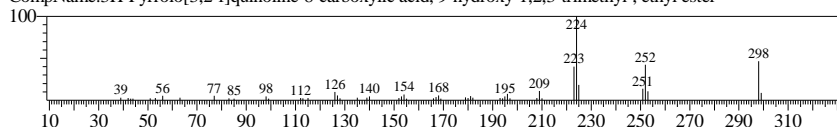
CompName:(-)-Perhydro-histronicotxin



Hit#:6 Entry:110494 Library:NIST08.LIB

SI:60 Formula:C17H18N2O3 CAS:0-00-0 MolWeight:298 RetIndex:2566

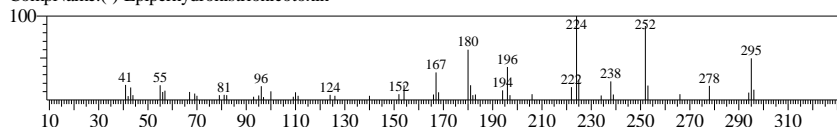
CompName:3H-Pyrrolo[3,2-f]quinoline-8-carboxylic acid, 9-hydroxy-1,2,3-trimethyl-, ethyl ester



Hit#:7 Entry:108469 Library:NIST08.LIB

SI:60 Formula:C19H37NO CAS:0-00-0 MolWeight:295 RetIndex:2385

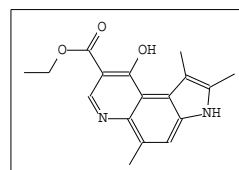
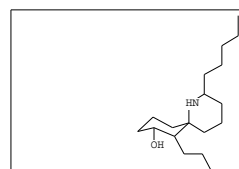
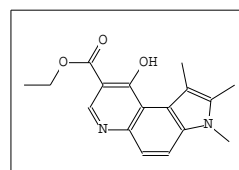
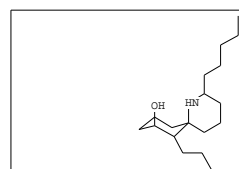
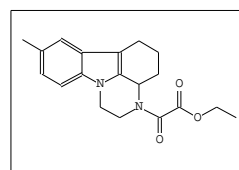
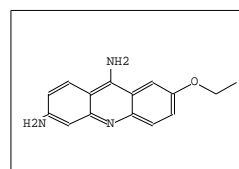
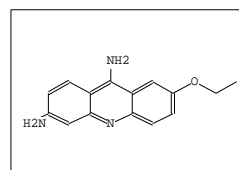
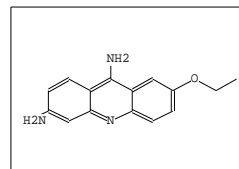
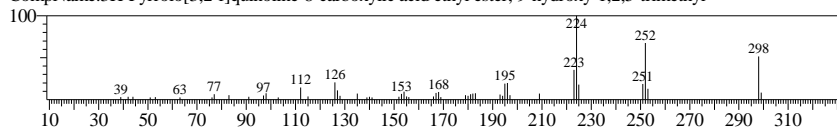
CompName:(-)-Epiperhydrohistrionicotxin



Hit#:8 Entry:110491 Library:NIST08.LIB

SI:60 Formula:C17H18N2O3 CAS:0-00-0 MolWeight:298 RetIndex:2711

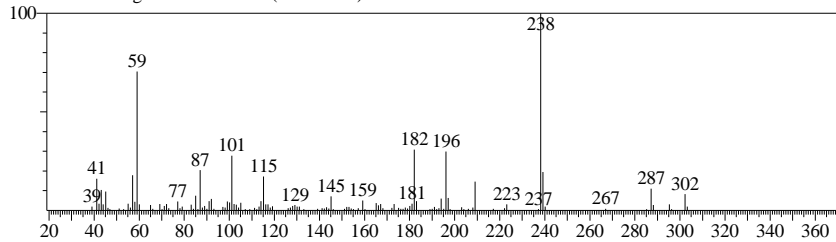
CompName:3H-Pyrrolo[3,2-f]quinoline-8-carboxylic acid ethyl ester, 9-hydroxy-1,2,5-trimethyl-



<< Target >>

Line#:47 R.Time:24.650(Scan#:2599) BasePeak:238.15(47507)

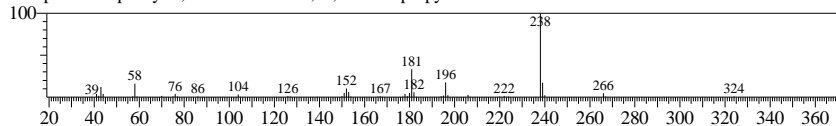
RawMode:Averaged 24.642-24.658(2598-2600) BG Mode:Calc. from Peak



Hit#:1 Entry:128703 Library:NIST08.LIB

SI:56 Formula:C20H24N2O2 CAS:348605-63-0 MolWeight:324 RetIndex:2862

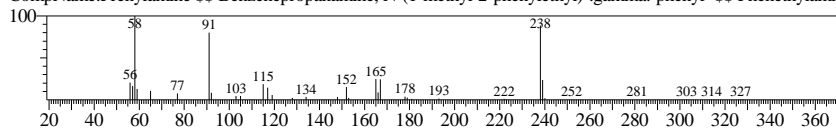
CompName:Biphenyl-2,2'-dicarboxamide, N,N'-bis(isopropyl-



Hit#:2 Entry:25564 Library:NIST08.LIB

SI:53 Formula:C24H27N CAS:390-64-7 MolWeight:329 RetIndex:2707

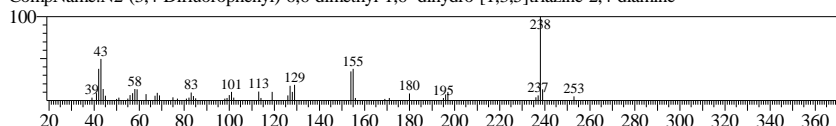
CompName:Prenylamine Phenethylamine, N-(1-methyl-2-phenylethyl)-, gamma-phenyl- Phenethylamine



Hit#:3 Entry:78043 Library:NIST08.LIB

SI:52 Formula:C11H13F2N5 CAS:0-00-0 MolWeight:253 RetIndex:2144

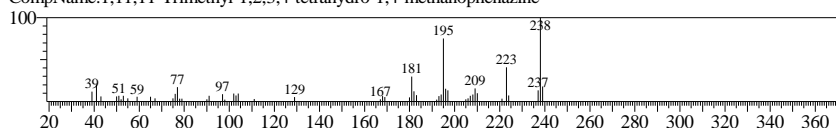
CompName:N2-(3,4-Difluorophenyl)-6,6-dimethyl-1,6-dihydro-[1,3,5]triazine-2,4-diamine



Hit#:4 Entry:68171 Library:NIST08.LIB

SI:52 Formula:C16H18N2 CAS:0-00-0 MolWeight:238 RetIndex:1812

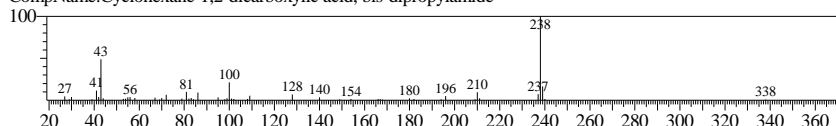
CompName:1,11,11-Trimethyl-1,2,3,4-tetrahydro-1,4-methanophenazine



Hit#:5 Entry:137865 Library:NIST08.LIB

SI:51 Formula:C20H38N2O2 CAS:0-00-0 MolWeight:338 RetIndex:2436

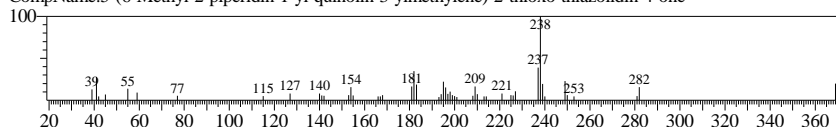
CompName:Cyclohexane-1,2-dicarboxylic acid, bis-dipropylamide



Hit#:6 Entry:154943 Library:NIST08.LIB

SI:51 Formula:C19H19N3OS2 CAS:0-00-0 MolWeight:369 RetIndex:3354

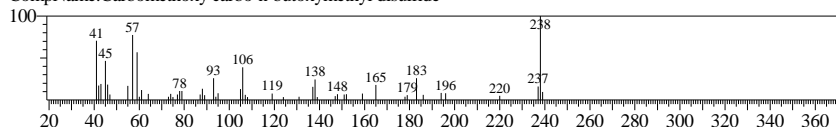
CompName:5-(6-Methyl-piperidin-1-yl-quinolin-3-ylmethylene)-2-thioxo-thiazolidin-4-one



Hit#:7 Entry:67436 Library:NIST08.LIB

SI:51 Formula:C8H14O4S2 CAS:0-00-0 MolWeight:238 RetIndex:1653

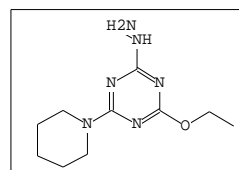
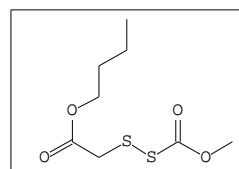
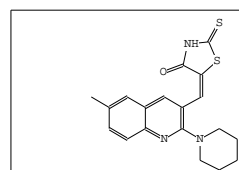
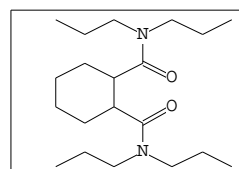
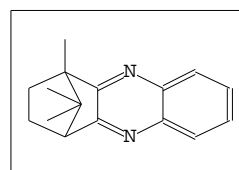
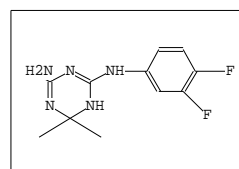
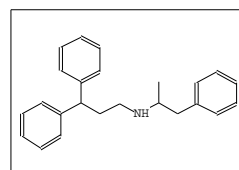
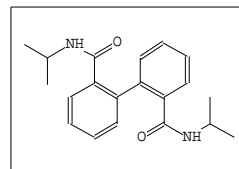
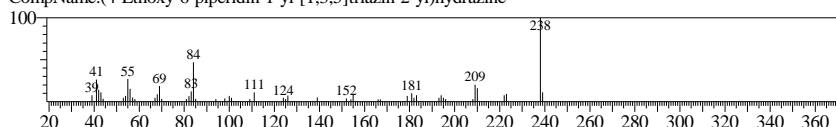
CompName:Carbomethoxy carbo-n-butoxymethyl disulfide



Hit#:8 Entry:67559 Library:NIST08.LIB

SI:51 Formula:C10H18N6O CAS:0-00-0 MolWeight:238 RetIndex:2131

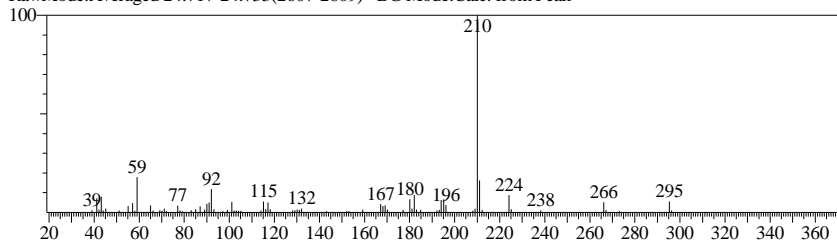
CompName:(4-Ethoxy-6-piperidin-1-yl-[1,3,5]triazin-2-yl)hydrazine



<< Target >>

Line#:48 R.Time:24.725(Scan#:2608) BasePeak:210.10(170049)

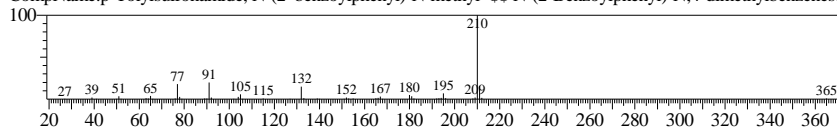
RawMode:Averaged 24.717-24.733(2607-2609) BG Mode:Calc. from Peak



Hit#:1 Entry:152984 Library:NIST08.LIB

SI:68 Formula:C21H19NO3S CAS:1859-73-0 MolWeight:365 RetIndex:3058

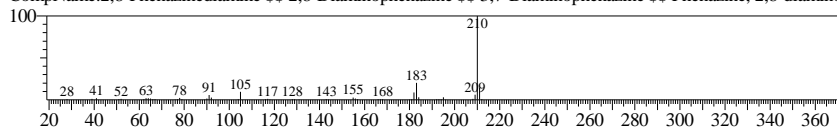
CompName:p-Tolylsulfonamide, N-(2'-benzoylphenyl)-N-methyl- \$N\$-(2-Benzoylphenyl)-N,4-dimethylbenzenesulfonamide



Hit#:2 Entry:49240 Library:NIST08.LIB

SI:68 Formula:C12H10N4 CAS:7704-40-7 MolWeight:210 RetIndex:2393

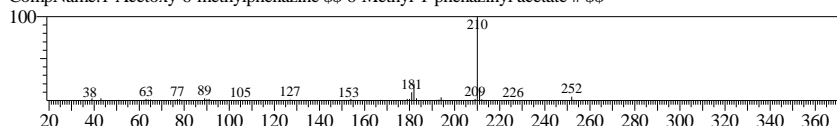
CompName:2,8-Phenazinediamine \$2,8\$-Diaminophenazine \$3,7\$-Diaminophenazine \$Phenazine, 2,8\$-diamino



Hit#:3 Entry:77607 Library:NIST08.LIB

SI:67 Formula:C15H12N2O2 CAS:14031-09-5 MolWeight:252 RetIndex:2263

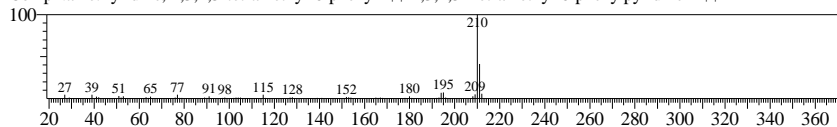
CompName:1-Acetoxy-6-methylphenazine \$6\$-Methyl-1-phenazinyl acetate # \$



Hit#:4 Entry:50086 Library:NIST08.LIB

SI:66 Formula:C15H17N CAS:80206-51-5 MolWeight:211 RetIndex:1814

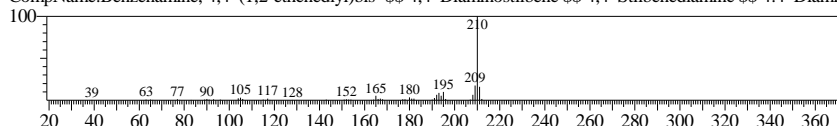
CompName:Pyridine, 2,3,4,5-tetramethyl-6-phenyl- \$2,3,4,5\$-Tetramethyl-6-phenylpyridine # \$



Hit#:5 Entry:49526 Library:NIST08.LIB

SI:66 Formula:C14H14N2 CAS:621-96-5 MolWeight:210 RetIndex:2197

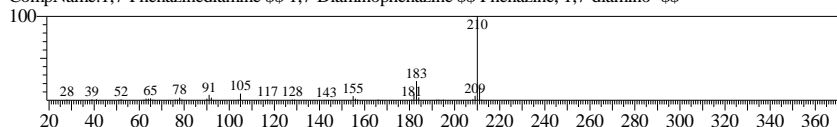
CompName:Benzenamine, 4,4'-(1,2-ethenediyl)bis- \$4,4'\$-Diaminostilbene \$4,4'\$-Diaminostilbene \$4,4'\$-Diaminostilbene



Hit#:6 Entry:49241 Library:NIST08.LIB

SI:66 Formula:C12H10N4 CAS:28124-29-0 MolWeight:210 RetIndex:2393

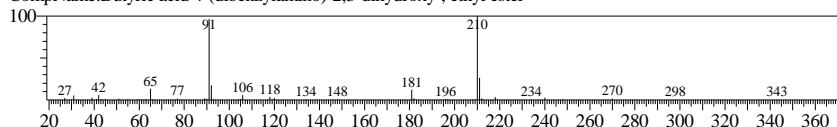
CompName:1,7-Phenazinediamine \$1,7\$-Diaminophenazine \$Phenazine, 1,7\$-diamino- \$



Hit#:7 Entry:141001 Library:NIST08.LIB

SI:66 Formula:C20H25NO4 CAS:0-00-0 MolWeight:343 RetIndex:2704

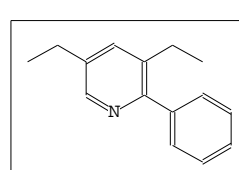
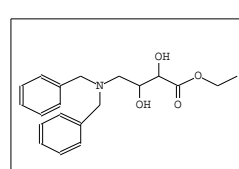
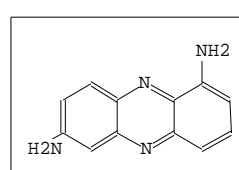
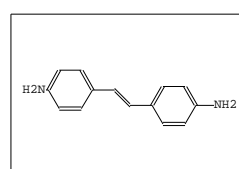
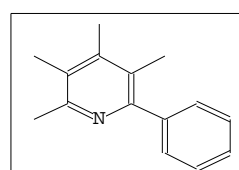
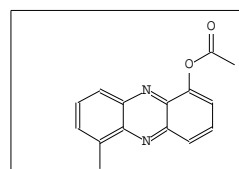
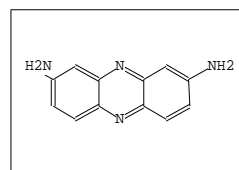
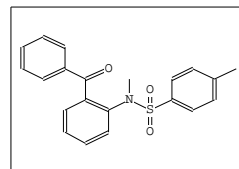
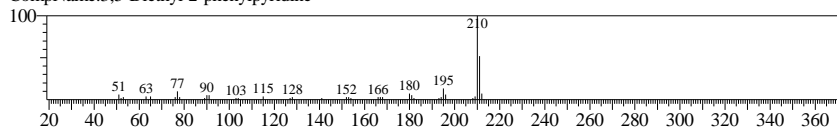
CompName:Butyric acid 4-(dibenzylamino)-2,3-dihydroxy-, ethyl ester



Hit#:8 Entry:50085 Library:NIST08.LIB

SI:66 Formula:C15H17N CAS:73669-43-9 MolWeight:211 RetIndex:1786

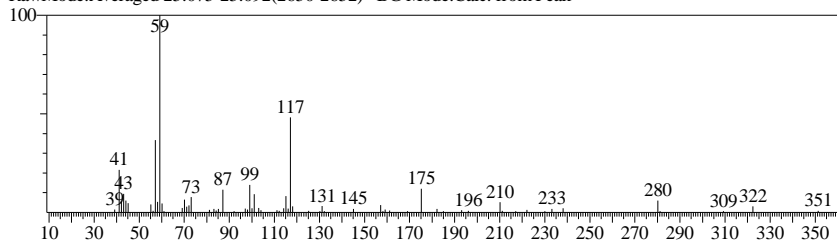
CompName:3,5-Diethyl-2-phenylpyridine



<< Target >>

Line#:49 R.Time:25.083(Scan#:2651) BasePeak:59.10(43505)

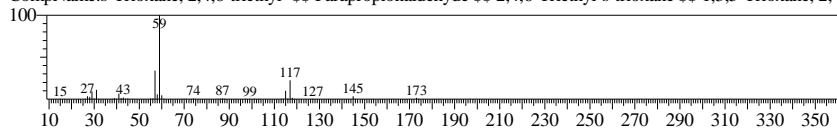
RawMode:Averaged 25.075-25.092(2650-2652) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

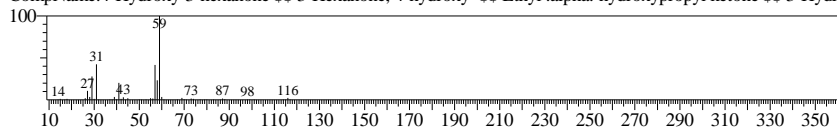
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:4568 Library:NIST08.LIB

SI:75 Formula:C6H12O2 CAS:4984-85-4 MolWeight:116 RetIndex:916

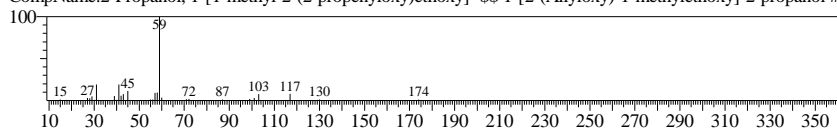
CompName:4-Hydroxy-3-hexanone \$\$ 3-Hexanone, 4-hydroxy- Ethyl .alpha.-hydroxypropyl ketone \$\$ 3-Hydr



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:75 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

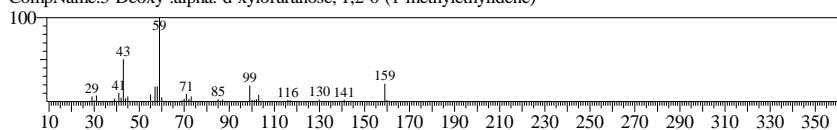
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:27412 Library:NIST08.LIB

SI:75 Formula:C8H14O4 CAS:0-00-0 MolWeight:174 RetIndex:1212

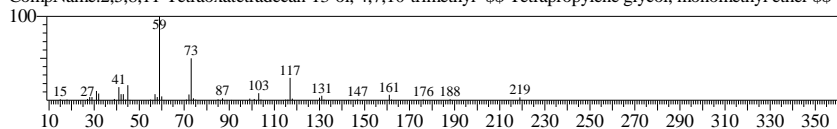
CompName:5-Deoxy-.alpha.-d-xylofuranose, 1,2-O-(1-methylethylidene)-



Hit#:5 Entry:85908 Library:NIST08.LIB

SI:75 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

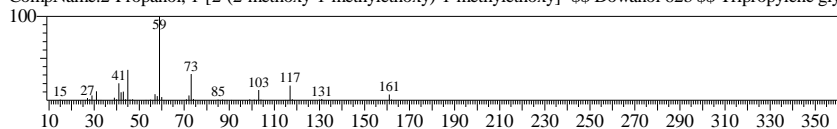
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#:6 Entry:17306 Library:NIST08s.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

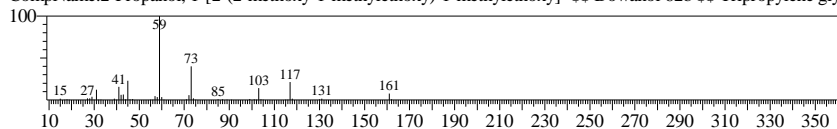
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:7 Entry:17307 Library:NIST08s.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

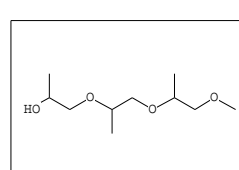
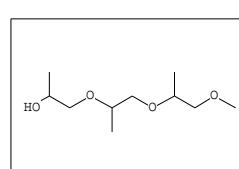
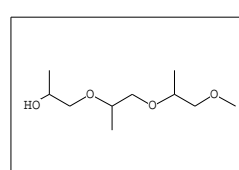
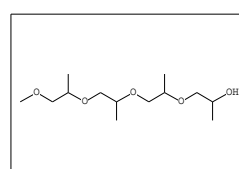
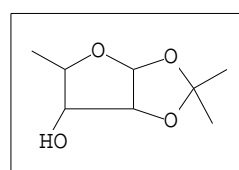
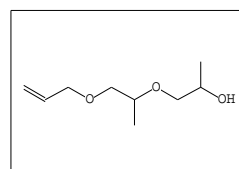
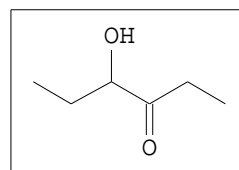
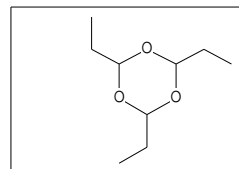
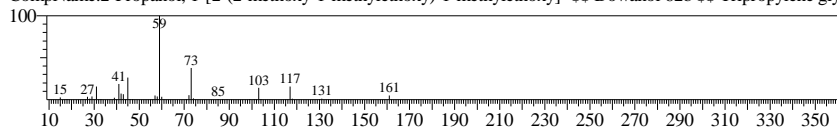
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:46317 Library:NIST08.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

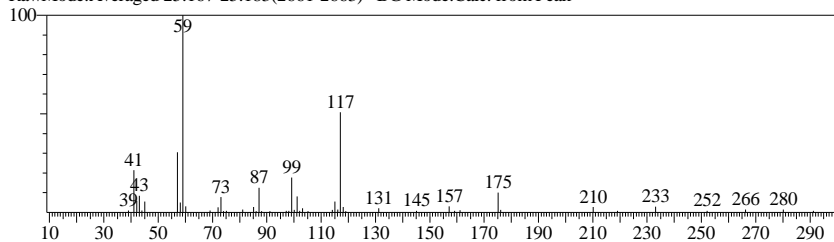
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line# 50 R.Time:25.175(Scan#:2662) BasePeak:59.10(88664)

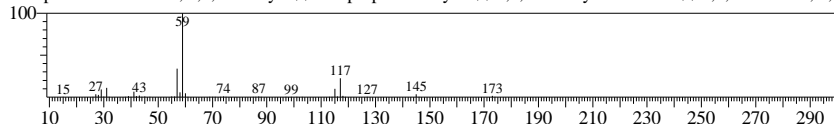
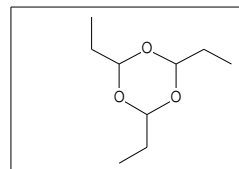
RawMode:Averaged 25.167-25.183(2661-2663) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:80 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

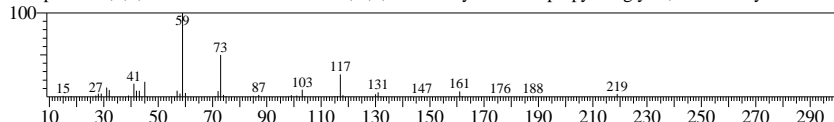
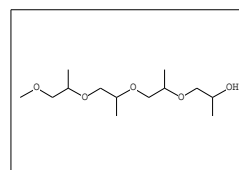
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:79 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

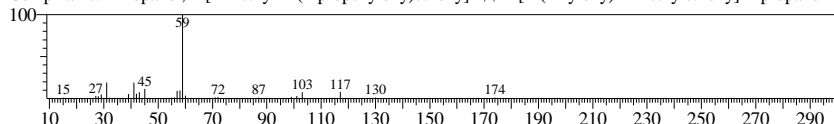
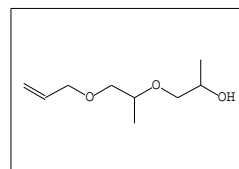
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#3 Entry:27550 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

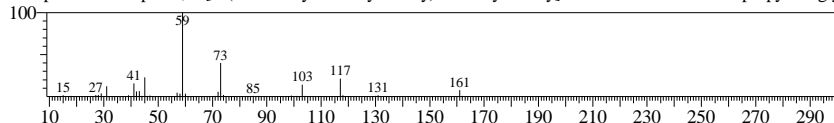
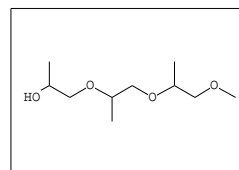
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:17307 Library:NIST08s.LIB

SI:78 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

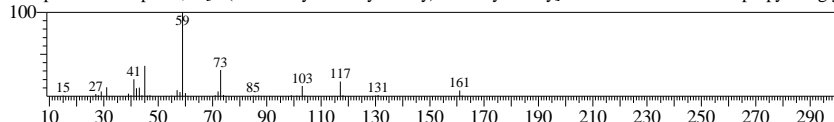
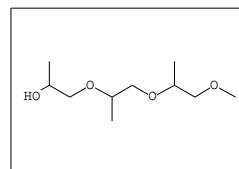
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#5 Entry:17306 Library:NIST08s.LIB

SI:78 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

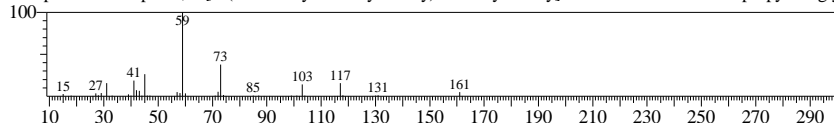
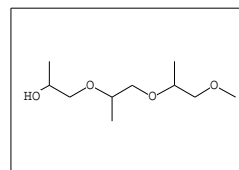
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#6 Entry:46317 Library:NIST08.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

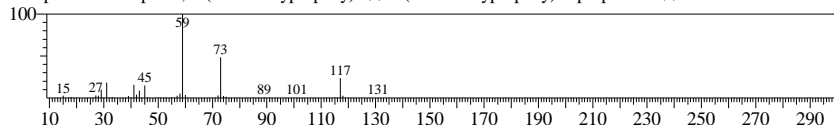
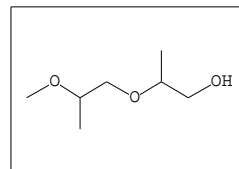
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#7 Entry:14297 Library:NIST08.LIB

SI:76 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

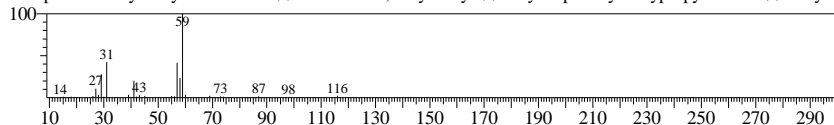
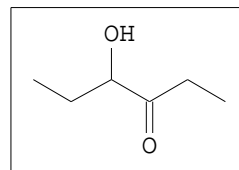
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#8 Entry:4568 Library:NIST08.LIB

SI:76 Formula:C6H12O2 CAS:4984-85-4 MolWeight:116 RetIndex:916

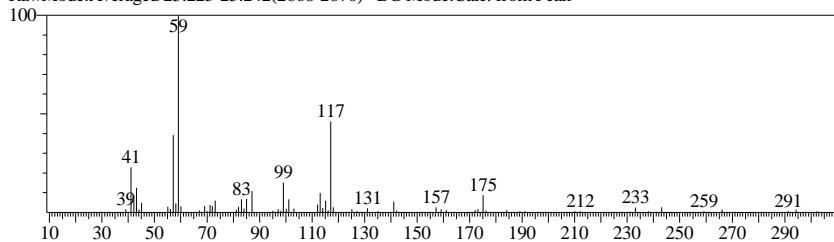
CompName:4-Hydroxy-3-hexanone \$\$ 3-Hexanone, 4-hydroxy- \$\$ Ethyl .alpha.-hydroxypropyl ketone \$\$ 3-Hydr



<< Target >>

Line#:51 R.Time:25.233(Scan#:2669) BasePeak:59.10(116970)

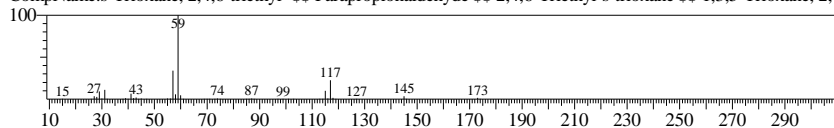
RawMode:Averaged 25.225-25.242(2668-2670) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:77 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

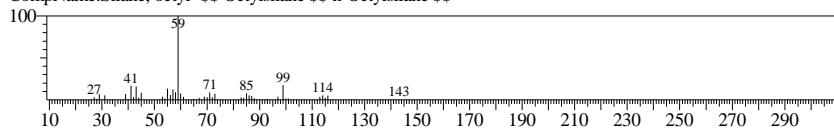
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:12984 Library:NIST08.LIB

SI:75 Formula:C8H20Si CAS:871-92-1 MolWeight:144 RetIndex:0

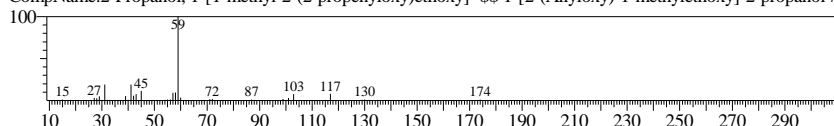
CompName:Silane, octyl- \$\$ Octylsilane \$\$ n-Octylsilane \$\$



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:75 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

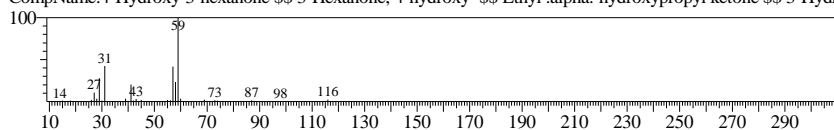
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:4568 Library:NIST08.LIB

SI:75 Formula:C6H12O2 CAS:4984-85-4 MolWeight:116 RetIndex:916

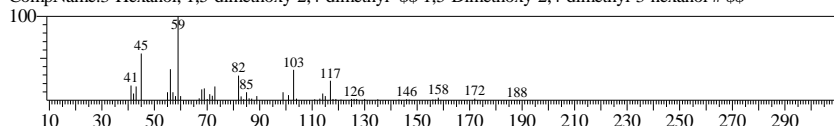
CompName:4-Hydroxy-3-hexanone \$\$ 3-Hexanone, 4-hydroxy- \$\$ Ethyl .alpha.-hydroxypropyl ketone \$\$ 3-Hydr



Hit#:5 Entry:36580 Library:NIST08.LIB

SI:74 Formula:C10H22O3 CAS:13897-22-8 MolWeight:190 RetIndex:1137

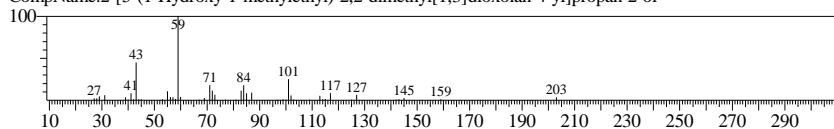
CompName:3-Hexanol, 1,5-dimethoxy-, 2,4-dimethyl- \$\$ 1,5-Dimethoxy-2,4-dimethyl-3-hexanol # \$\$



Hit#:6 Entry:54251 Library:NIST08.LIB

SI:74 Formula:C11H22O4 CAS:0-00-0 MolWeight:218 RetIndex:1415

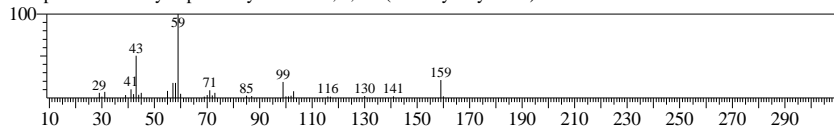
CompName:2-[5-(1-Hydroxy-1-methylethyl)-2,2-dimethyl[1,3]dioxolan-4-yl]propan-2-ol



Hit#:7 Entry:27412 Library:NIST08.LIB

SI:74 Formula:C8H14O4 CAS:0-00-0 MolWeight:174 RetIndex:1212

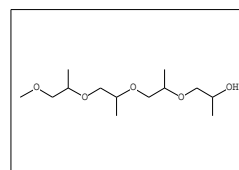
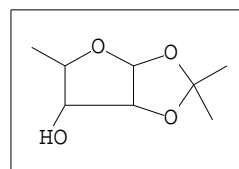
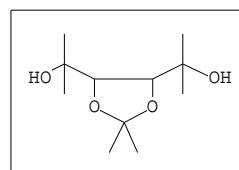
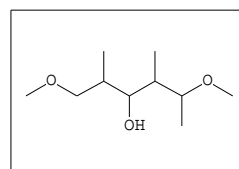
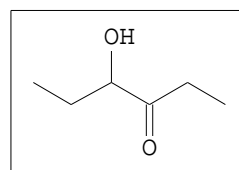
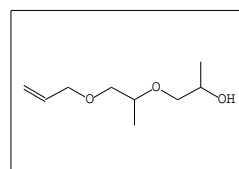
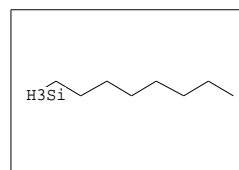
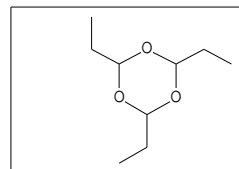
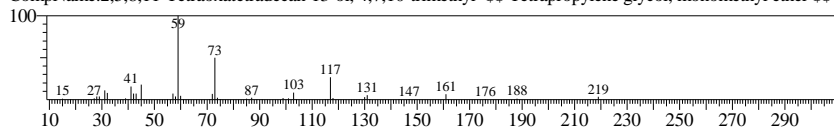
CompName:5-Deoxy-.alpha.-d-xylofuranose, 1,2-O-(1-methylethylidene)-



Hit#:8 Entry:85908 Library:NIST08.LIB

SI:74 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

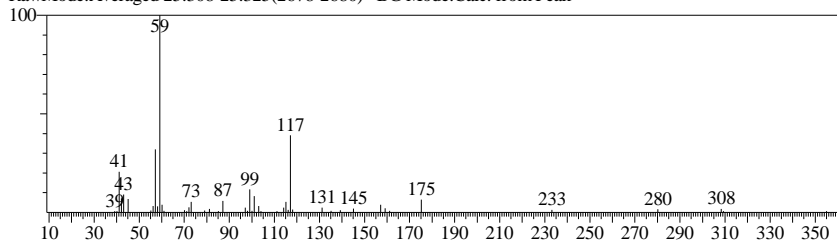
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



<< Target >>

Line#:52 R.Time:25.317(Scan#:2679) BasePeak:59.10(51813)

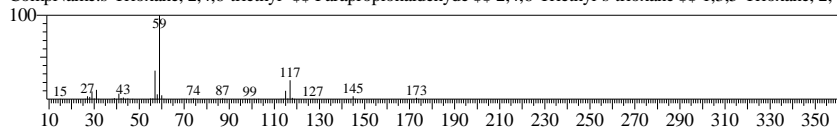
RawMode:Averaged 25.308-25.325(2678-2680) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:83 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

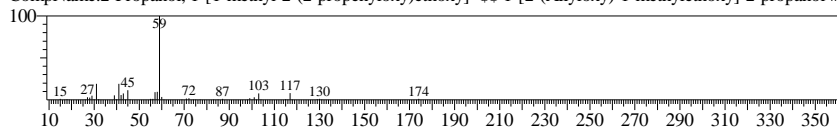
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:81 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

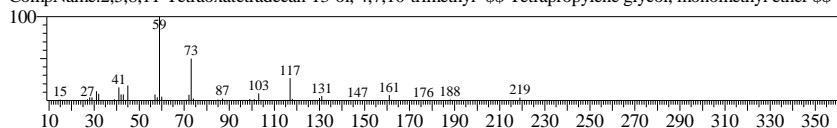
CompName:1-[1-methyl-2-(2-propenyloxy)ethoxy]-1-methylethoxy-2-propanol #



Hit#:3 Entry:85908 Library:NIST08.LIB

SI:80 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

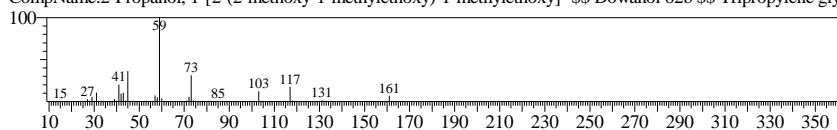
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#:4 Entry:17306 Library:NIST08s.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

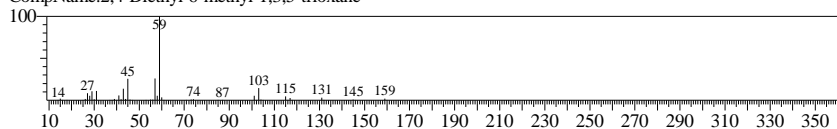
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:5 Entry:19968 Library:NIST08.LIB

SI:79 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

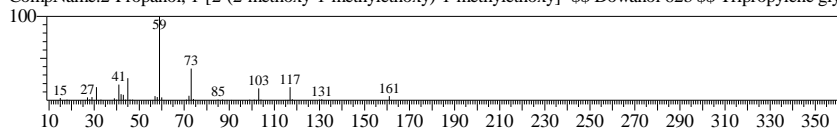
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:6 Entry:46317 Library:NIST08.LIB

SI:79 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

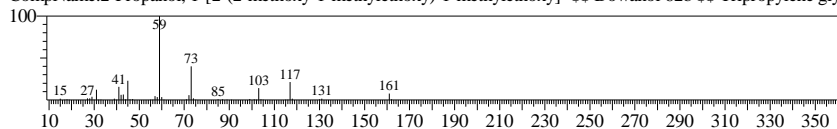
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:7 Entry:17307 Library:NIST08s.LIB

SI:79 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

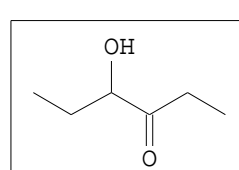
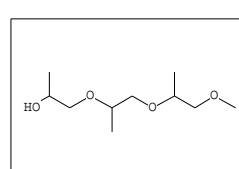
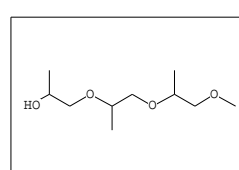
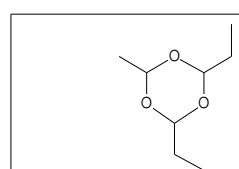
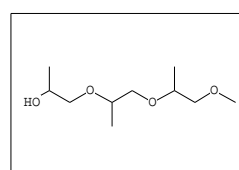
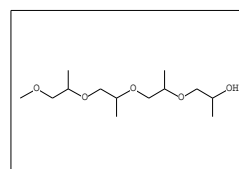
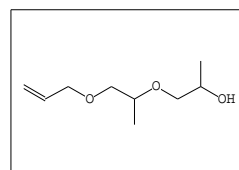
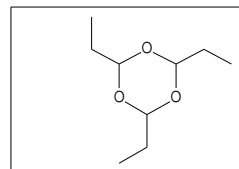
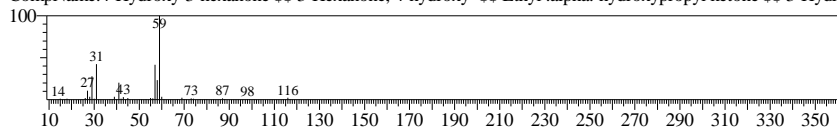
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:4568 Library:NIST08.LIB

SI:79 Formula:C6H12O2 CAS:4984-85-4 MolWeight:116 RetIndex:916

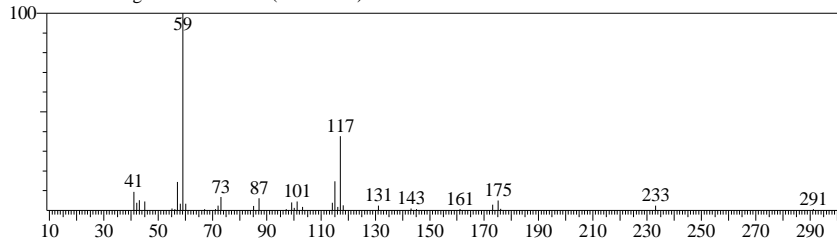
CompName:4-Hydroxy-3-hexanone \$\$ 3-Hexanone, 4-hydroxy- \$\$ Ethyl .alpha.-hydroxypropyl ketone \$\$ 3-Hydr



<< Target >>

Line# 53 R.Time:25.725(Scan#:2728) BasePeak:59.10(487907)

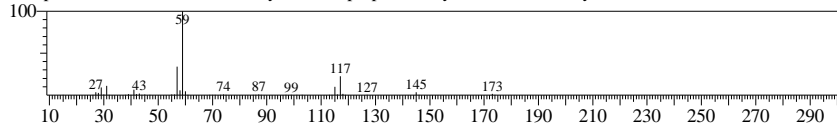
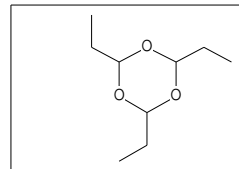
RawMode:Averaged 25.717-25.733(2727-2729) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:86 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

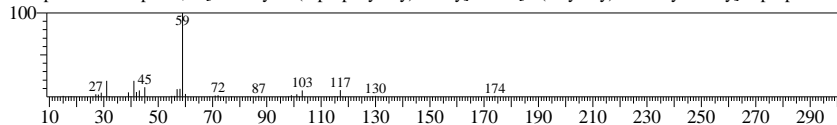
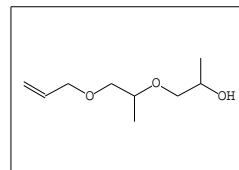
CompName:s-Trioxane, 2,4,6-triethyl- \$ \$ Parapropionaldehyde \$ \$ 2,4,6-Triethyl-s-trioxane \$ \$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

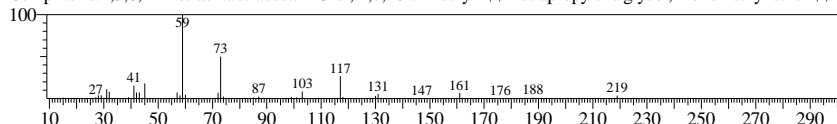
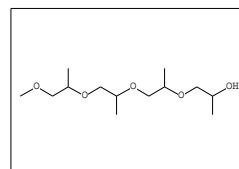
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:85908 Library:NIST08.LIB

SI:82 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

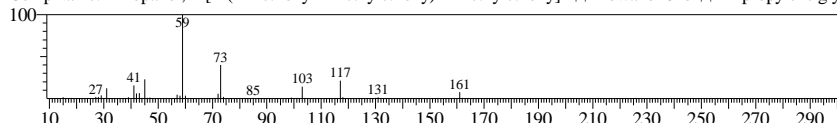
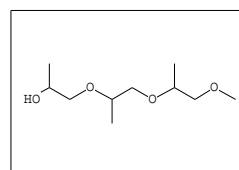
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$ \$ Tetrapropylene glycol, monomethyl ether \$ \$



Hit#:4 Entry:17307 Library:NIST08s.LIB

SI:81 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

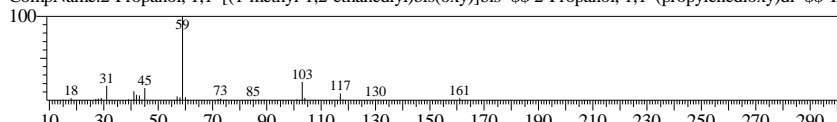
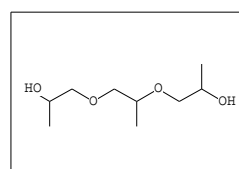
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ \$ Dowanol 62b \$ \$ Tripropylene gly



Hit#:5 Entry:37499 Library:NIST08.LIB

SI:81 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

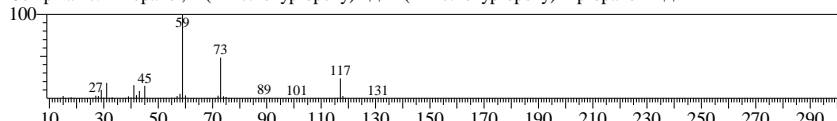
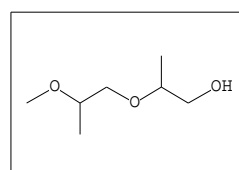
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$ \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ \$ T



Hit#:6 Entry:14297 Library:NIST08.LIB

SI:81 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

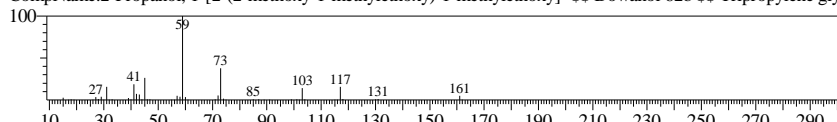
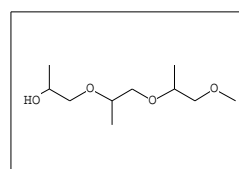
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$ \$ 2-(2-Methoxypropoxy)-1-propanol # \$ \$



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

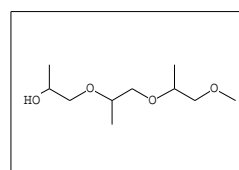
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ \$ Dowanol 62b \$ \$ Tripropylene gly



Hit#:8 Entry:17306 Library:NIST08s.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

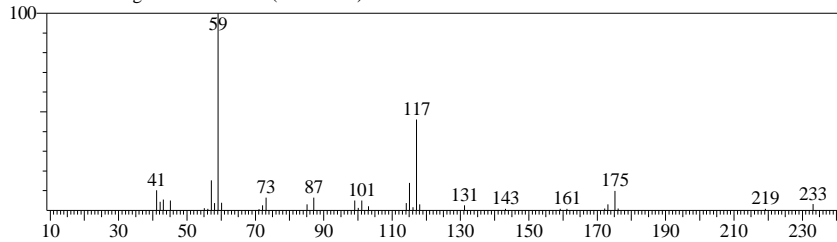
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ \$ Dowanol 62b \$ \$ Tripropylene gly



<< Target >>

Line# 54 R.Time:25.808(Scan#:2738) BasePeak:59.10(974667)

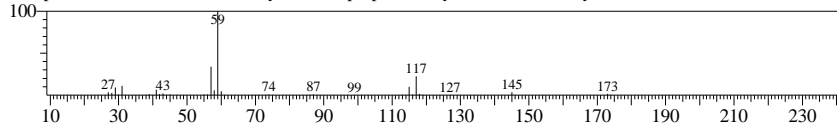
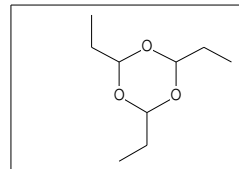
RawMode:Averaged 25.800-25.817(2737-2739) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:83 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

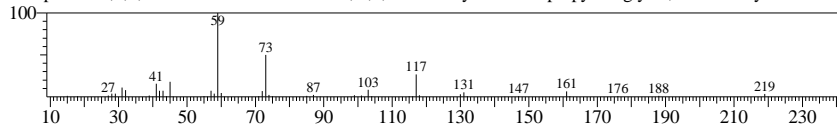
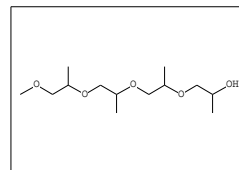
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:81 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

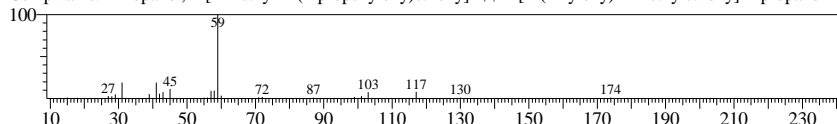
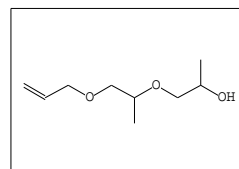
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#3 Entry:27550 Library:NIST08.LIB

SI:81 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

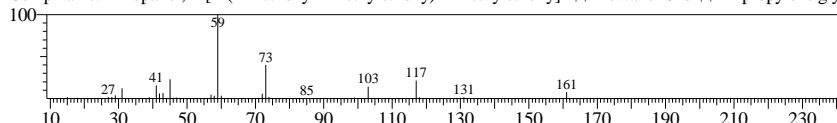
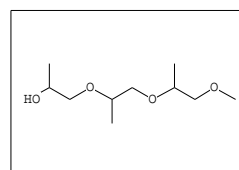
CompName:2-Propanol, 1-[2-(1-methyl-2-(2-propenyloxy)ethoxy)]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:17307 Library:NIST08s.LIB

SI:80 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

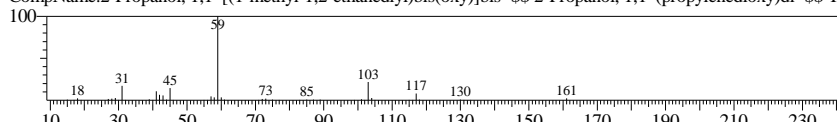
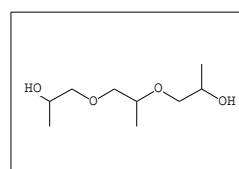
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#5 Entry:37499 Library:NIST08.LIB

SI:80 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

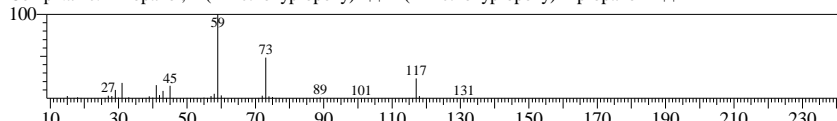
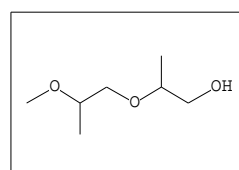
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#6 Entry:14297 Library:NIST08.LIB

SI:80 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

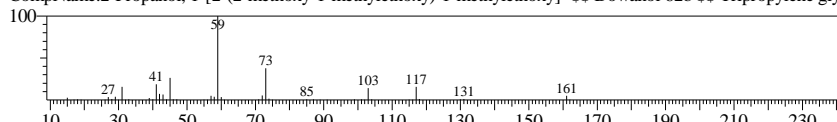
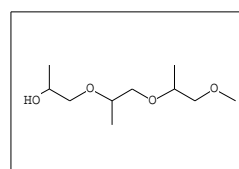
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#7 Entry:46317 Library:NIST08.LIB

SI:79 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

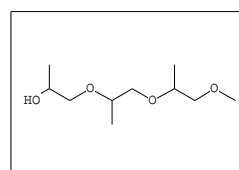
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#8 Entry:17306 Library:NIST08s.LIB

SI:79 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

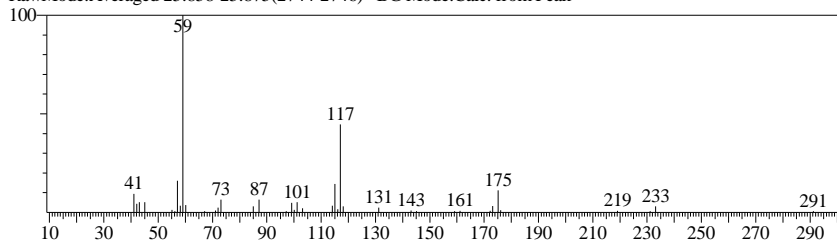
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line#55 R.Time:25.867(Scan#:2745) BasePeak:59.10(819386)

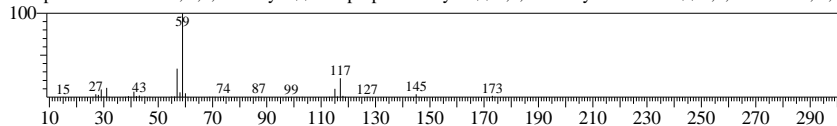
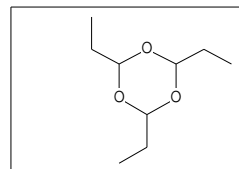
RawMode:Averaged 25.858-25.875(2744-2746) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:83 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

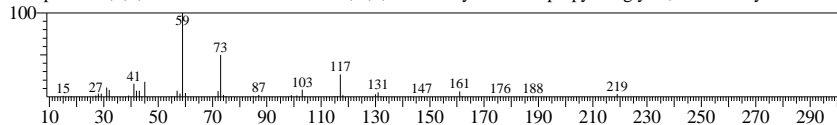
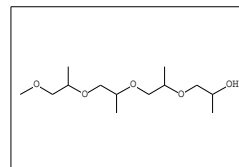
CompName:s-Trioxane, 2,4,6-triethyl- \$S\$ Parapropionaldehyde \$S\$ 2,4,6-Triethyl-s-trioxane \$S\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:80 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

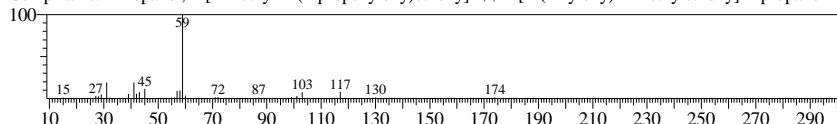
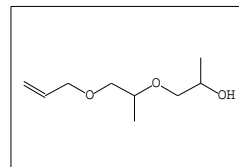
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$S\$ Tetrapropylene glycol, monomethyl ether \$S\$ 4



Hit#3 Entry:27550 Library:NIST08.LIB

SI:80 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

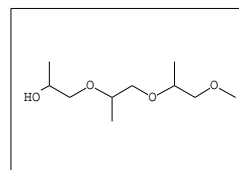
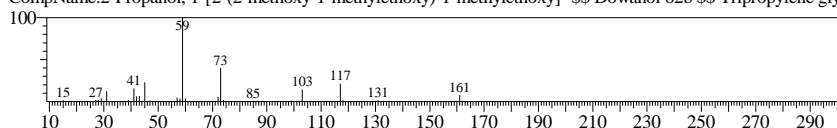
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$S\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:17307 Library:NIST08s.LIB

SI:79 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

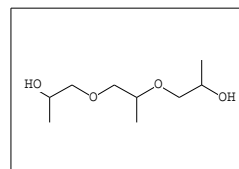
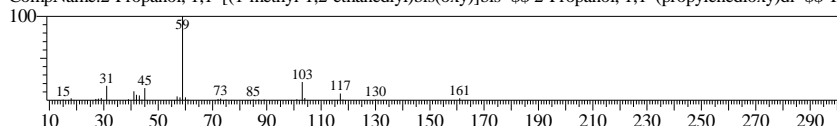
CompName:2-Propanol, 1,1'-[(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$S\$ Dowanol 62b \$S\$ Tripropylene gly



Hit#5 Entry:37499 Library:NIST08.LIB

SI:79 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

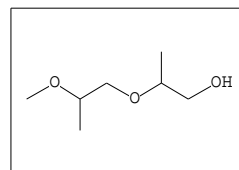
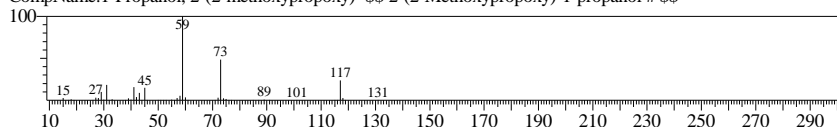
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$S\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$S\$ T



Hit#6 Entry:14297 Library:NIST08.LIB

SI:79 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

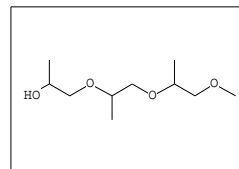
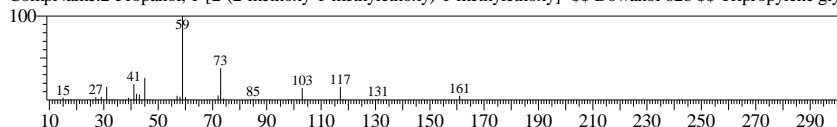
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$S\$ 2-(2-Methoxypropoxy)-1-propanol # \$S\$



Hit#7 Entry:46317 Library:NIST08.LIB

SI:78 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

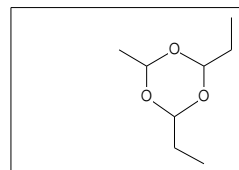
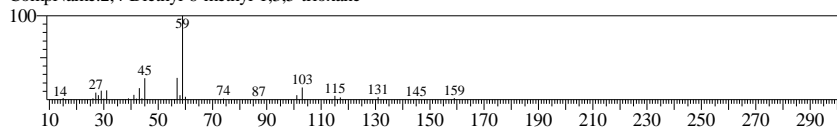
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$S\$ Dowanol 62b \$S\$ Tripropylene gly



Hit#8 Entry:19968 Library:NIST08.LIB

SI:78 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

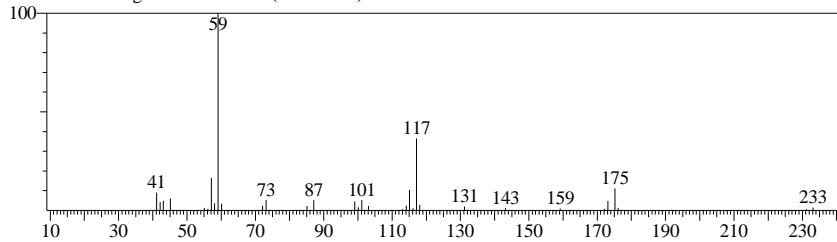
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line# 56 R.Time:25.950(Scan#:2755) BasePeak:59.10(295894)

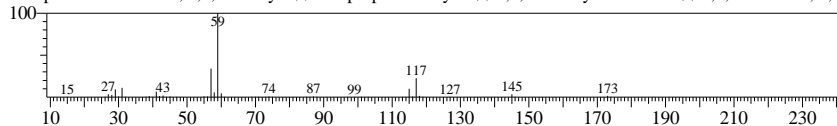
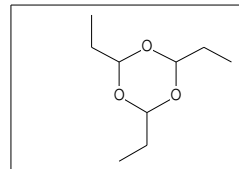
RawMode:Averaged 25.942-25.958(2754-2756) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:85 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

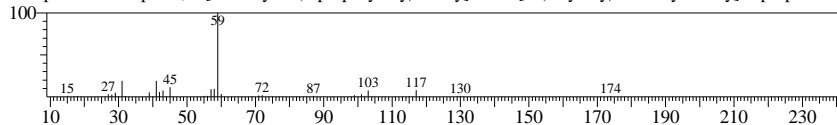
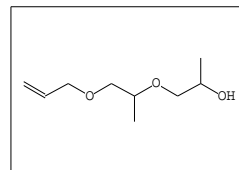
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:81 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

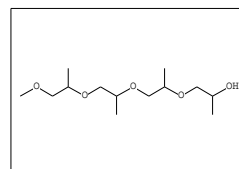
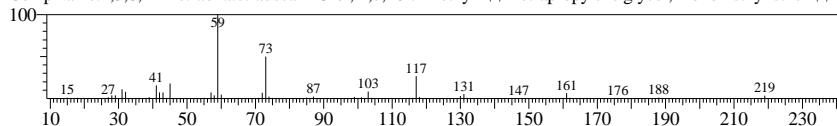
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:85908 Library:NIST08.LIB

SI:80 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

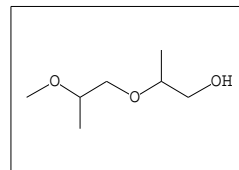
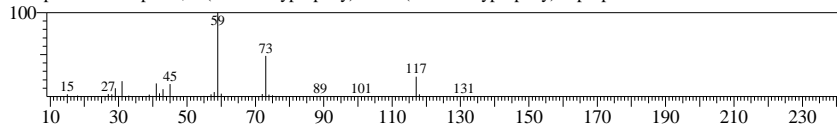
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 2



Hit#:4 Entry:14297 Library:NIST08.LIB

SI:80 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

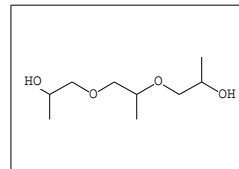
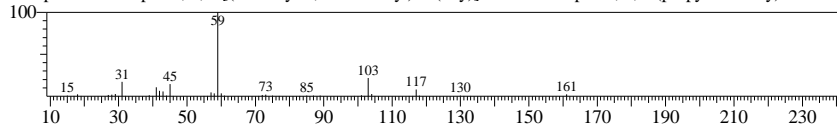
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:5 Entry:37499 Library:NIST08.LIB

SI:80 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

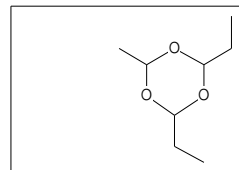
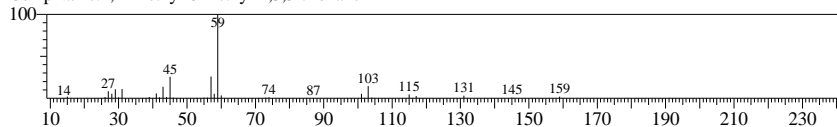
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediy)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:6 Entry:19968 Library:NIST08.LIB

SI:80 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

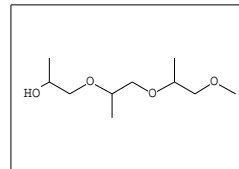
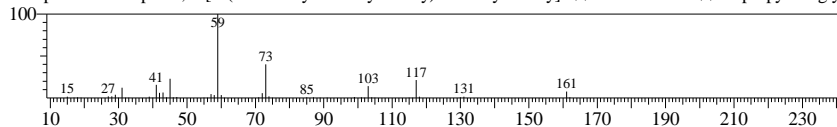
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:7 Entry:17307 Library:NIST08s.LIB

SI:79 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

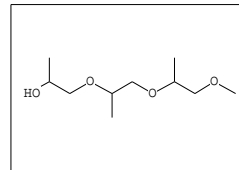
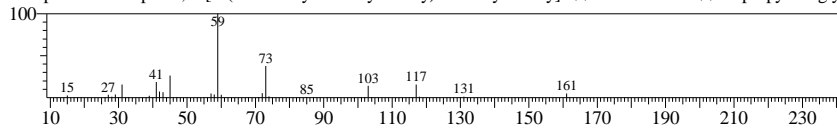
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:46317 Library:NIST08.LIB

SI:78 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

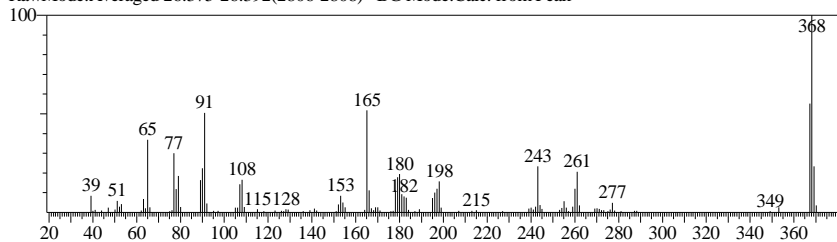
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line#:57 R.Time:26.383(Scan#:2807) BasePeak:368.15(250569)

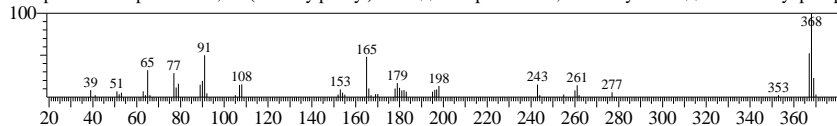
RawMode:Averaged 26.375-26.392(2806-2808) BG Mode:Calc. from Peak



Hit#:1 Entry:154440 Library:NIST08s.LIB

SI:96 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

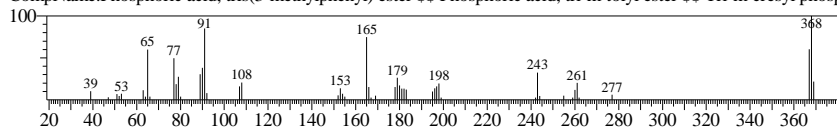
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:2 Entry:26661 Library:NIST08s.LIB

SI:91 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

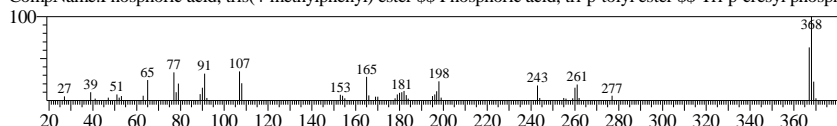
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:3 Entry:154441 Library:NIST08s.LIB

SI:91 Formula:C₂₁H₂₁O₄P CAS:78-32-0 MolWeight:368 RetIndex:0

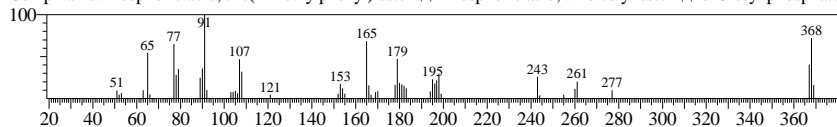
CompName:Phosphoric acid, tris(4-methylphenyl) ester \$\$ Phosphoric acid, tri-p-tolyl ester \$\$ Tri-p-cresyl phosph



Hit#:4 Entry:26660 Library:NIST08s.LIB

SI:83 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

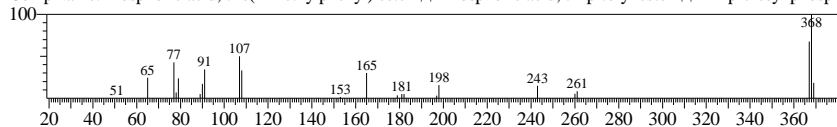
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



Hit#:5 Entry:26663 Library:NIST08s.LIB

SI:80 Formula:C₂₁H₂₁O₄P CAS:78-32-0 MolWeight:368 RetIndex:0

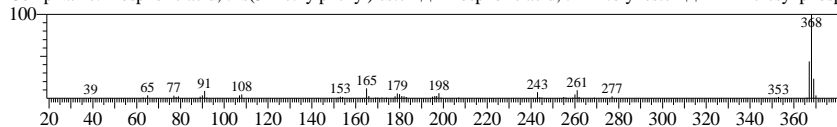
CompName:Phosphoric acid, tris(4-methylphenyl) ester \$\$ Phosphoric acid, tri-p-tolyl ester \$\$ Tri-p-cresyl phosph



Hit#:6 Entry:26664 Library:NIST08s.LIB

SI:77 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

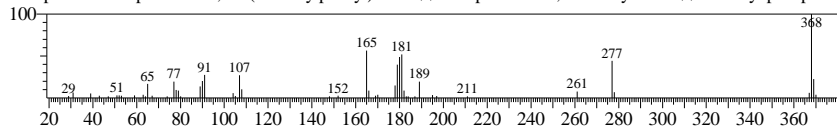
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:7 Entry:26662 Library:NIST08s.LIB

SI:76 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

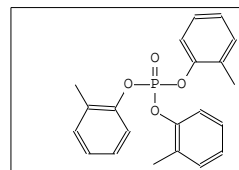
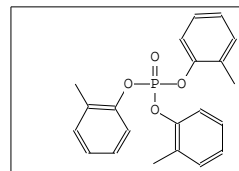
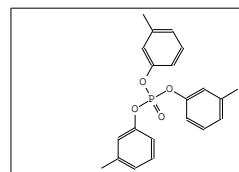
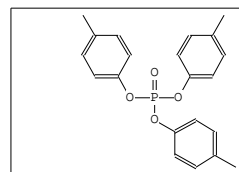
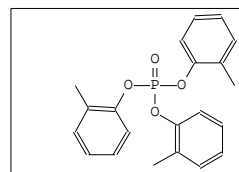
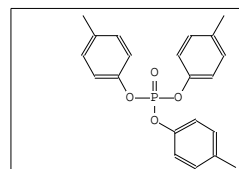
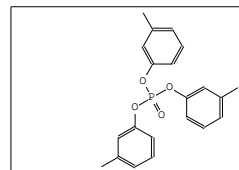
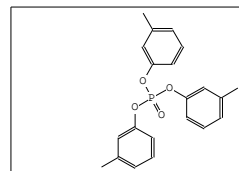
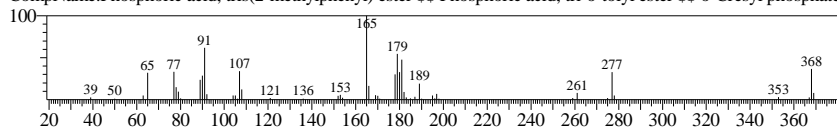
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



Hit#:8 Entry:154438 Library:NIST08s.LIB

SI:76 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

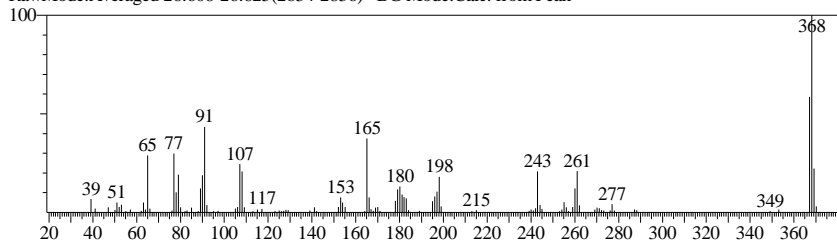
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



<< Target >>

Line#:58 R.Time:26.617(Scan#:2835) BasePeak:368.15(467668)

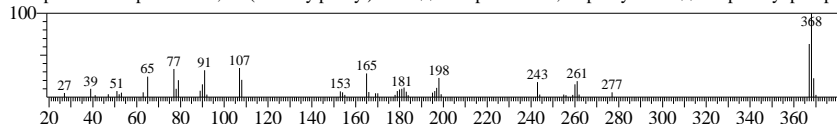
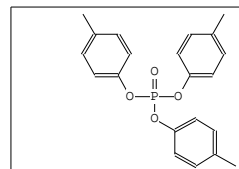
RawMode:Averaged 26.608-26.625(2834-2836) BG Mode:Calc. from Peak



Hit#:1 Entry:154441 Library:NIST08.LIB

SI:94 Formula:C₂₁H₂₁O₄P CAS:78-32-0 MolWeight:368 RetIndex:0

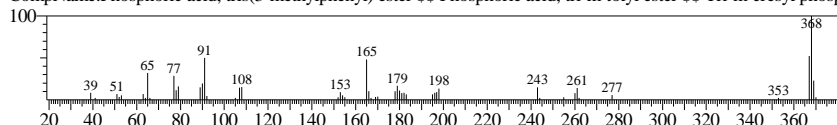
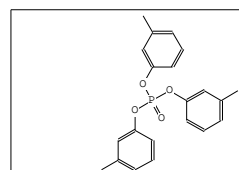
CompName:Phosphoric acid, tris(4-methylphenyl) ester \$\$ Phosphoric acid, tri-p-tolyl ester \$\$ Tri-p-cresyl phosph



Hit#:2 Entry:154440 Library:NIST08.LIB

SI:93 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

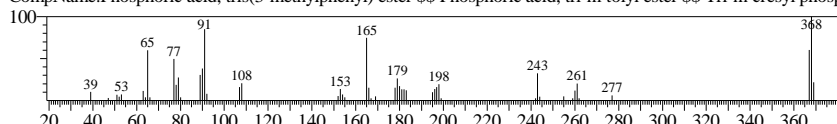
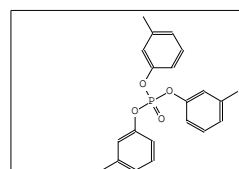
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:3 Entry:26661 Library:NIST08s.LIB

SI:89 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

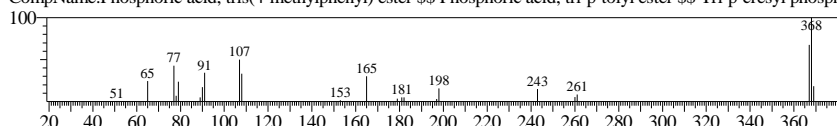
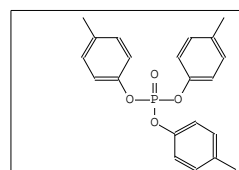
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:4 Entry:26663 Library:NIST08s.LIB

SI:84 Formula:C₂₁H₂₁O₄P CAS:78-32-0 MolWeight:368 RetIndex:0

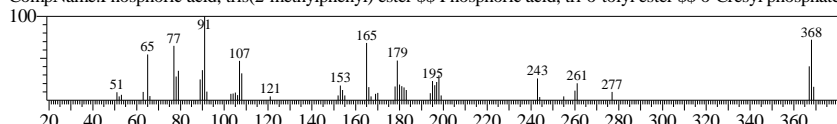
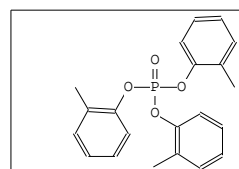
CompName:Phosphoric acid, tris(4-methylphenyl) ester \$\$ Phosphoric acid, tri-p-tolyl ester \$\$ Tri-p-cresyl phosph



Hit#:5 Entry:26660 Library:NIST08s.LIB

SI:81 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

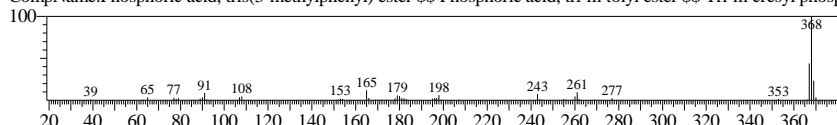
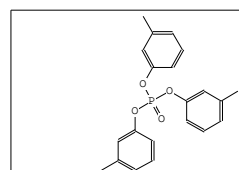
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



Hit#:6 Entry:26664 Library:NIST08s.LIB

SI:78 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

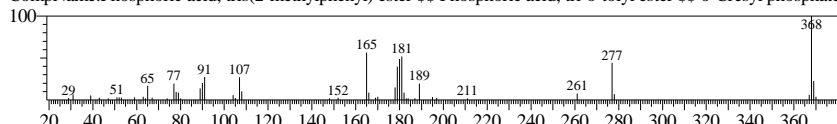
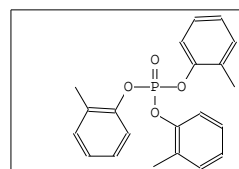
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:7 Entry:26662 Library:NIST08s.LIB

SI:75 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

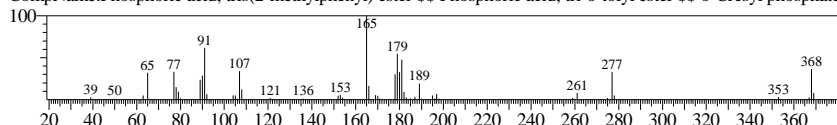
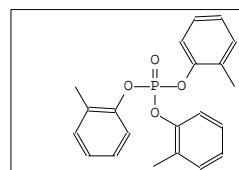
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



Hit#:8 Entry:154438 Library:NIST08.LIB

SI:73 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

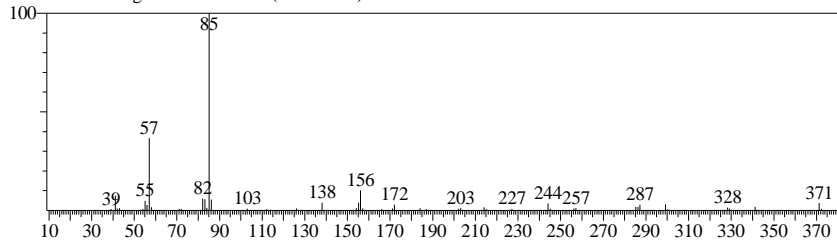
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



<< Target >>

Line#:59 R.Time:26.750(Scan#:2851) BasePeak:85.10(1326984)

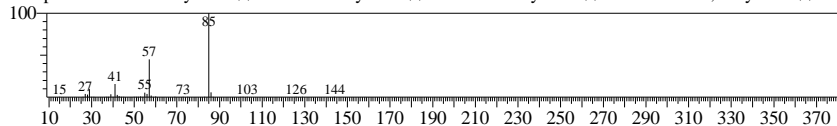
RawMode:Averaged 26.742-26.758(2850-2852) BG Mode:Calc. from Peak



Hit#:1 Entry:34397 Library:NIST08s.LIB

SI:80 Formula:C10H18O3 CAS:2082-59-9 MolWeight:186 RetIndex:1319

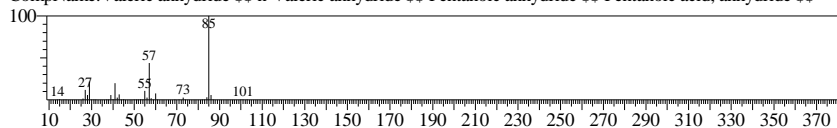
CompName:Valeric anhydride \$\$ n-Valeric anhydride \$\$ Pentanoic anhydride \$\$ Pentanoic acid, anhydride \$\$



Hit#:2 Entry:14624 Library:NIST08s.LIB

SI:78 Formula:C10H18O3 CAS:2082-59-9 MolWeight:186 RetIndex:1319

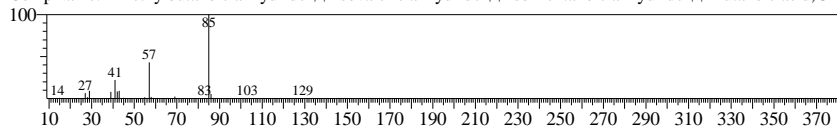
CompName:Valeric anhydride \$\$ n-Valeric anhydride \$\$ Pentanoic anhydride \$\$ Pentanoic acid, anhydride \$\$



Hit#:3 Entry:34398 Library:NIST08s.LIB

SI:77 Formula:C10H18O3 CAS:1468-39-9 MolWeight:186 RetIndex:1190

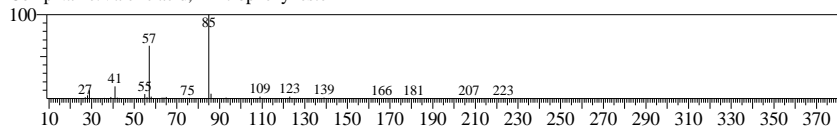
CompName:2-Methylbutanoic anhydride \$\$ Isovaleric anhydride \$\$ iso-Pentanoic anhydride \$\$ Butanoic acid, 3-



Hit#:4 Entry:57757 Library:NIST08s.LIB

SI:76 Formula:C11H13NO4 CAS:0-00-0 MolWeight:223 RetIndex:1754

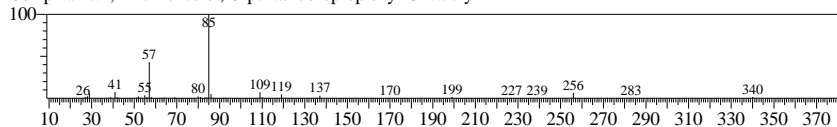
CompName:Valeric acid, 4-nitrophenyl ester



Hit#:5 Entry:138645 Library:NIST08s.LIB

SI:76 Formula:C14H13F5O4 CAS:0-00-0 MolWeight:340 RetIndex:1533

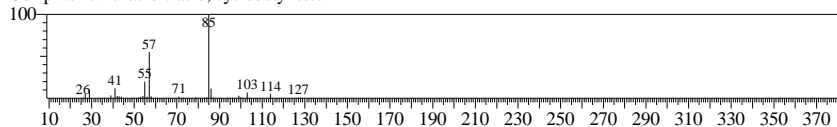
CompName:1,2-Benzenediol, O-pentafluoropropionyl-O'-valeryl-



Hit#:6 Entry:18296 Library:NIST08s.LIB

SI:75 Formula:C9H16O2 CAS:0-00-0 MolWeight:156 RetIndex:1106

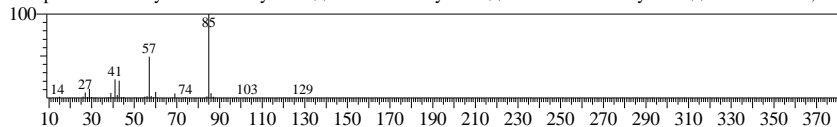
CompName:Valeric acid, cyclobutyl ester



Hit#:7 Entry:14625 Library:NIST08s.LIB

SI:75 Formula:C10H18O3 CAS:1468-39-9 MolWeight:186 RetIndex:1190

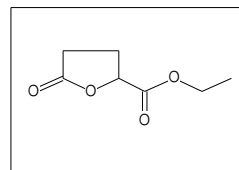
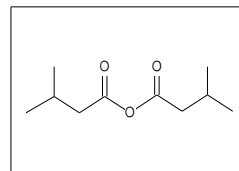
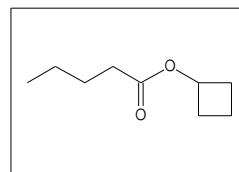
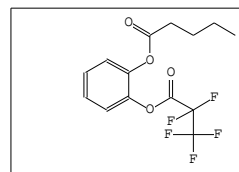
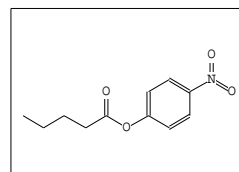
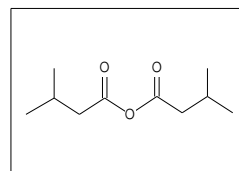
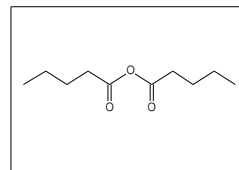
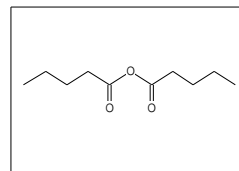
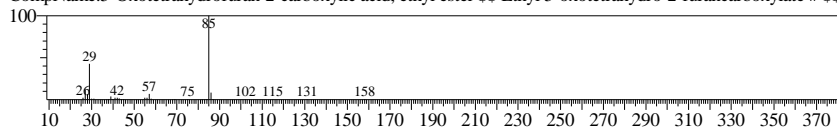
CompName:2-Methylbutanoic anhydride \$\$ Isovaleric anhydride \$\$ iso-Pentanoic anhydride \$\$ Butanoic acid, 3-



Hit#:8 Entry:18953 Library:NIST08s.LIB

SI:74 Formula:C7H10O4 CAS:1126-51-8 MolWeight:158 RetIndex:1252

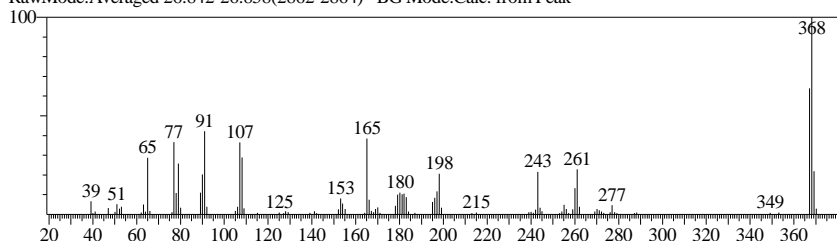
CompName:5-Oxotetrahydrofuran-2-carboxylic acid, ethyl ester \$\$ Ethyl 5-oxotetrahydro-2-furancarboxylate # \$\$



<< Target >>

Line#:60 R.Time:26.850(Scan#:2863) BasePeak:368.20(250283)

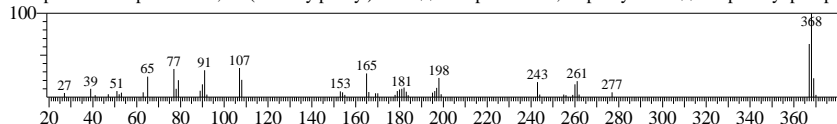
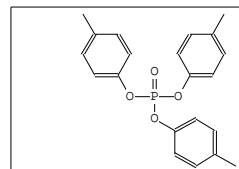
RawMode:Averaged 26.842-26.858(2862-2864) BG Mode:Calc. from Peak



Hit#:1 Entry:154441 Library:NIST08.LIB

SI:95 Formula:C₂₁H₂₁O₄P CAS:78-32-0 MolWeight:368 RetIndex:0

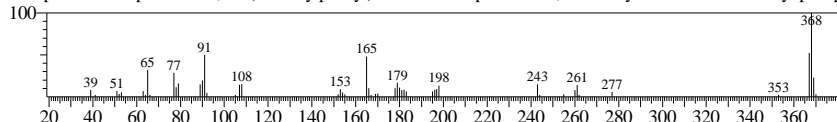
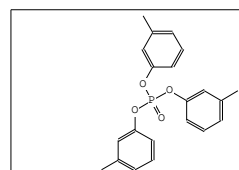
CompName:Phosphoric acid, tris(4-methylphenyl) ester \$\$ Phosphoric acid, tri-p-tolyl ester \$\$ Tri-p-cresyl phosph



Hit#:2 Entry:154440 Library:NIST08.LIB

SI:91 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

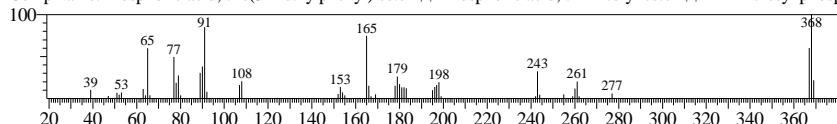
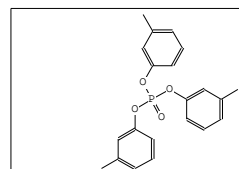
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:3 Entry:26661 Library:NIST08s.LIB

SI:88 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

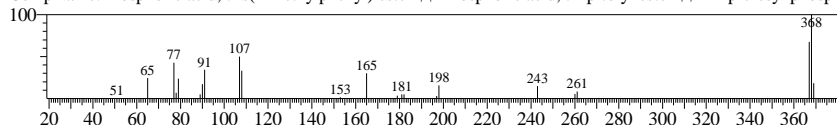
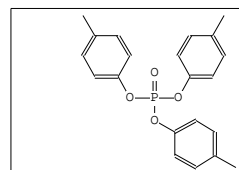
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:4 Entry:26663 Library:NIST08s.LIB

SI:85 Formula:C₂₁H₂₁O₄P CAS:78-32-0 MolWeight:368 RetIndex:0

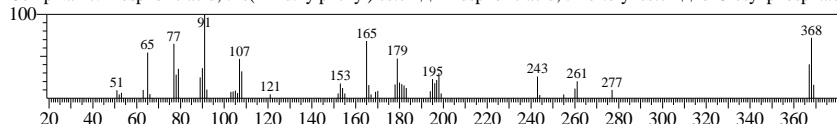
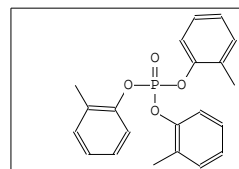
CompName:Phosphoric acid, tris(4-methylphenyl) ester \$\$ Phosphoric acid, tri-p-tolyl ester \$\$ Tri-p-cresyl phosph



Hit#:5 Entry:26660 Library:NIST08s.LIB

SI:83 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

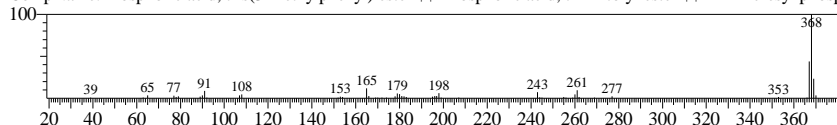
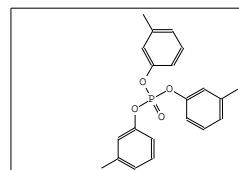
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



Hit#:6 Entry:26664 Library:NIST08s.LIB

SI:76 Formula:C₂₁H₂₁O₄P CAS:563-04-2 MolWeight:368 RetIndex:0

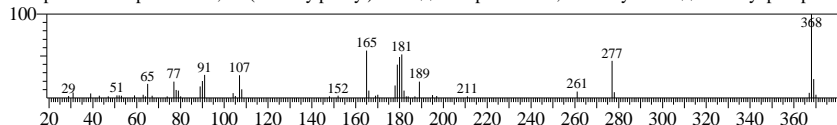
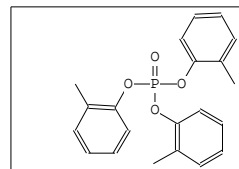
CompName:Phosphoric acid, tris(3-methylphenyl) ester \$\$ Phosphoric acid, tri-m-tolyl ester \$\$ Tri-m-cresyl phosph



Hit#:7 Entry:26662 Library:NIST08s.LIB

SI:74 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

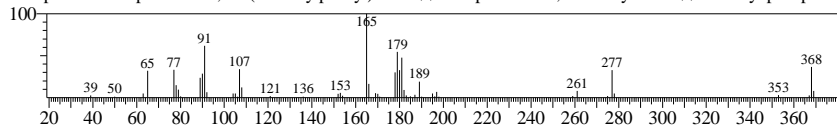
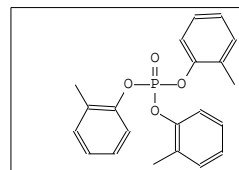
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



Hit#:8 Entry:154438 Library:NIST08.LIB

SI:73 Formula:C₂₁H₂₁O₄P CAS:78-30-8 MolWeight:368 RetIndex:0

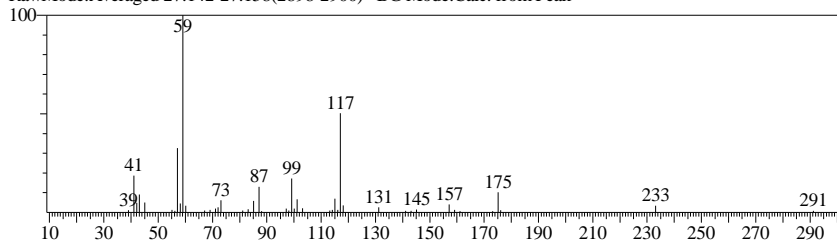
CompName:Phosphoric acid, tris(2-methylphenyl) ester \$\$ Phosphoric acid, tri-o-tolyl ester \$\$ o-Cresyl phosphate



<< Target >>

Line# 61 R.Time:27.150(Scan#:2899) BasePeak:59.10(284679)

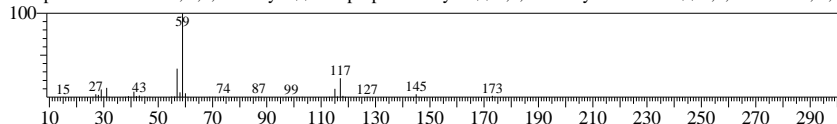
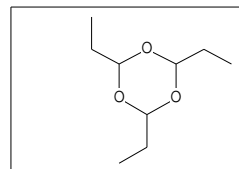
RawMode:Averaged 27.142-27.158(2898-2900) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

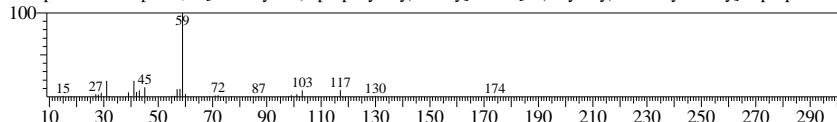
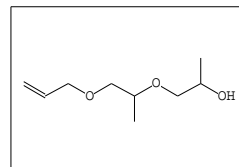
CompName:s-Trioxane, 2,4,6-triethyl- \$- Parapropionaldehyde \$- 2,4,6-Triethyl-s-trioxane \$- 1,3,5-Trioxane, 2,4



Hit#2 Entry:27550 Library:NIST08.LIB

SI:79 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

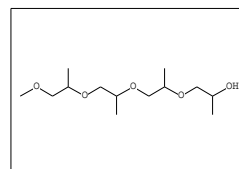
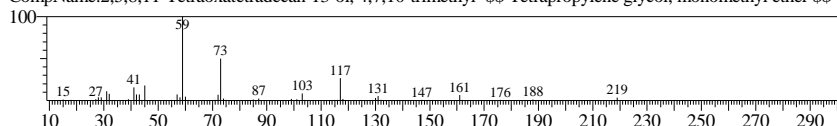
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#3 Entry:85908 Library:NIST08.LIB

SI:78 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

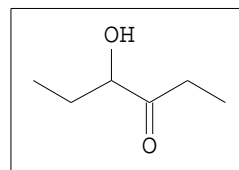
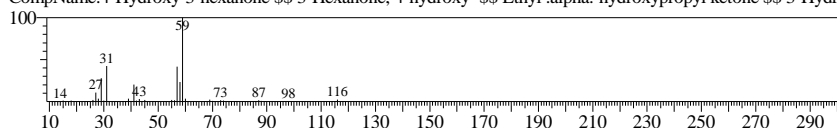
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$- Tetrapropylene glycol, monomethyl ether \$- 4



Hit#4 Entry:4568 Library:NIST08.LIB

SI:77 Formula:C6H12O2 CAS:4984-85-4 MolWeight:116 RetIndex:916

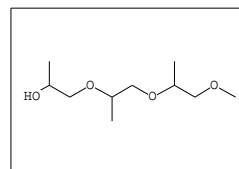
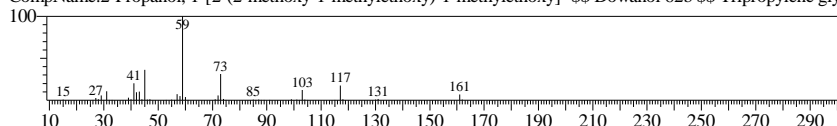
CompName:4-Hydroxy-3-hexanone \$- 3-Hexanone, 4-hydroxy- \$- Ethyl .alpha.-hydroxypropyl ketone \$- 3-Hydr



Hit#5 Entry:17306 Library:NIST08s.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

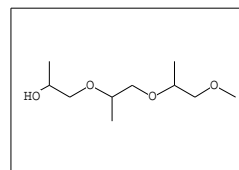
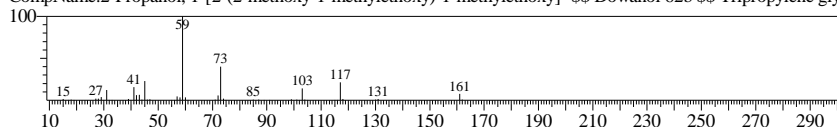
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#6 Entry:17307 Library:NIST08s.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

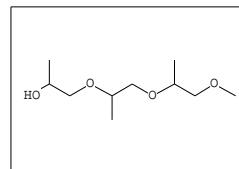
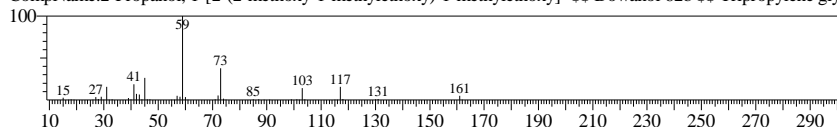
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#7 Entry:46317 Library:NIST08.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

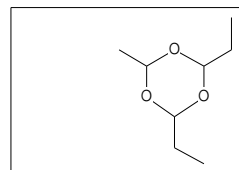
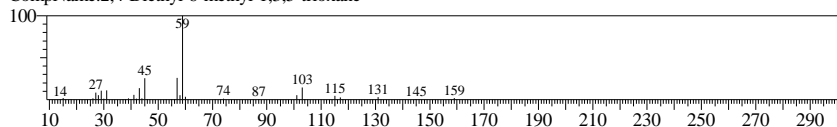
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#8 Entry:19968 Library:NIST08.LIB

SI:76 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

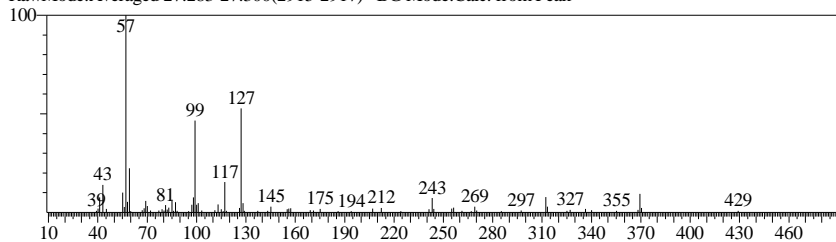
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line#:62 R.Time:27.292(Scan#:2916) BasePeak:57.10(51646)

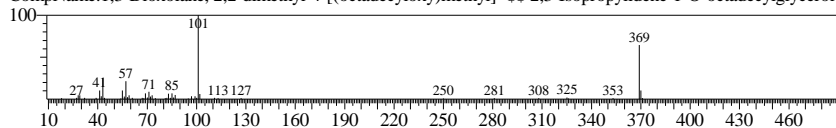
RawMode:Averaged 27.283-27.300(2915-2917) BG Mode:Calc. from Peak



Hit#:1 Entry:161964 Library:NIST08.LIB

SI:57 Formula:C24H48O3 CAS:16725-43-2 MolWeight:384 RetIndex:2638

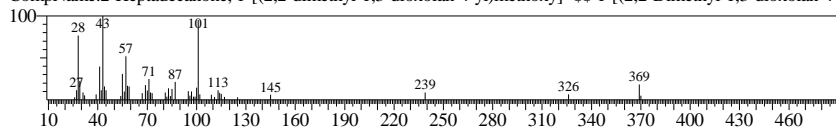
CompName:1,3-Dioxolane, 2,2-dimethyl-4-[(octadecyloxy)methyl]- \$\$ 2,3-Isopropylidene-1-O-octadecylglycerol



Hit#:2 Entry:161852 Library:NIST08.LIB

SI:56 Formula:C23H44O4 CAS:39033-40-4 MolWeight:384 RetIndex:2675

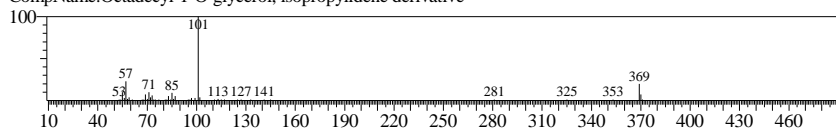
CompName:2-Heptadecanone, 1-[(2,2-dimethyl-1,3-dioxolan-4-yl)methoxy]- \$ 1-[(2,2-Dimethyl-1,3-dioxolan-4-



Hit#:3 Entry:161963 Library:NIST08.LIB

SI:55 Formula:C24H48O3 CAS:0-00-0 MolWeight:384 RetIndex:2638

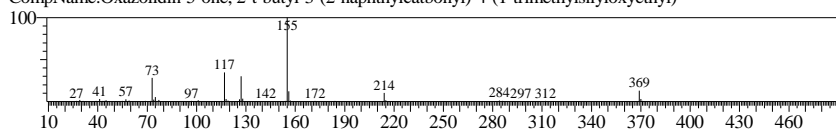
CompName:Octadecyl-1-O-glycerol, isopropylidene derivative



Hit#:4 Entry:171390 Library:NIST08.LIB

SI:48 Formula:C23H31NO4Si CAS:0-00-0 MolWeight:413 RetIndex:2953

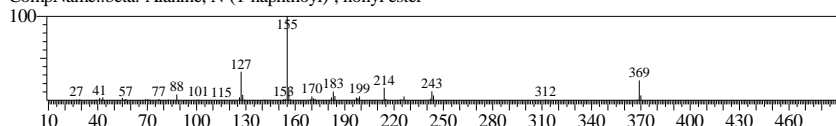
CompName:Oxazolidin-5-one, 2-t-butyl-3-(2-naphthylcarbonyl)-4-(1-trimethylsilyloxyethyl)-



Hit#:5 Entry:155138 Library:NIST08.LIB

SI:45 Formula:C23H31NO3 CAS:0-00-0 MolWeight:369 RetIndex:3105

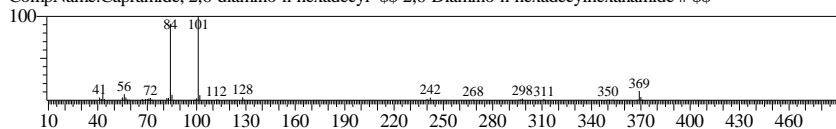
CompName:beta-Alanine, N-(1-naphthoyl)-, nonyl ester



Hit#:6 Entry:155123 Library:NIST08.LIB

SI:42 Formula:C22H47N3O CAS:133831-22-8 MolWeight:369 RetIndex:3030

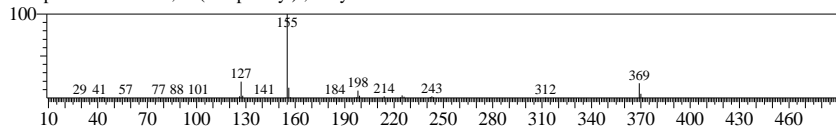
CompName:Capramide, 2,6-diamino-n-hexadecyl- \$\$ 2,6-Diamino-n-hexadecylhexanamide # \$\$



Hit#:7 Entry:155137 Library:NIST08.LIB

SI:41 Formula:C23H31NO3 CAS:0-00-0 MolWeight:369 RetIndex:2906

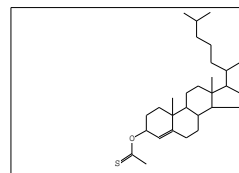
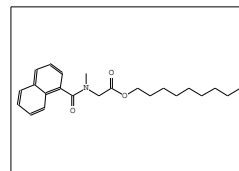
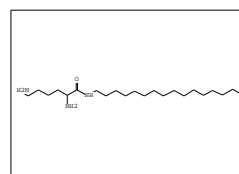
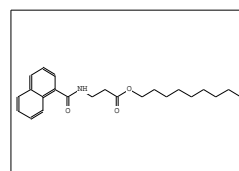
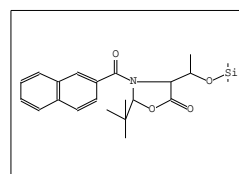
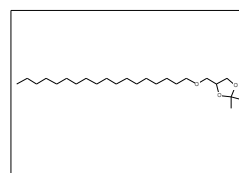
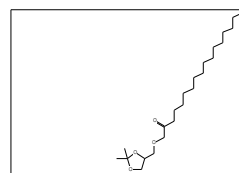
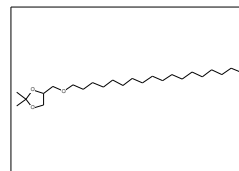
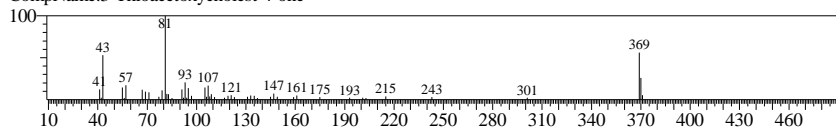
CompName:Sarcosine, N-(1-naphthoyl)-, nonyl ester



Hit#:8 Entry:178091 Library:NIST08.LIB

SI:40 Formula:C29H48OS CAS:0-00-0 MolWeight:444 RetIndex:3031

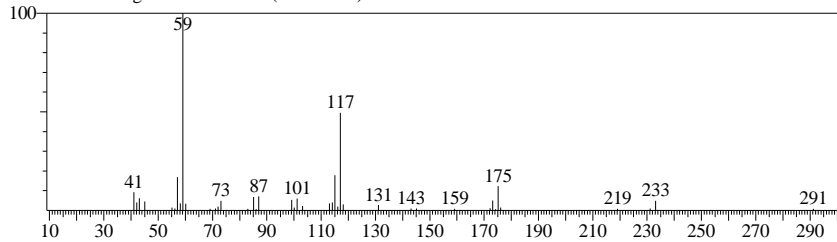
CompName:3-Thioacetoxcholest-4-one



<< Target >>

Line# 63 R.Time:27.650(Scan#:2959) BasePeak:59.10(653205)

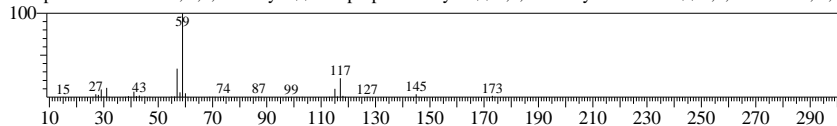
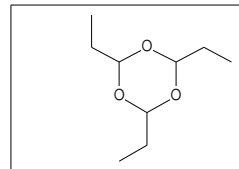
RawMode:Averaged 27.642-27.658(2958-2960) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

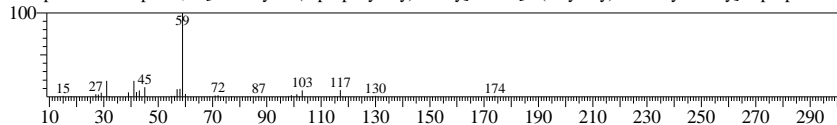
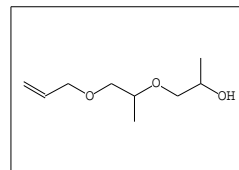
CompName:s-Trioxane, 2,4,6-triethyl- \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:27550 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

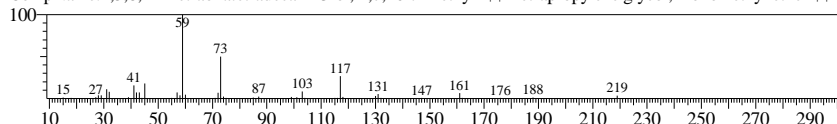
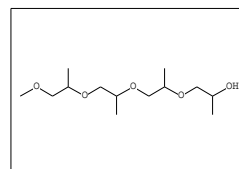
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:3 Entry:85908 Library:NIST08.LIB

SI:77 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

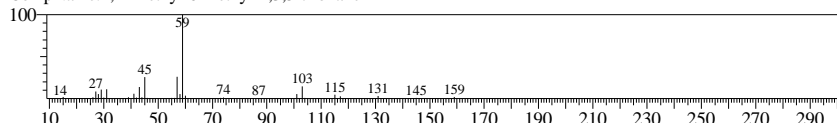
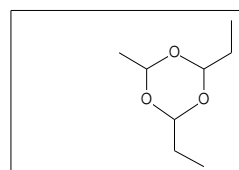
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$ Tetrapropylene glycol, monomethyl ether \$



Hit#:4 Entry:19968 Library:NIST08.LIB

SI:76 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

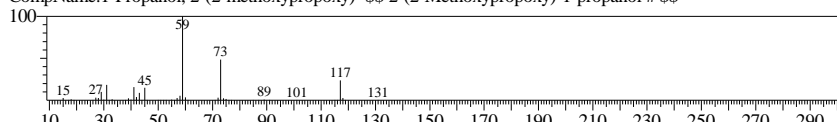
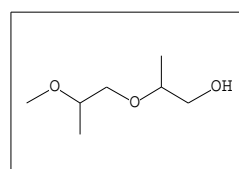
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:5 Entry:14297 Library:NIST08.LIB

SI:76 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

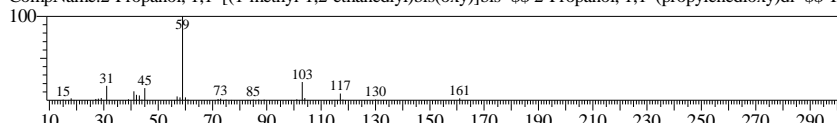
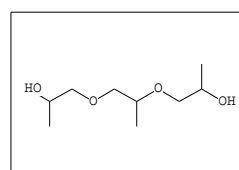
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$ 2-(2-Methoxypropoxy)-1-propanol # \$



Hit#:6 Entry:37499 Library:NIST08.LIB

SI:76 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

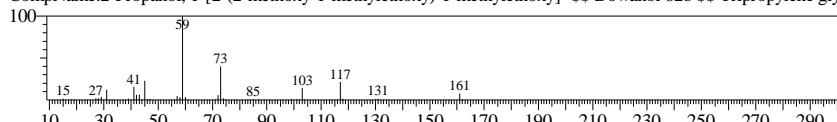
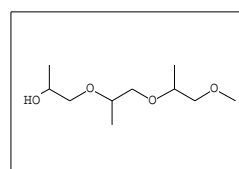
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$ 2-Propanol, 1,1'-(propylenedioxy)di- \$ T



Hit#:7 Entry:17307 Library:NIST08s.LIB

SI:75 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

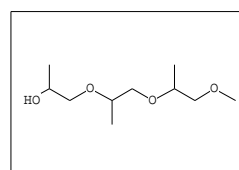
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



Hit#:8 Entry:46317 Library:NIST08.LIB

SI:75 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

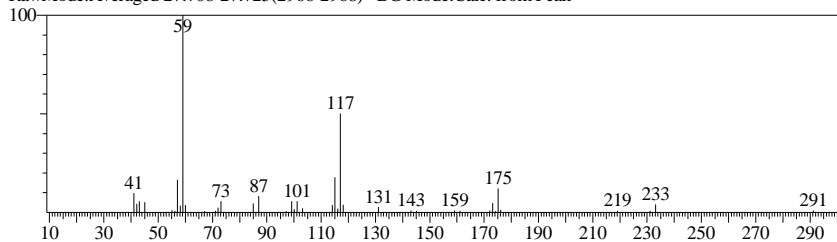
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$ Dowanol 62b \$ Tripropylene gly



<< Target >>

Line# 64 R.Time:27.717(Scan#:2967) BasePeak:59.10(1233123)

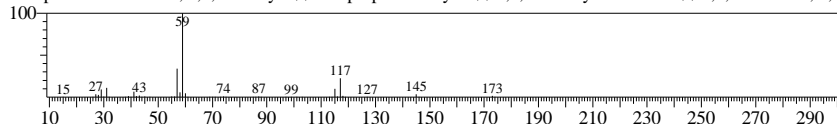
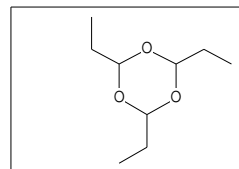
RawMode:Averaged 27.708-27.725(2966-2968) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:82 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

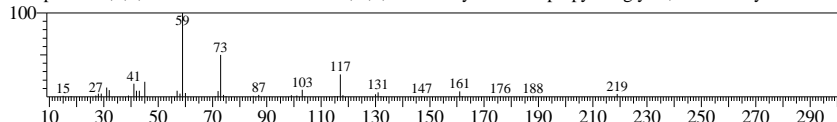
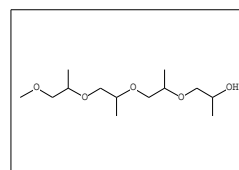
CompName:s-Trioxane, 2,4,6-triethyl- \$S\$ Parapropionaldehyde \$S\$ 2,4,6-Triethyl-s-trioxane \$S\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:79 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

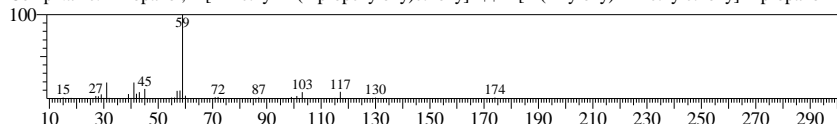
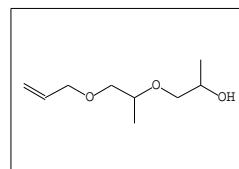
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$S\$ Tetrapropylene glycol, monomethyl ether \$S\$ 4



Hit#3 Entry:27550 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

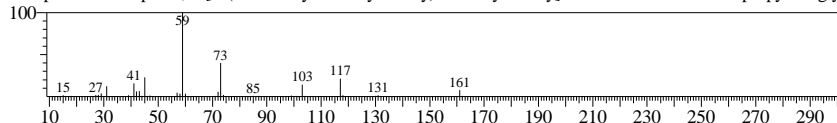
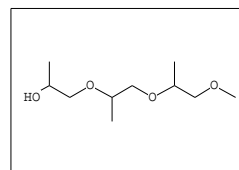
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$S\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:17307 Library:NIST08s.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

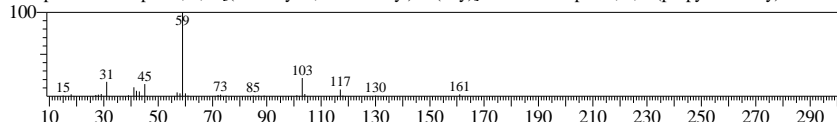
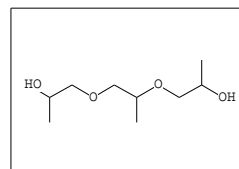
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$S\$ Dowanol 62b \$S\$ Tripropylene gly



Hit#5 Entry:37499 Library:NIST08.LIB

SI:77 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

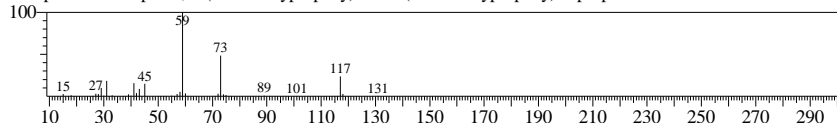
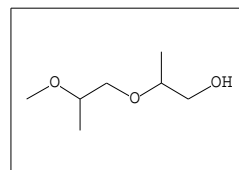
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$S\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$S\$ T



Hit#6 Entry:14297 Library:NIST08.LIB

SI:77 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

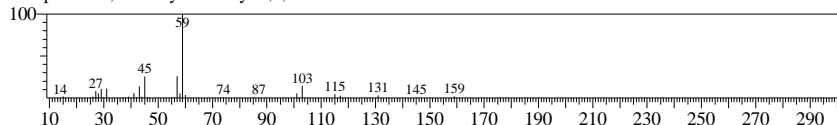
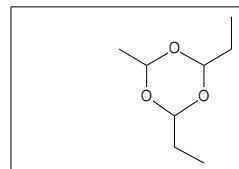
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$S\$ 2-(2-Methoxypropoxy)-1-propanol # \$S\$



Hit#7 Entry:19968 Library:NIST08.LIB

SI:77 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

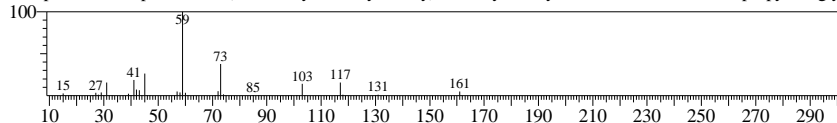
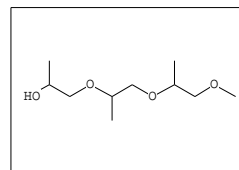
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#8 Entry:46317 Library:NIST08.LIB

SI:77 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

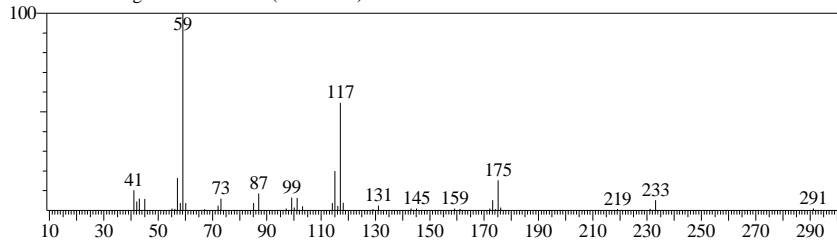
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$S\$ Dowanol 62b \$S\$ Tripropylene gly



<< Target >>

Line#:65 R.Time:27.783(Scan#:2975) BasePeak:59.10(880629)

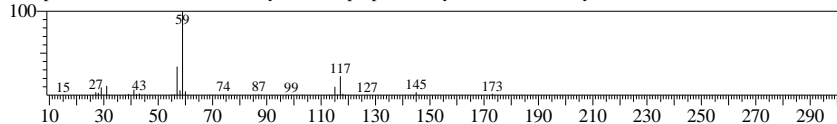
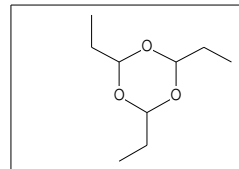
RawMode:Averaged 27.775-27.792(2974-2976) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:80 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

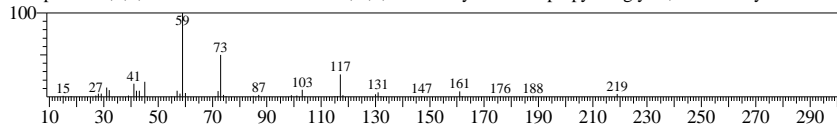
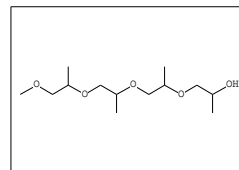
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:85908 Library:NIST08.LIB

SI:78 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

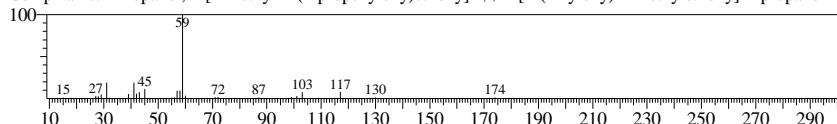
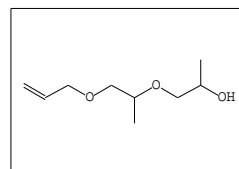
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl-



Hit#:3 Entry:27550 Library:NIST08.LIB

SI:77 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

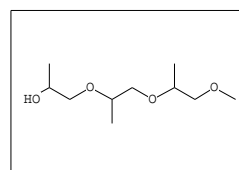
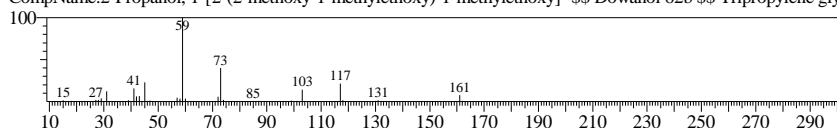
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:4 Entry:17307 Library:NIST08s.LIB

SI:76 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

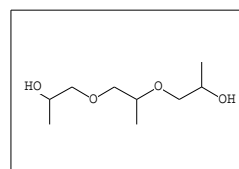
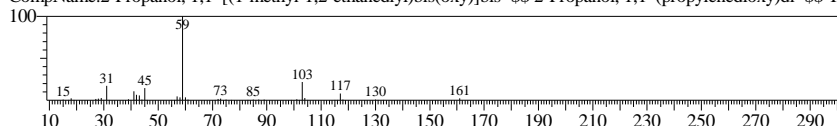
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:5 Entry:37499 Library:NIST08.LIB

SI:76 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

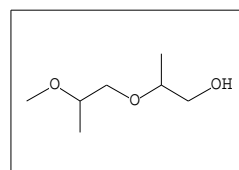
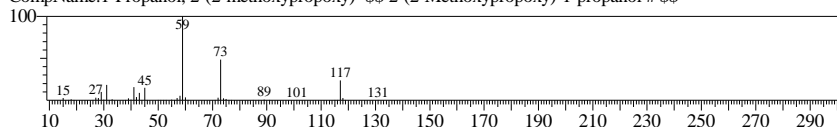
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#:6 Entry:14297 Library:NIST08.LIB

SI:76 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

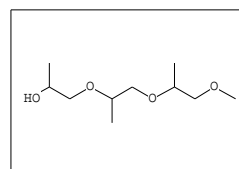
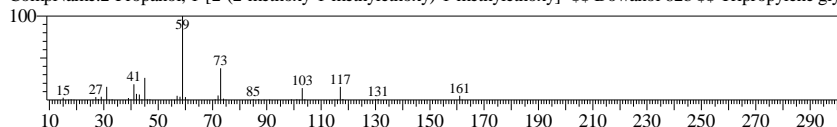
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#:7 Entry:46317 Library:NIST08.LIB

SI:76 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

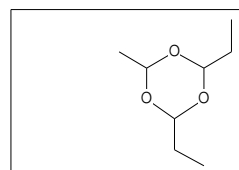
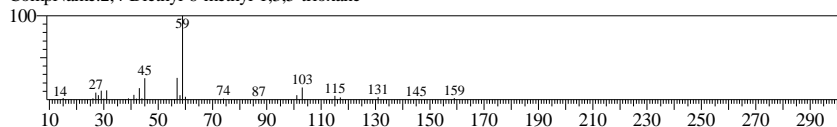
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#:8 Entry:19968 Library:NIST08.LIB

SI:76 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

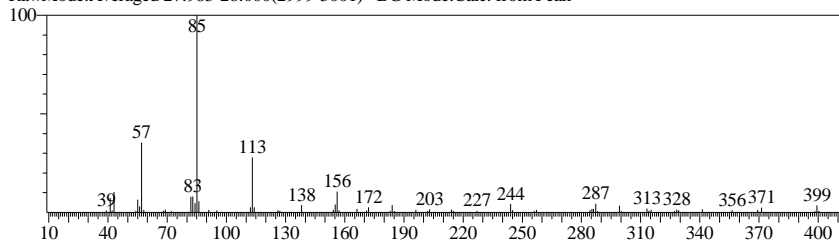
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line#:66 R.Time:27.992(Scan#:3000) BasePeak:85.10(1052779)

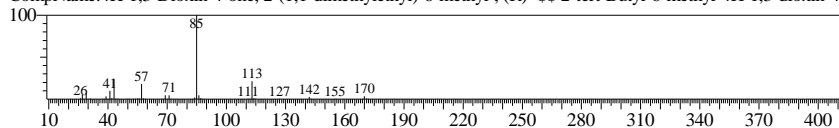
RawMode:Averaged 27.983-28.000(2999-3001) BG Mode:Calc. from Peak



Hit#:1 Entry:25344 Library:NIST08.LIB

SI:76 Formula:C9H14O3 CAS:107289-20-3 MolWeight:170 RetIndex:1280

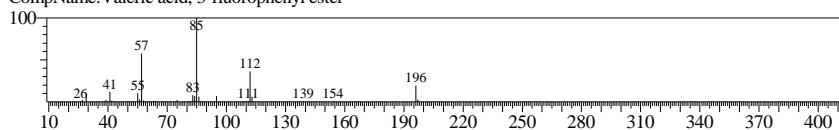
CompName:4H-1,3-Dioxin-4-one, 2-(1,1-dimethylethyl)-6-methyl-, (R)- \$2-tert-Butyl-6-methyl-4H-1,3-dioxin-4



Hit#:2 Entry:40362 Library:NIST08.LIB

SI:74 Formula:C11H13FO2 CAS:0-00-0 MolWeight:196 RetIndex:1333

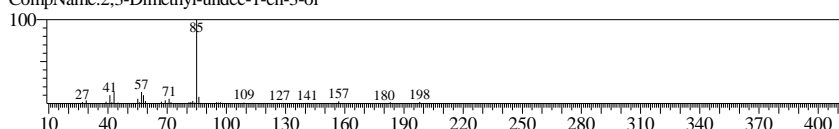
CompName:Valeric acid, 3-fluorophenyl ester



Hit#:3 Entry:41887 Library:NIST08.LIB

SI:72 Formula:C13H26O CAS:0-00-0 MolWeight:198 RetIndex:1372

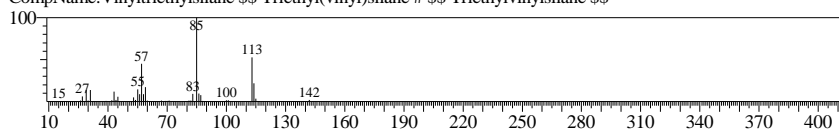
CompName:2,3-Dimethyl-undec-1-en-3-ol



Hit#:4 Entry:12147 Library:NIST08.LIB

SI:71 Formula:C8H18Si CAS:1112-54-5 MolWeight:142 RetIndex:717

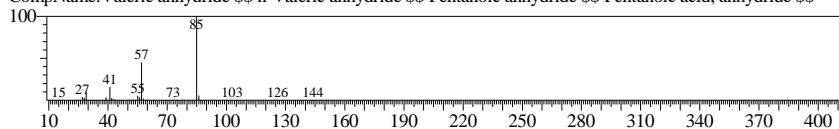
CompName:Vinyltriethylsilane \$Triethyl(vinyl)silane # \$Triethylvinylsilane \$



Hit#:5 Entry:34397 Library:NIST08.LIB

SI:71 Formula:C10H18O3 CAS:2082-59-9 MolWeight:186 RetIndex:1319

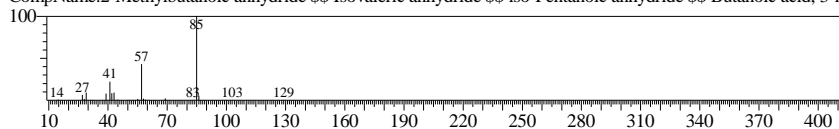
CompName:Valeric anhydride \$n-Valeric anhydride \$Pentanoic anhydride \$Pentanoic acid, anhydride \$



Hit#:6 Entry:34398 Library:NIST08.LIB

SI:70 Formula:C10H18O3 CAS:1468-39-9 MolWeight:186 RetIndex:1190

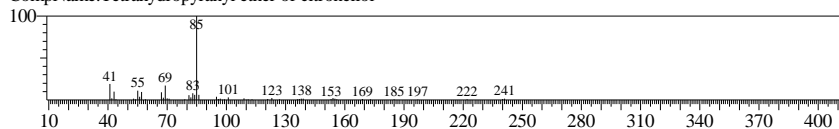
CompName:2-Methylbutanoic anhydride \$Isovaleric anhydride \$iso-Pentanoic anhydride \$Butanoic acid, 3-methyl-



Hit#:7 Entry:69538 Library:NIST08.LIB

SI:70 Formula:C15H28O2 CAS:0-00-0 MolWeight:240 RetIndex:1661

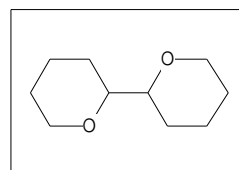
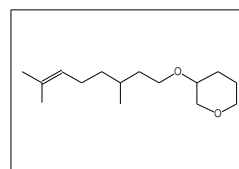
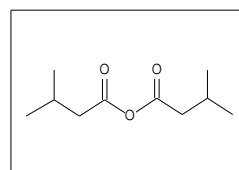
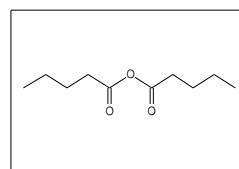
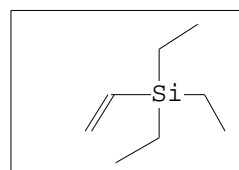
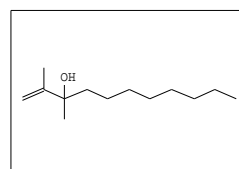
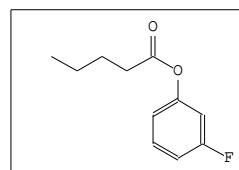
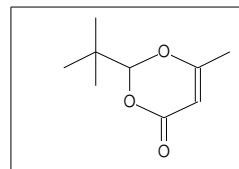
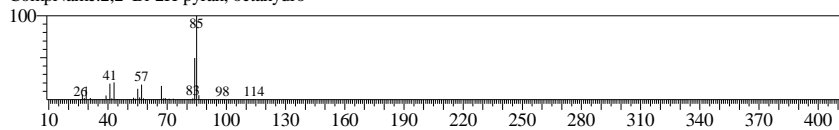
CompName:Tetrahydropyranyl ether of citronellol



Hit#:8 Entry:25558 Library:NIST08.LIB

SI:69 Formula:C10H18O2 CAS:16282-29-4 MolWeight:170 RetIndex:1319

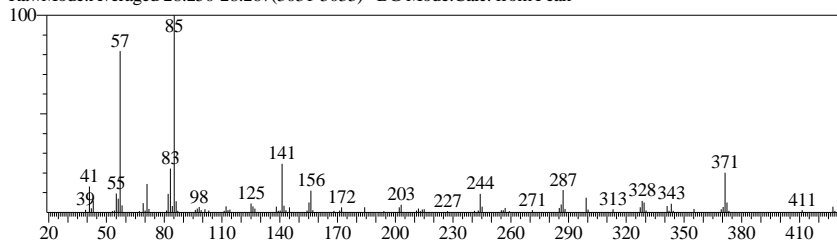
CompName:2,2'-Bi-2H-pyran, octahydro-



<< Target >>

Line#:67 R.Time:28.258(Scan#:3032) BasePeak:85.10(1020510)

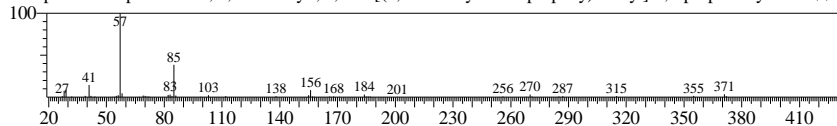
RawMode:Averaged 28.250-28.267(3031-3033) BG Mode:Calc. from Peak



Hit#:1 Entry:181922 Library:NIST08.LIB

SI:67 Formula:C25H44O8 CAS:5178-17-6 MolWeight:472 RetIndex:2753

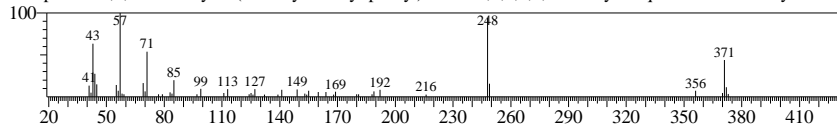
CompName:Propanoic acid, 2,2-dimethyl-, 2,2-bis[(2,2-dimethyl-1-oxopropoxy)methyl]-1,3-propanediyl ester \$S I



Hit#:2 Entry:156076 Library:NIST08.LIB

SI:52 Formula:C21H25NO3S CAS:0-00-0 MolWeight:371 RetIndex:3018

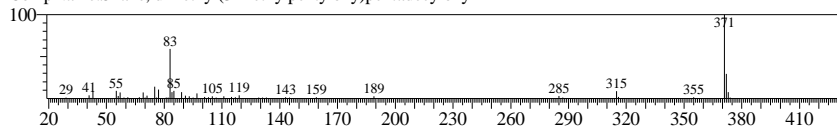
CompName:2,7,7-Trimethyl-4-(4-methylsulfanyl-phenyl)-5-oxo-1,4,5,6,7,8-hexahydro-quinoline-3-carboxylic acid



Hit#:3 Entry:162793 Library:NIST08.LIB

SI:49 Formula:C23H50O2Si CAS:0-00-0 MolWeight:386 RetIndex:2305

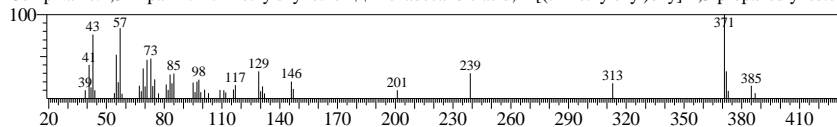
CompName:Silane, dimethyl(3-methylpentyl)oxy)pentadecyloxy-



Hit#:4 Entry:28229 Library:NIST08s.LIB

SI:46 Formula:C38H76O5Si CAS:53212-95-6 MolWeight:640 RetIndex:4055

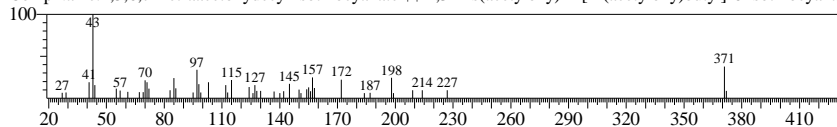
CompName:1,3-Dipalmitin trimethylsilyl ether \$S Hexadecanoic acid, 2-[(trimethylsilyl)oxy]-1,3-propanediyl ester



Hit#:5 Entry:175569 Library:NIST08.LIB

SI:46 Formula:C19H29NO8S CAS:57103-44-3 MolWeight:431 RetIndex:0

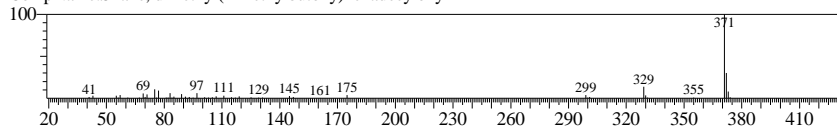
CompName:4,5,6,7-Tetraacetoxyldecyl isothiocyanate \$S 2,3-Bis(acetyloxy)-1-[1-(acetyloxy)butyl]-6-isothiocyana



Hit#:6 Entry:162800 Library:NIST08.LIB

SI:46 Formula:C23H50O2Si CAS:0-00-0 MolWeight:386 RetIndex:2305

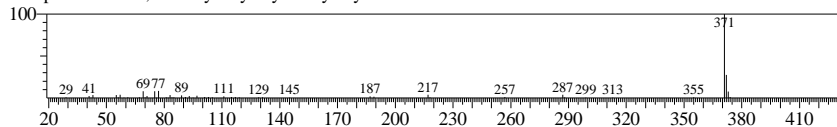
CompName:Silane, dimethyl(2-methylbutoxy)hexadecyloxy-



Hit#:7 Entry:162795 Library:NIST08.LIB

SI:46 Formula:C23H50O2Si CAS:0-00-0 MolWeight:386 RetIndex:2369

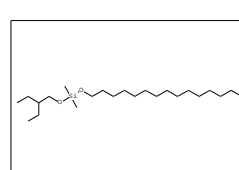
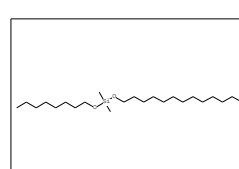
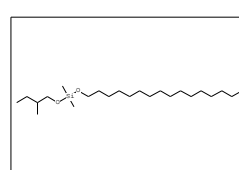
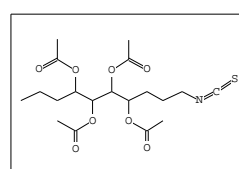
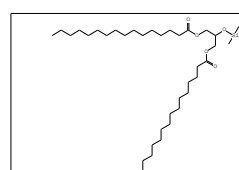
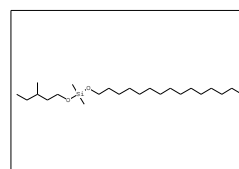
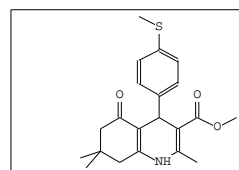
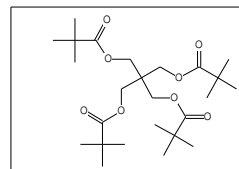
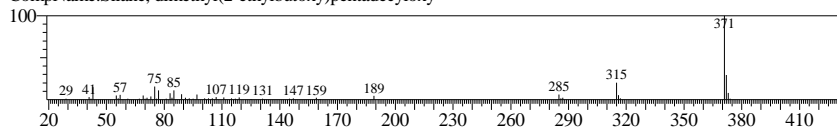
CompName:Silane, dimethyloctyloxytridecyloxy-



Hit#:8 Entry:162799 Library:NIST08.LIB

SI:45 Formula:C23H50O2Si CAS:0-00-0 MolWeight:386 RetIndex:2305

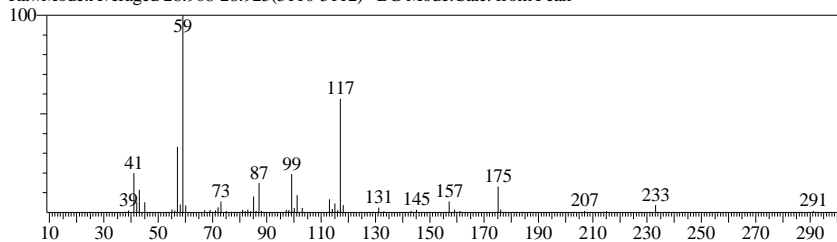
CompName:Silane, dimethyl(2-ethylbutoxy)pentadecyloxy-



<< Target >>

Line# 68 R.Time:28.917(Scan#:3111) BasePeak:59.10(299099)

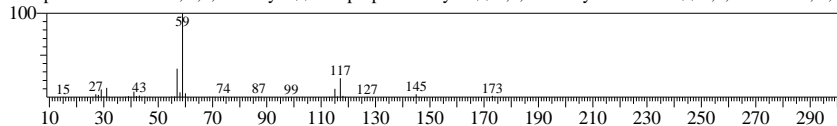
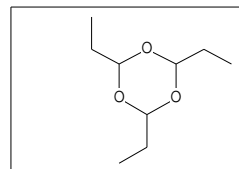
RawMode:Averaged 28.908-28.925(3110-3112) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:79 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

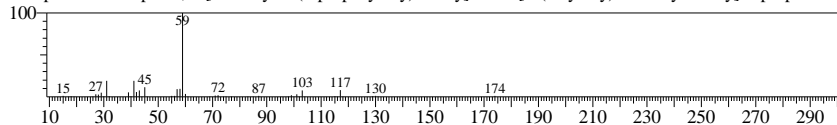
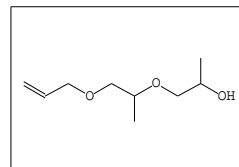
CompName:s-Trioxane, 2,4,6-triethyl- \$- Parapropionaldehyde \$- 2,4,6-Triethyl-s-trioxane \$- 1,3,5-Trioxane, 2,4



Hit#2 Entry:27550 Library:NIST08.LIB

SI:76 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

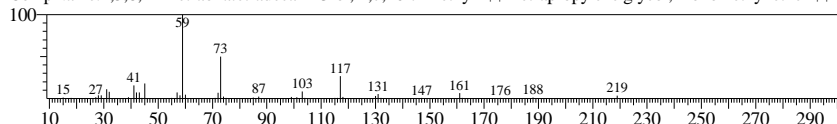
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#3 Entry:85908 Library:NIST08.LIB

SI:76 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

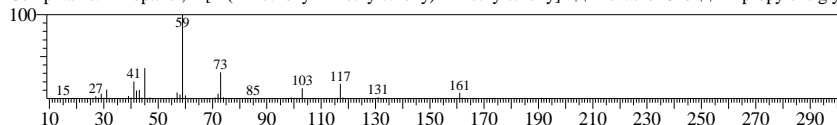
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$- Tetrapropylene glycol, monomethyl ether \$- 4



Hit#4 Entry:17306 Library:NIST08.LIB

SI:75 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

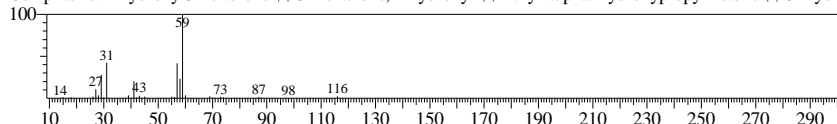
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#5 Entry:4568 Library:NIST08.LIB

SI:75 Formula:C6H12O2 CAS:4984-85-4 MolWeight:116 RetIndex:916

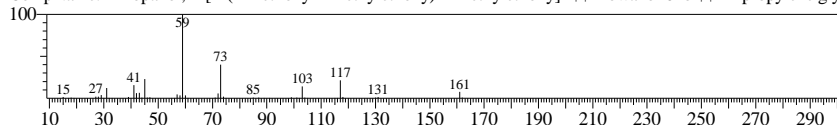
CompName:4-Hydroxy-3-hexanone \$- 3-Hexanone, 4-hydroxy- \$- Ethyl. alpha.-hydroxypropyl ketone \$- 3-Hydr



Hit#6 Entry:17307 Library:NIST08.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

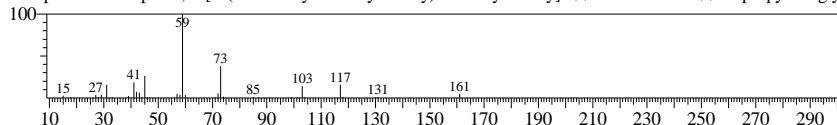
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#7 Entry:46317 Library:NIST08.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

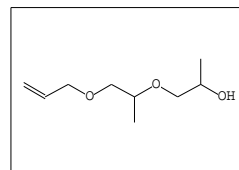
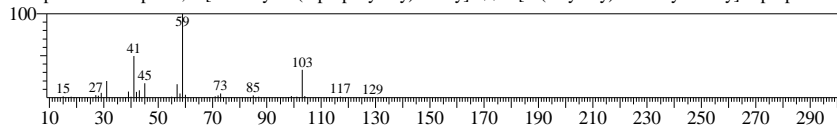
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$- Dowanol 62b \$- Tripropylene gly



Hit#8 Entry:12792 Library:NIST08.LIB

SI:74 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

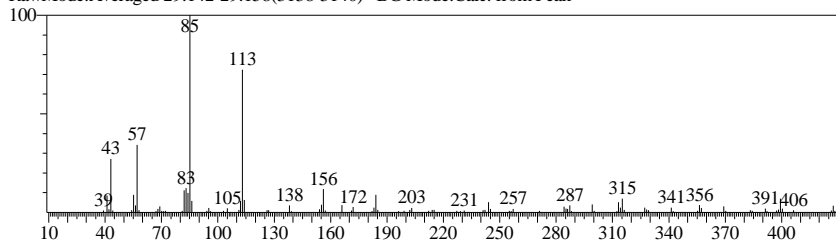
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$- 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



<< Target >>

Line#:69 R.Time:29.150(Scan#:3139) BasePeak:85.10(1054189)

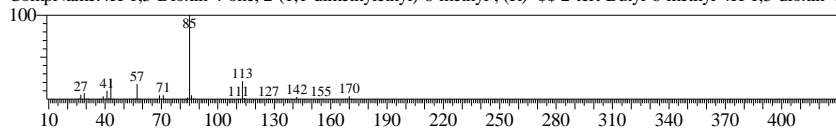
RawMode:Averaged 29.142-29.158(3138-3140) BG Mode:Calc. from Peak



Hit#:1 Entry:25344 Library:NIST08.LIB

SI:70 Formula:C₉H₁₄O₃ CAS:107289-20-3 MolWeight:170 RetIndex:1280

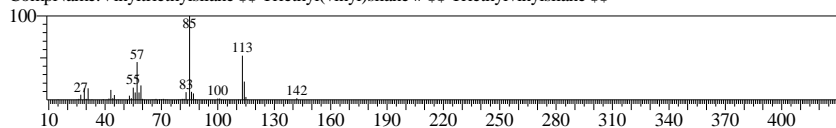
CompName:4H-1,3-Dioxin-4-one, 2-(1,1-dimethylethyl)-6-methyl-, (R)-



Hit#:2 Entry:12147 Library:NIST08.LIB

SI:68 Formula:C₈H₁₈Si CAS:1112-54-5 MolWeight:142 RetIndex:717

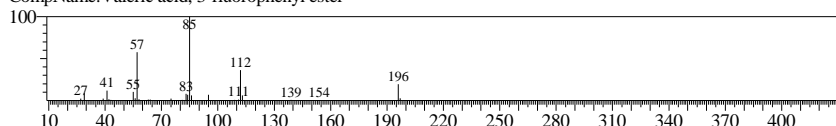
CompName:Vinyltriethylsilane



Hit#:3 Entry:40362 Library:NIST08.LIB

SI:67 Formula:C₁₁H₁₃FO₂ CAS:0-00-0 MolWeight:196 RetIndex:1333

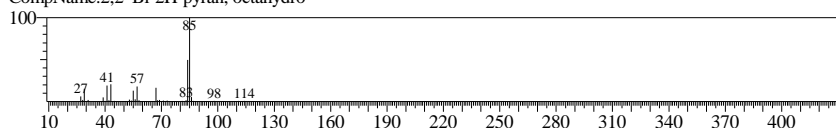
CompName:Valeric acid, 3-fluorophenyl ester



Hit#:4 Entry:25558 Library:NIST08.LIB

SI:66 Formula:C₁₀H₁₈O₂ CAS:16282-29-4 MolWeight:170 RetIndex:1319

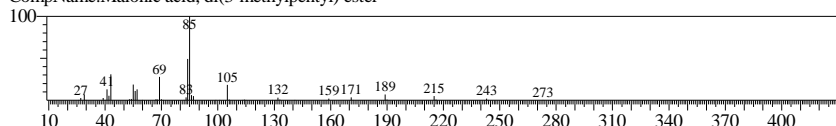
CompName:2,2'-Bi-2H-pyran, octahydro-



Hit#:5 Entry:91967 Library:NIST08.LIB

SI:65 Formula:C₁₅H₂₈O₄ CAS:0-00-0 MolWeight:272 RetIndex:1719

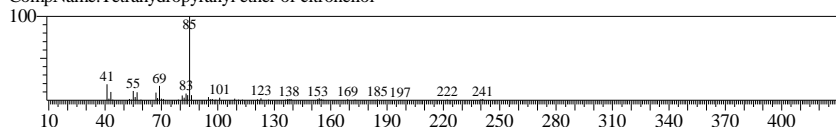
CompName:Malonic acid, di(3-methylpentyl) ester



Hit#:6 Entry:69538 Library:NIST08.LIB

SI:65 Formula:C₁₅H₂₈O₂ CAS:0-00-0 MolWeight:240 RetIndex:1661

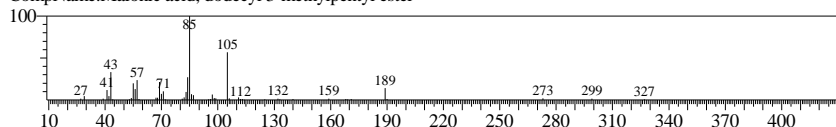
CompName:Tetrahydropyranyl ether of citronellol



Hit#:7 Entry:148358 Library:NIST08.LIB

SI:65 Formula:C₂₁H₄₀O₄ CAS:0-00-0 MolWeight:356 RetIndex:2379

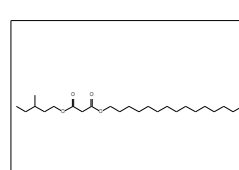
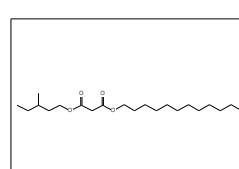
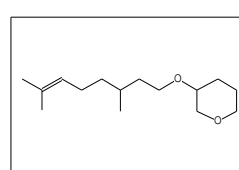
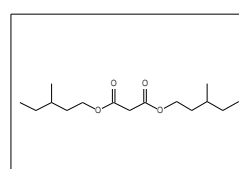
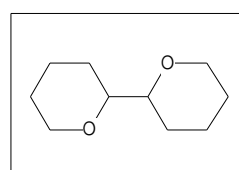
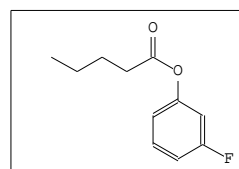
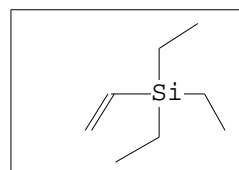
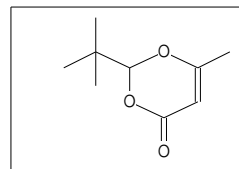
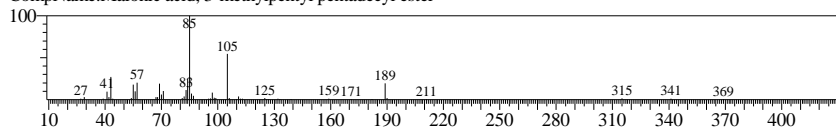
CompName:Malonic acid, dodecyl 3-methylpentyl ester



Hit#:8 Entry:167197 Library:NIST08.LIB

SI:65 Formula:C₂₄H₄₆O₄ CAS:0-00-0 MolWeight:398 RetIndex:2677

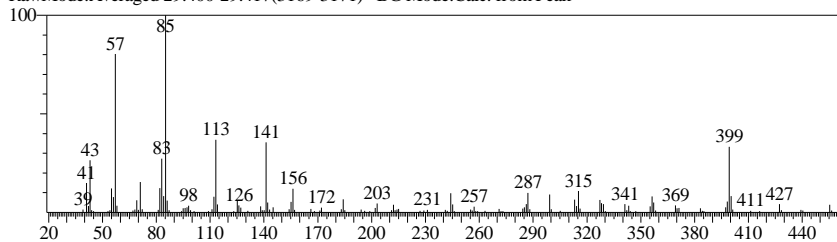
CompName:Malonic acid, 3-methylpentyl pentadecyl ester



<< Target >>

Line#:70 R.Time:29.408(Scan#:3170) BasePeak:85.10(1966373)

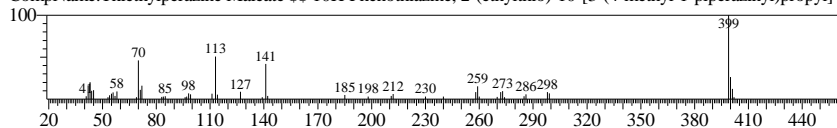
RawMode:Averaged 29.400-29.417(3169-3171) BG Mode:Calc. from Peak



Hit#:1 Entry:167503 Library:NIST08.LIB

SI:52 Formula:C22H29N3S2 CAS:1179-69-7 MolWeight:399 RetIndex:3355

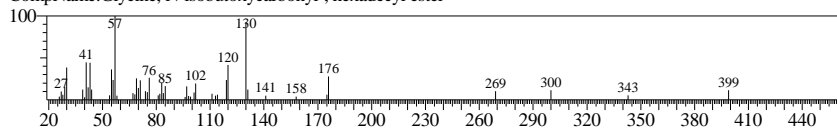
CompName:Thiethylpiperazine Maleate SS 10H-Phenothiazine, 2-(ethylthio)-10-[3-(4-methyl-1-piperazinyl)propyl]-



Hit#:2 Entry:167540 Library:NIST08.LIB

SI:48 Formula:C23H45NO4 CAS:0-00-0 MolWeight:399 RetIndex:2776

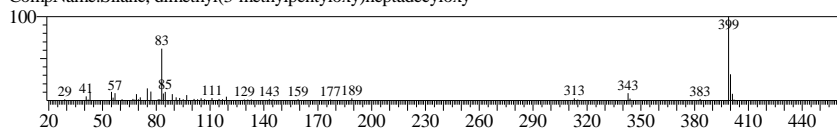
CompName:Glycine, N-isobutoxycarbonyl-, hexadecyl ester



Hit#:3 Entry:171775 Library:NIST08.LIB

SI:48 Formula:C25H54O2Si CAS:0-00-0 MolWeight:414 RetIndex:2504

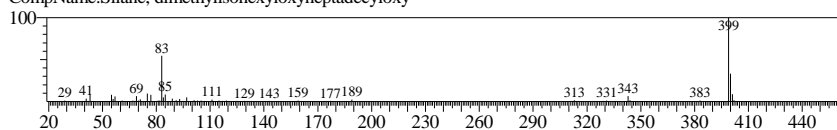
CompName:Silane, dimethyl(3-methylpentyl)oxyheptadecyloxy-



Hit#:4 Entry:171774 Library:NIST08.LIB

SI:47 Formula:C25H54O2Si CAS:0-00-0 MolWeight:414 RetIndex:2504

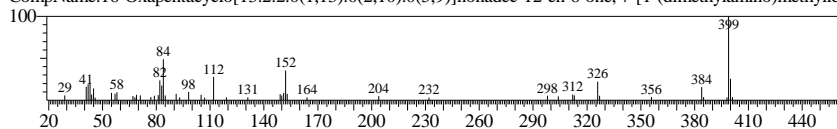
CompName:Silane, dimethylisohexyloxyheptadecyloxy-



Hit#:5 Entry:167658 Library:NIST08.LIB

SI:47 Formula:C25H37NO3 CAS:0-00-0 MolWeight:399 RetIndex:2637

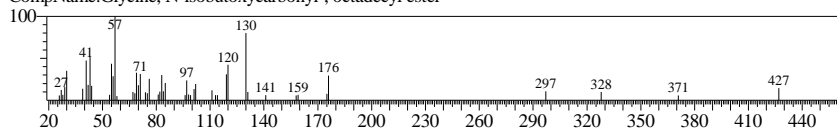
CompName:16-Oxapentacyclo[13.2.2.0(1,13).0(2,10).0(5,9)]nonadec-12-en-6-one, 7-[1-(dimethylamino)methylid]



Hit#:6 Entry:174750 Library:NIST08.LIB

SI:47 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975

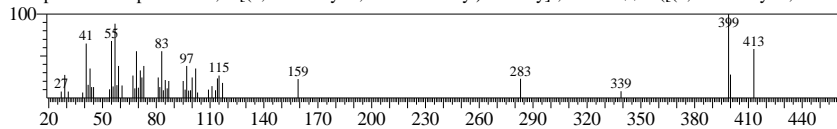
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



Hit#:7 Entry:175018 Library:NIST08.LIB

SI:45 Formula:C25H48O5 CAS:57346-64-2 MolWeight:428 RetIndex:2841

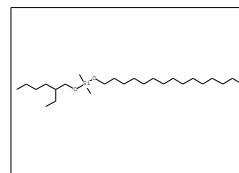
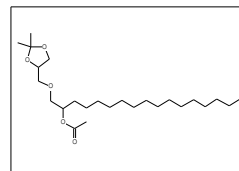
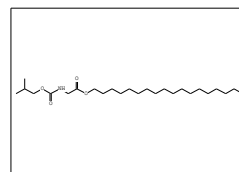
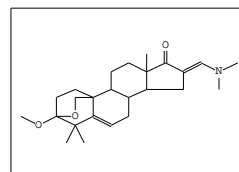
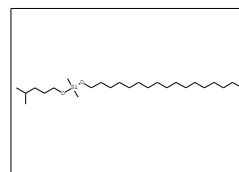
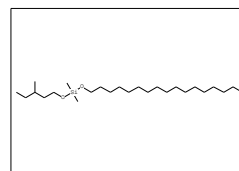
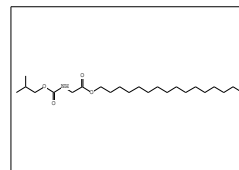
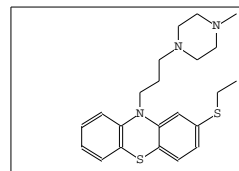
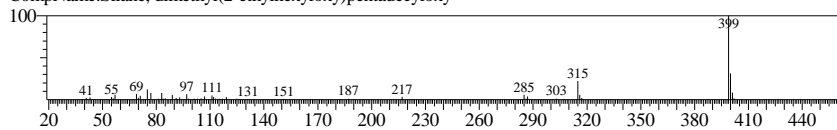
CompName:2-Heptadecanol, 1-[(2,2-dimethyl-1,3-dioxolan-4-yl)methoxy]-, acetate SS 1-[(2,2-Dimethyl-1,3-diox



Hit#:8 Entry:171776 Library:NIST08.LIB

SI:45 Formula:C25H54O2Si CAS:0-00-0 MolWeight:414 RetIndex:2504

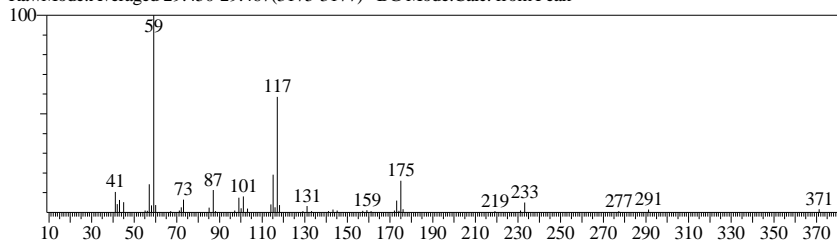
CompName:Silane, dimethyl(2-ethylhexyloxy)pentadecyloxy-



<< Target >>

Line# 71 R.Time:29.458(Scan#:3176) BasePeak:59.10(720179)

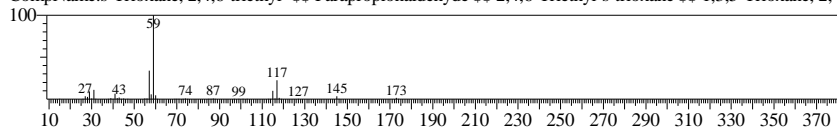
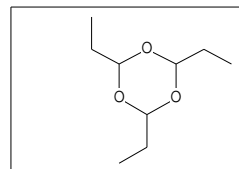
RawMode:Averaged 29.450-29.467(3175-3177) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:78 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

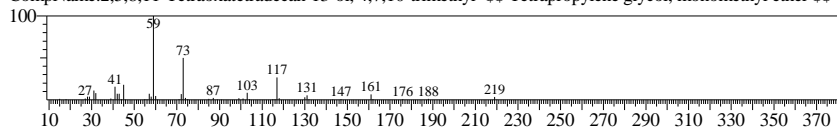
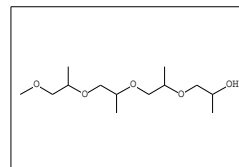
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:77 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

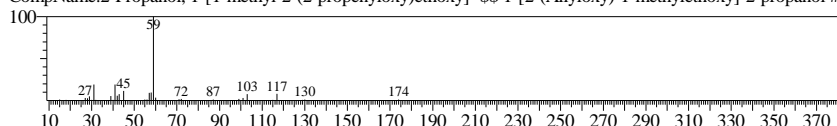
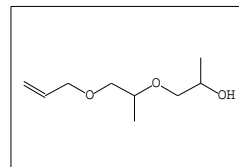
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#3 Entry:27550 Library:NIST08.LIB

SI:76 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

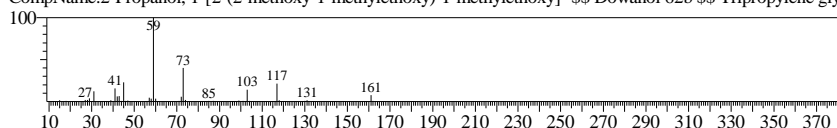
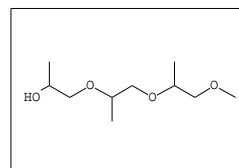
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:17307 Library:NIST08s.LIB

SI:75 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

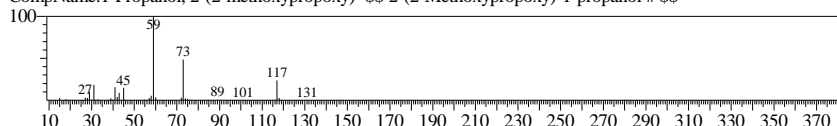
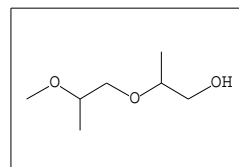
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#5 Entry:14297 Library:NIST08.LIB

SI:74 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

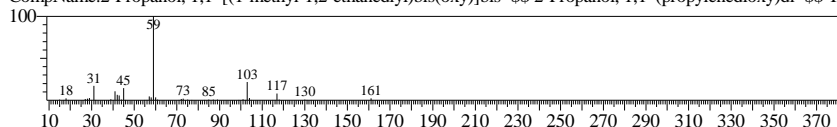
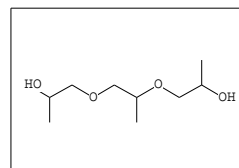
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#6 Entry:37499 Library:NIST08.LIB

SI:74 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

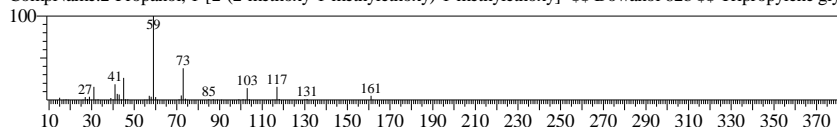
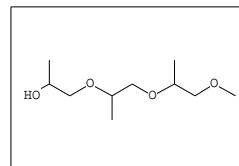
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#7 Entry:46317 Library:NIST08.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

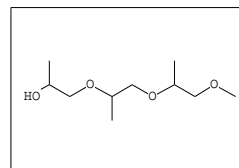
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#8 Entry:17306 Library:NIST08s.LIB

SI:74 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

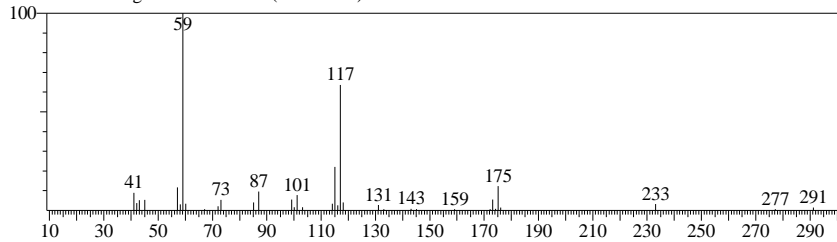
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line# 72 R.Time:29.517(Scan#:3183) BasePeak:59.10(503044)

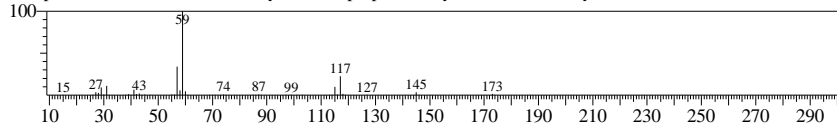
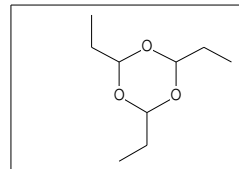
RawMode:Averaged 29.508-29.525(3182-3184) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:80 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

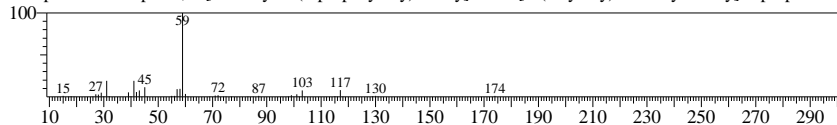
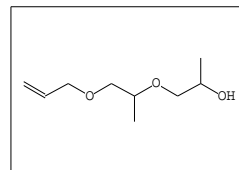
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:27550 Library:NIST08.LIB

SI:77 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

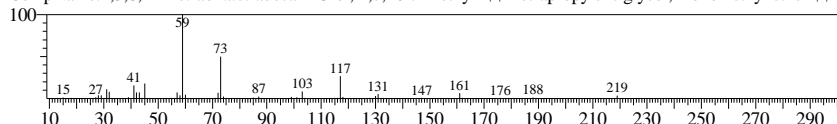
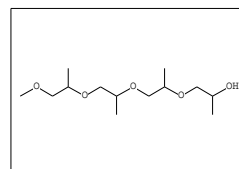
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#3 Entry:85908 Library:NIST08.LIB

SI:76 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

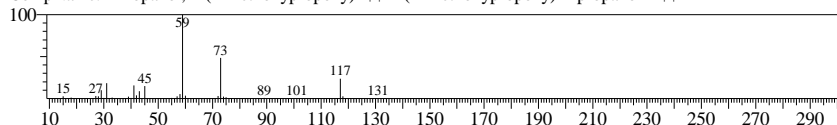
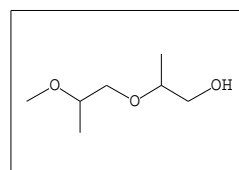
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#4 Entry:14297 Library:NIST08.LIB

SI:76 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

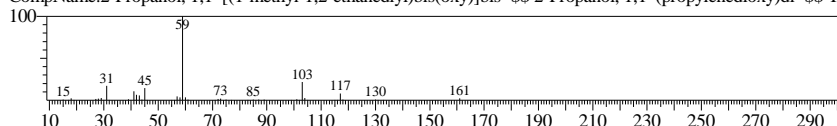
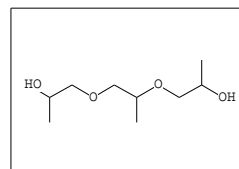
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol #



Hit#5 Entry:37499 Library:NIST08.LIB

SI:76 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

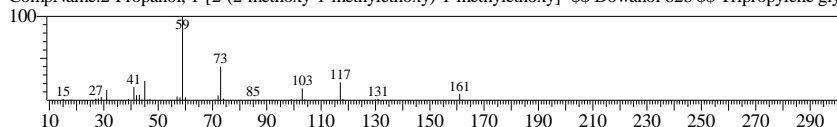
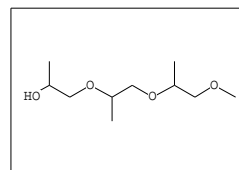
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#6 Entry:17307 Library:NIST08s.LIB

SI:75 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

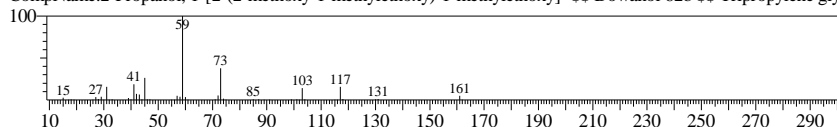
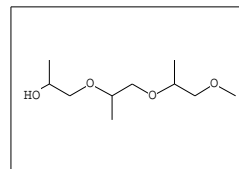
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#7 Entry:46317 Library:NIST08.LIB

SI:75 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

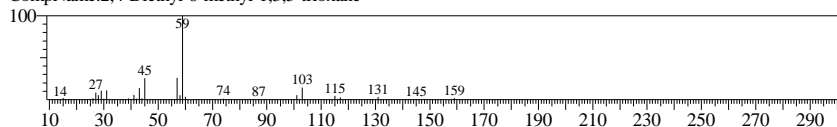
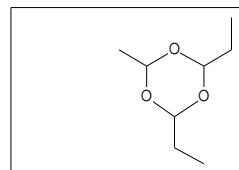
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#8 Entry:19968 Library:NIST08.LIB

SI:75 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

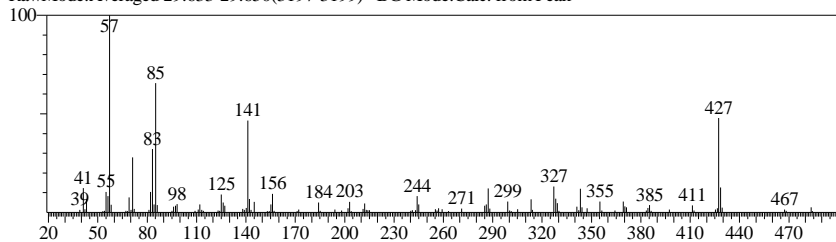
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



<< Target >>

Line#:73 R.Time:29.642(Scan#:3198) BasePeak:57.10(1794928)

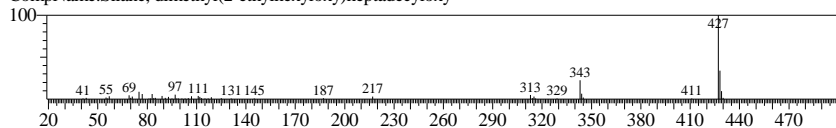
RawMode:Averaged 29.633-29.650(3197-3199) BG Mode:Calc. from Peak



Hit#:1 Entry:177733 Library:NIST08.LIB

SI:49 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2703

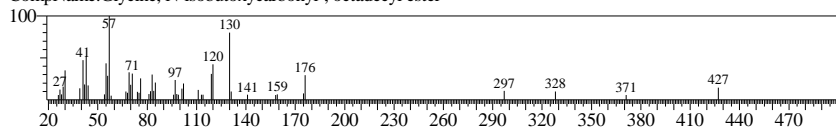
CompName:Silane, dimethyl(2-ethylhexyloxy)heptadecyloxy-



Hit#:2 Entry:174750 Library:NIST08.LIB

SI:49 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975

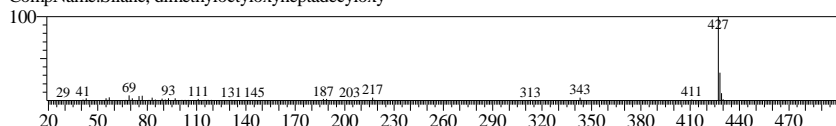
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



Hit#:3 Entry:177734 Library:NIST08.LIB

SI:47 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

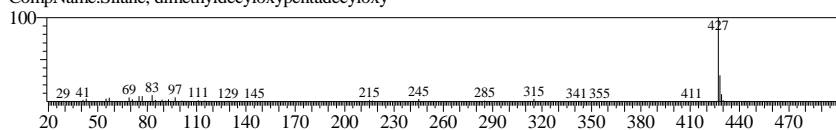
CompName:Silane, dimethyloctyloxyheptadecyloxy-



Hit#:4 Entry:177735 Library:NIST08.LIB

SI:47 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

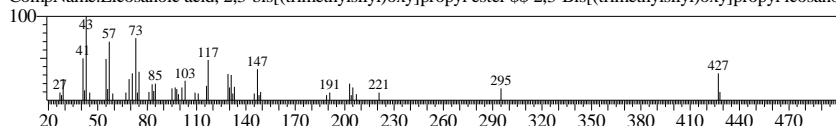
CompName:Silane, dimethyldecyloxy pentadecyloxy-



Hit#:5 Entry:186484 Library:NIST08.LIB

SI:42 Formula:C29H62O4Si2 CAS:55517-94-7 MolWeight:530 RetIndex:2979

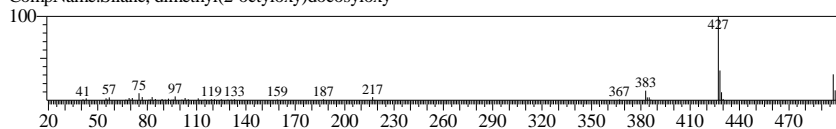
CompName:Eicosanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$ 2,3-Bis[(trimethylsilyl)oxy]propyl icosano



Hit#:6 Entry:185523 Library:NIST08.LIB

SI:41 Formula:C32H68O2Si CAS:0-00-0 MolWeight:512 RetIndex:3200

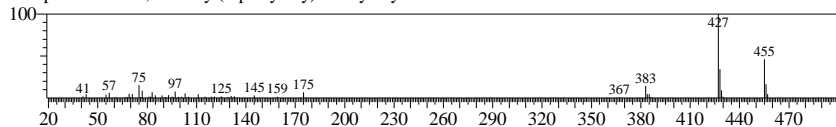
CompName:Silane, dimethyl(2-octyloxy)docosyloxy-



Hit#:7 Entry:181721 Library:NIST08.LIB

SI:41 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902

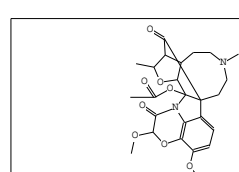
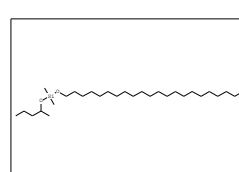
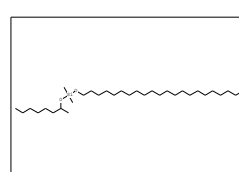
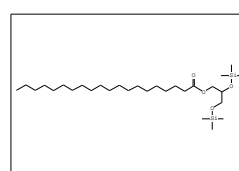
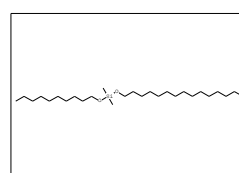
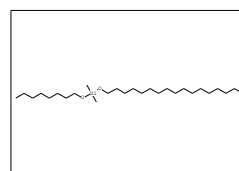
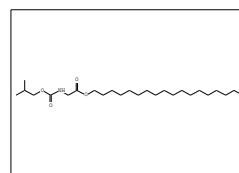
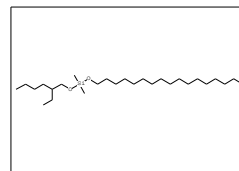
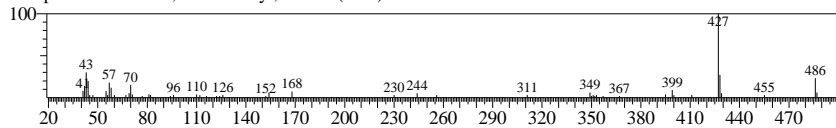
CompName:Silane, dimethyl(2-pentyloxy)docosyloxy-



Hit#:8 Entry:183457 Library:NIST08.LIB

SI:39 Formula:C25H30N2O8 CAS:29474-86-0 MolWeight:486 RetIndex:3525

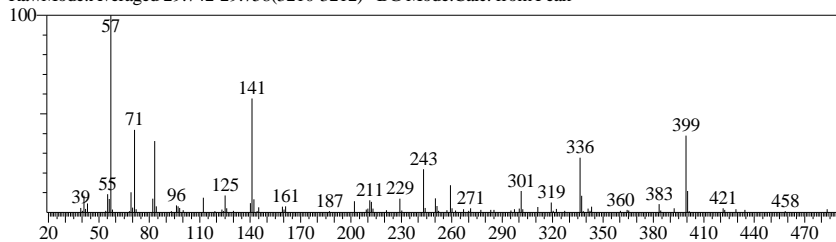
CompName:Dichotine, 11-methoxy-, acetate (ester)



<< Target >>

Line#:74 R.Time:29.750(Scan#:3211) BasePeak:57.10(58753)

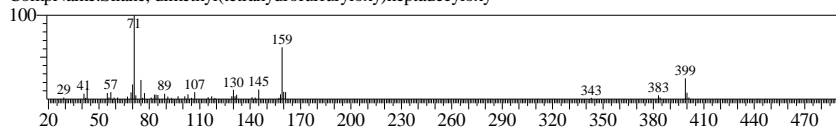
RawMode:Averaged 29.742-29.758(3210-3212) BG Mode:Calc. from Peak



Hit#:1 Entry:171746 Library:NIST08.LIB

SI:48 Formula:C24H50O3Si CAS:0-00-0 MolWeight:414 RetIndex:2600

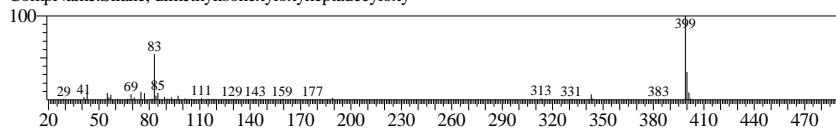
CompName:Silane, dimethyl(tetrahydrofurfuryloxy)heptadecyloxy-



Hit#:2 Entry:171774 Library:NIST08.LIB

SI:48 Formula:C25H54O2Si CAS:0-00-0 MolWeight:414 RetIndex:2504

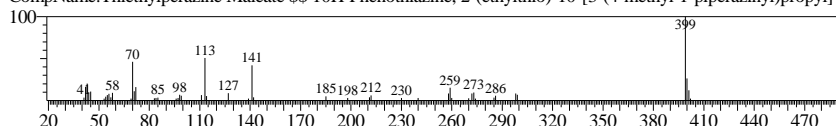
CompName:Silane, dimethylisohexyloxyheptadecyloxy-



Hit#:3 Entry:167503 Library:NIST08.LIB

SI:48 Formula:C22H29N3S2 CAS:1179-69-7 MolWeight:399 RetIndex:3355

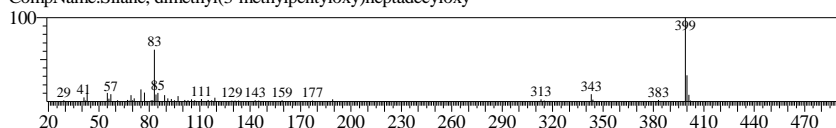
CompName:Thiethylperazine Maleate \$S 10H-Phenothiazine, 2-(ethylthio)-10-[3-(4-methyl-1-piperazinyl)propyl]-



Hit#:4 Entry:171775 Library:NIST08.LIB

SI:47 Formula:C25H54O2Si CAS:0-00-0 MolWeight:414 RetIndex:2504

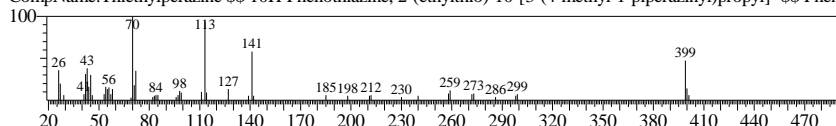
CompName:Silane, dimethyl(3-methylpentyloxy)heptadecyloxy-



Hit#:5 Entry:167502 Library:NIST08.LIB

SI:46 Formula:C22H29N3S2 CAS:1420-55-9 MolWeight:399 RetIndex:3355

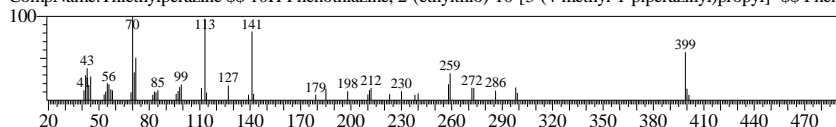
CompName:Thiethylperazine \$S 10H-Phenothiazine, 2-(ethylthio)-10-[3-(4-methyl-1-piperazinyl)propyl]- \$S Phen



Hit#:6 Entry:27289 Library:NIST08s.LIB

SI:46 Formula:C22H29N3S2 CAS:1420-55-9 MolWeight:399 RetIndex:3355

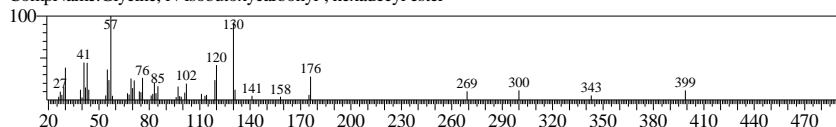
CompName:Thiethylperazine \$S 10H-Phenothiazine, 2-(ethylthio)-10-[3-(4-methyl-1-piperazinyl)propyl]- \$S Phen



Hit#:7 Entry:167540 Library:NIST08.LIB

SI:45 Formula:C23H45NO4 CAS:0-00-0 MolWeight:399 RetIndex:2776

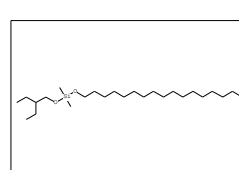
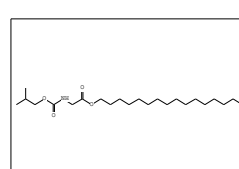
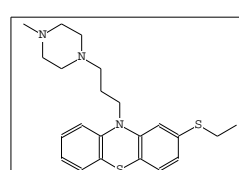
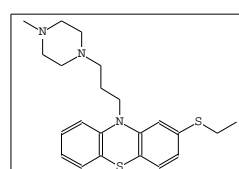
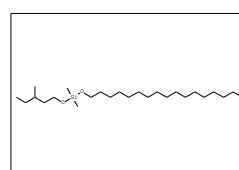
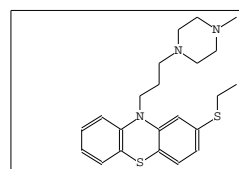
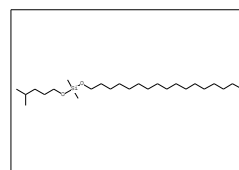
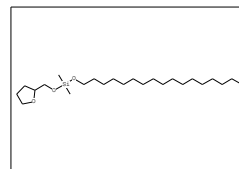
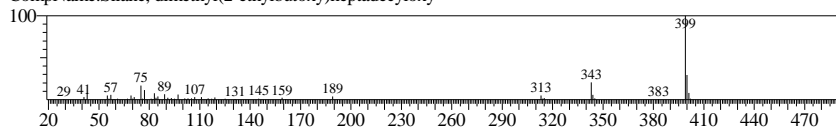
CompName:Glycine, N-isobutoxycarbonyl-, hexadecyl ester



Hit#:8 Entry:171777 Library:NIST08.LIB

SI:43 Formula:C25H54O2Si CAS:0-00-0 MolWeight:414 RetIndex:2504

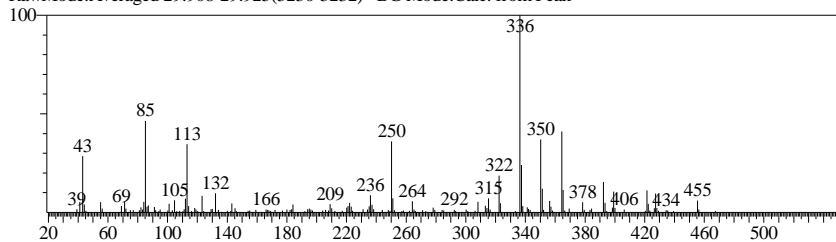
CompName:Silane, dimethyl(2-ethylbutoxy)heptadecyloxy-



<< Target >>

Line#:75 R.Time:29.917(Scan#:3231) BasePeak:336.30(50121)

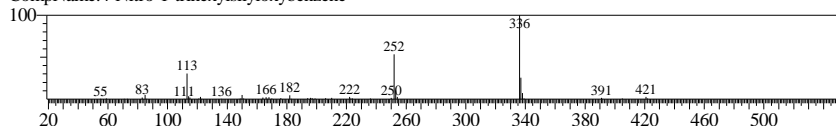
RawMode:Averaged 29.908-29.925(3230-3232) BG Mode:Calc. from Peak



Hit#:1 Entry:173409 Library:NIST08.LIB

SI:47 Formula:C24H43NO3Si CAS:0-00-0 MolWeight:421 RetIndex:2865

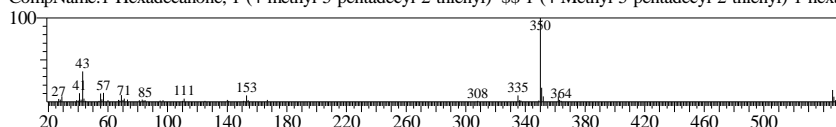
CompName:4-Nitro-1-trihexylsilyloxybenzene



Hit#:2 Entry:187267 Library:NIST08.LIB

SI:40 Formula:C36H66OS CAS:59782-69-3 MolWeight:546 RetIndex:3992

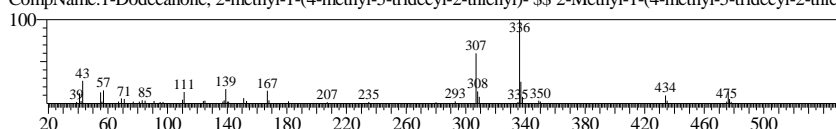
CompName:1-Hexadecanone, 1-(4-methyl-5-pentadecyl-2-thienyl)- \$- 1-(4-Methyl-5-pentadecyl-2-thienyl)-1-hexa



Hit#:3 Entry:182423 Library:NIST08.LIB

SI:40 Formula:C31H56OS CAS:60089-65-8 MolWeight:476 RetIndex:3431

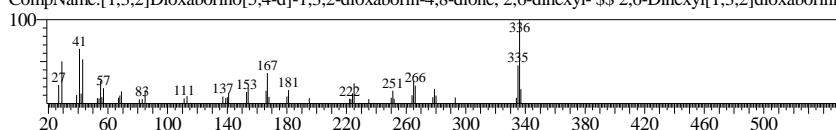
CompName:1-Dodecanone, 2-methyl-1-(4-methyl-5-tridecyl-2-thienyl)- \$- 2-Methyl-1-(4-methyl-5-tridecyl-2-thie



Hit#:4 Entry:136427 Library:NIST08.LIB

SI:39 Formula:C16H26B2O6 CAS:87841-15-4 MolWeight:336 RetIndex:0

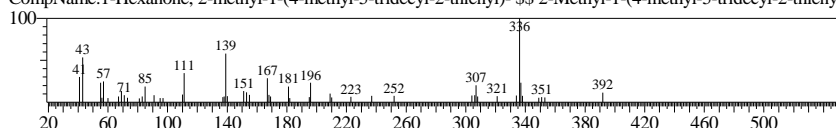
CompName:[1,3,2]Dioxaborino[5,4-d]-1,3,2-dioxaborin-4,8-dione, 2,6-dihexyl- \$- 2,6-Dihexyl[1,3,2]dioxaborini



Hit#:5 Entry:165121 Library:NIST08.LIB

SI:39 Formula:C25H44OS CAS:60089-62-5 MolWeight:392 RetIndex:2835

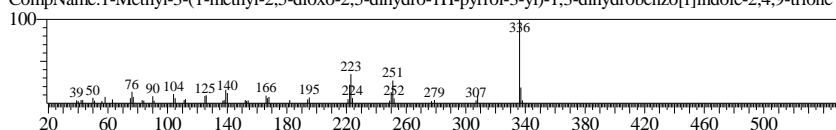
CompName:1-Hexanone, 2-methyl-1-(4-methyl-5-tridecyl-2-thienyl)- \$- 2-Methyl-1-(4-methyl-5-tridecyl-2-thienyl)



Hit#:6 Entry:136504 Library:NIST08.LIB

SI:39 Formula:C18H12N2O5 CAS:0-00-0 MolWeight:336 RetIndex:3135

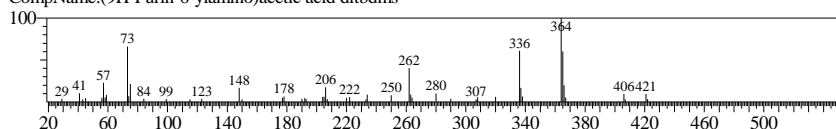
CompName:1-Methyl-3-(1-methyl-2,5-dioxo-2,5-dihydro-1H-pyrrol-3-yl)-1,3-dihydrobenzo[f]indole-2,4,9-trione



Hit#:7 Entry:173363 Library:NIST08.LIB

SI:38 Formula:C19H35N5O2Si2 CAS:0-00-0 MolWeight:421 RetIndex:2490

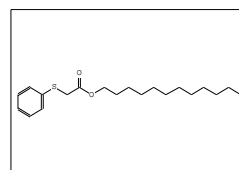
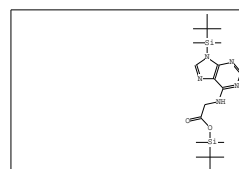
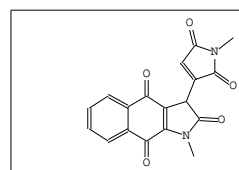
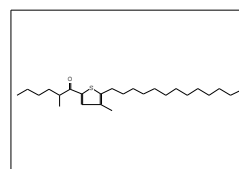
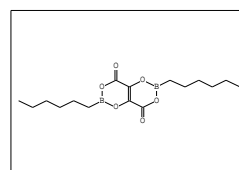
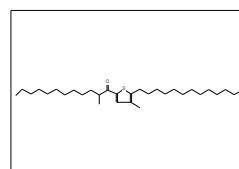
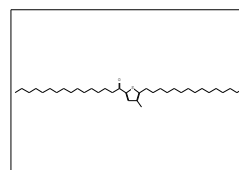
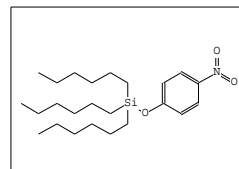
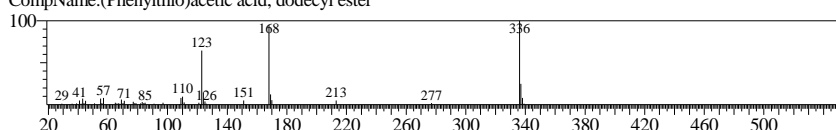
CompName:(9H-Purin-6-ylamino)acetic acid ditbdms



Hit#:8 Entry:136681 Library:NIST08.LIB

SI:38 Formula:C20H32O2S CAS:0-00-0 MolWeight:336 RetIndex:2504

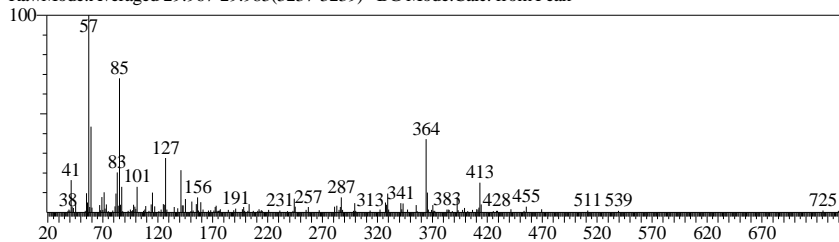
CompName:(Phenylthio)acetic acid, dodecyl ester



<< Target >>

Line#:76 R.Time:29.975(Scan#:3238) BasePeak:57.10(44796)

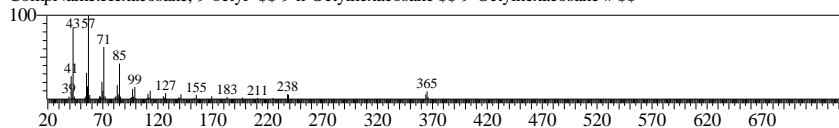
RawMode:Averaged 29.967-29.983(3237-3239) BG Mode:Calc. from Peak



Hit#:1 Entry:182631 Library:NIST08.LIB

SI:54 Formula:C34H70 CAS:55429-83-9 MolWeight:478 RetIndex:3337

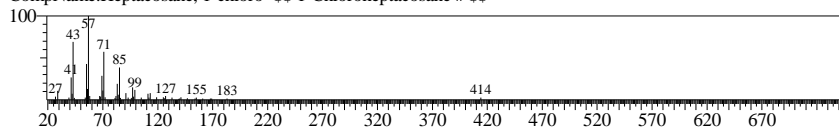
CompName:Hexacosane, 9-octyl- \$\$ 9-n-Octylhexacosane \$\$ 9-Octylhexacosane # \$\$



Hit#:2 Entry:171834 Library:NIST08.LIB

SI:52 Formula:C27H55Cl CAS:62016-79-9 MolWeight:414 RetIndex:2930

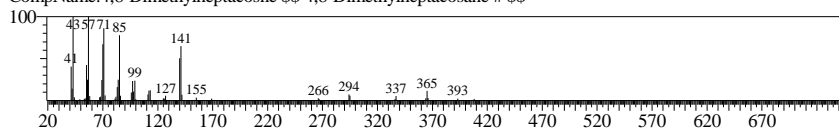
CompName:Heptacosane, 1-chloro- \$\$ 1-Chloroheptacosane # \$\$



Hit#:3 Entry:170192 Library:NIST08.LIB

SI:50 Formula:C29H60 CAS:61295-60-1 MolWeight:408 RetIndex:2776

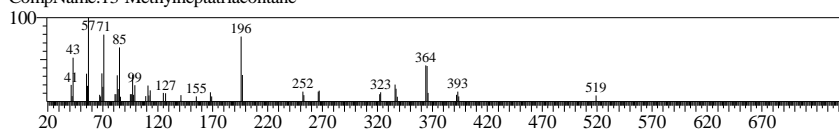
CompName:4,8-Dimethylheptacosane \$\$ 4,8-Dimethylheptacosane # \$\$



Hit#:4 Entry:186716 Library:NIST08.LIB

SI:50 Formula:C38H78 CAS:56987-87-2 MolWeight:534 RetIndex:3734

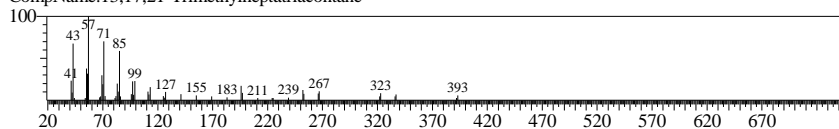
CompName:13-Methylheptatriacontane



Hit#:5 Entry:187922 Library:NIST08.LIB

SI:50 Formula:C40H82 CAS:58668-40-9 MolWeight:562 RetIndex:3805

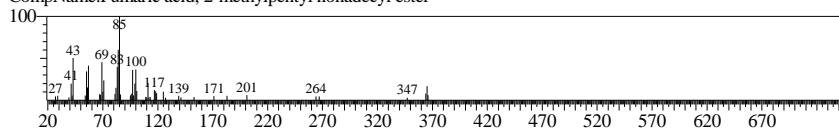
CompName:13,17,21-Trimethylheptatriacontane



Hit#:6 Entry:181206 Library:NIST08.LIB

SI:49 Formula:C29H54O4 CAS:0-00-0 MolWeight:466 RetIndex:3182

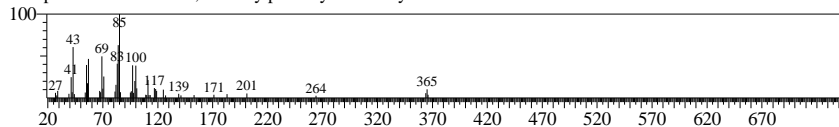
CompName:Fumaric acid, 2-methylpentyl nonadecyl ester



Hit#:7 Entry:181207 Library:NIST08.LIB

SI:49 Formula:C29H54O4 CAS:0-00-0 MolWeight:466 RetIndex:3118

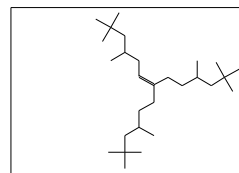
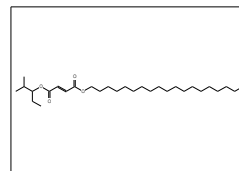
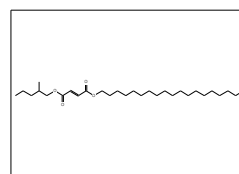
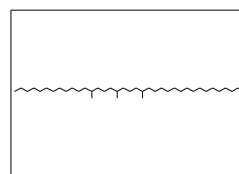
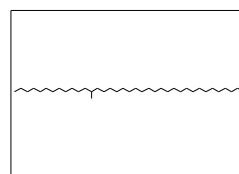
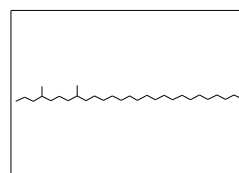
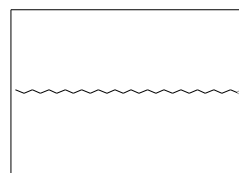
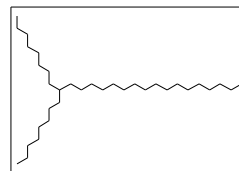
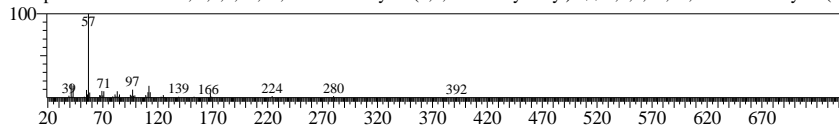
CompName:Fumaric acid, 2-methylpent-3-yl nonadecyl ester



Hit#:8 Entry:165157 Library:NIST08.LIB

SI:48 Formula:C28H56 CAS:55255-73-7 MolWeight:392 RetIndex:2344

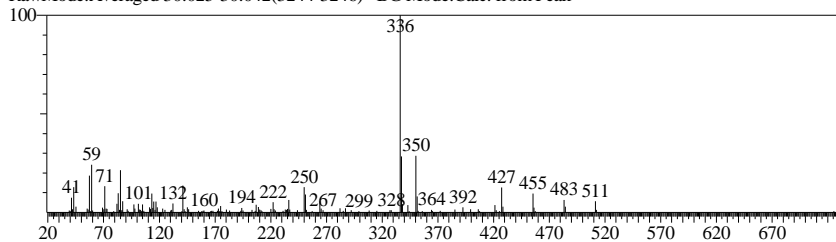
CompName:6-Tridecene, 2,2,4,10,12,12-hexamethyl-7-(3,5,5-trimethylhexyl)- \$\$ 2,2,4,10,12,12-Hexamethyl-7-(3



<< Target >>

Line#:77 R.Time:30.033(Scan#:3245) BasePeak:336.30(72720)

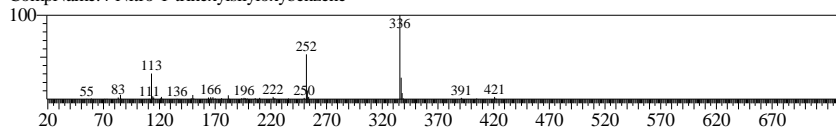
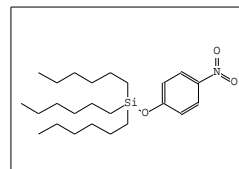
RawMode:Averaged 30.025-30.042(3244-3246) BG Mode:Calc. from Peak



Hit#:1 Entry:173409 Library:NIST08.LIB

SI:51 Formula:C24H43NO3Si CAS:0-00-0 MolWeight:421 RetIndex:2865

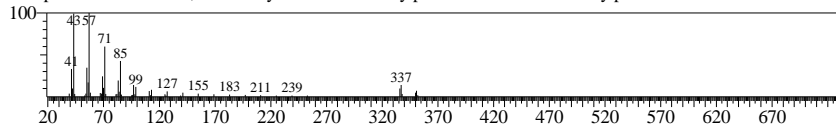
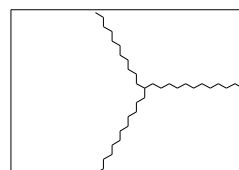
CompName:4-Nitro-1-trihexylsilyloxybenzene



Hit#:2 Entry:185134 Library:NIST08.LIB

SI:48 Formula:C36H74 CAS:55517-89-0 MolWeight:506 RetIndex:3536

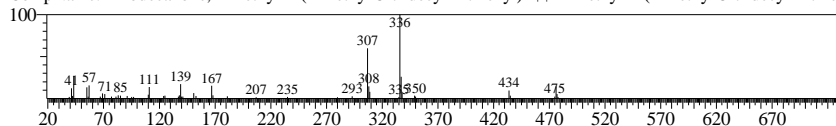
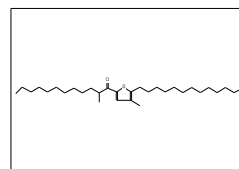
CompName:Pentacosane, 13-undecyl- \$\$ 13-n-Undecylpentacosane \$\$ 13-Undecylpentacosane # \$\$



Hit#:3 Entry:182423 Library:NIST08.LIB

SI:48 Formula:C31H56OS CAS:60089-65-8 MolWeight:476 RetIndex:3431

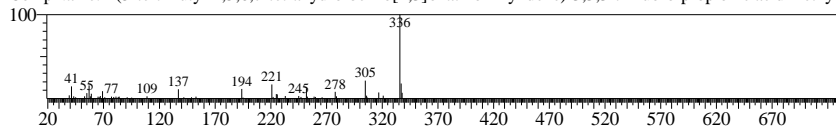
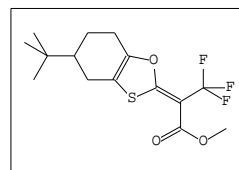
CompName:1-Dodecanone, 2-methyl-1-(4-methyl-5-tridecyl-2-thienyl)- \$\$ 2-Methyl-1-(4-methyl-5-tridecyl-2-thi



Hit#:4 Entry:136374 Library:NIST08.LIB

SI:47 Formula:C15H19F3O3S CAS:0-00-0 MolWeight:336 RetIndex:1787

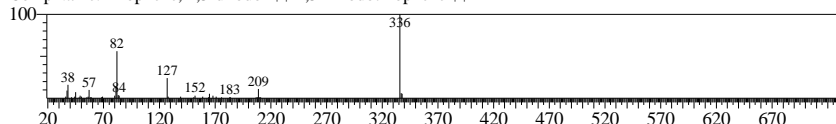
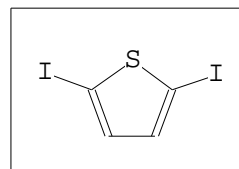
CompName:2-(5-tert-Butyl-4,5,6,7-tetrahydro-benzo[1,3]oxathiol-2-ylidene)-3,3,3-trifluoro-propionic acid methyl



Hit#:5 Entry:136222 Library:NIST08.LIB

SI:46 Formula:C4H2I2S CAS:625-88-7 MolWeight:336 RetIndex:1514

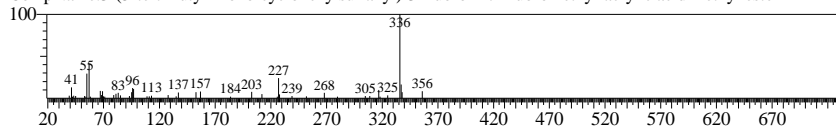
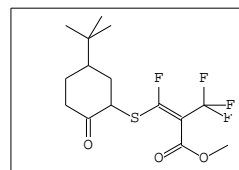
CompName:Thiophene, 2,5-diiodo- \$\$ 2,5-Diiodothiophene \$\$



Hit#:6 Entry:148020 Library:NIST08.LIB

SI:46 Formula:C15H20F4O3S CAS:0-00-0 MolWeight:356 RetIndex:1812

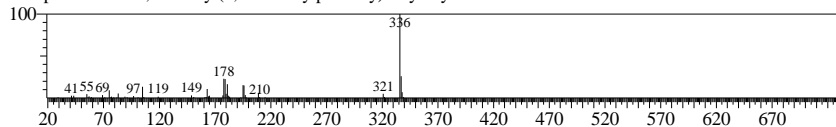
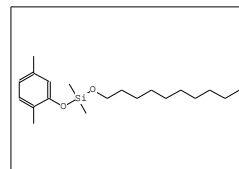
CompName:3-(5-tert-Butyl-2-oxo-cyclohexylsulfanyl)-3-fluoro-2-trifluoromethyl-acrylic acid methyl ester



Hit#:7 Entry:136719 Library:NIST08.LIB

SI:46 Formula:C20H36O2Si CAS:0-00-0 MolWeight:336 RetIndex:2175

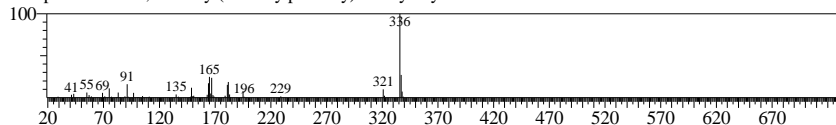
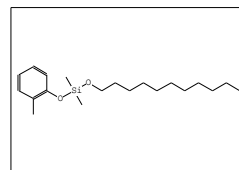
CompName:Silane, dimethyl(2,5-dimethylphenoxy)decyloxy-



Hit#:8 Entry:136718 Library:NIST08.LIB

SI:45 Formula:C20H36O2Si CAS:0-00-0 MolWeight:336 RetIndex:2161

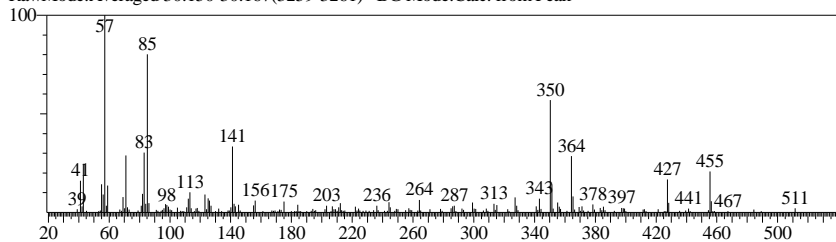
CompName:Silane, dimethyl(2-methylphenoxy)undecyloxy-



<< Target >>

Line#:78 R.Time:30.158(Scan#:3260) BasePeak:57.10(63549)

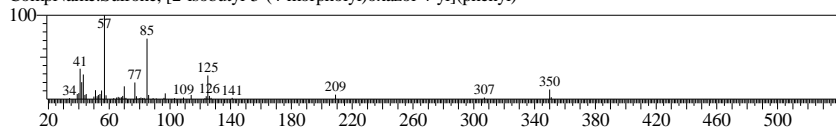
RawMode:Averaged 30.150-30.167(3259-3261) BG Mode:Calc. from Peak



Hit#:1 Entry:144859 Library:NIST08.LIB

SI:58 Formula:C17H22N2O4S CAS:0-00-0 MolWeight:350 RetIndex:2775

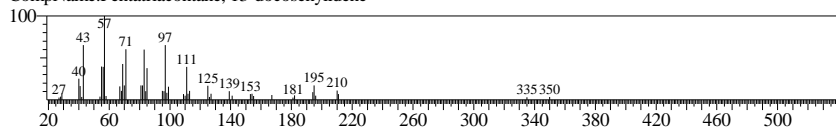
CompName:Sulfone, [2-isobutyl-5-(4-morpholylo)oxazol-4-yl](phenyl)-



Hit#:2 Entry:190082 Library:NIST08.LIB

SI:56 Formula:C47H94 CAS:0-00-0 MolWeight:658 RetIndex:4678

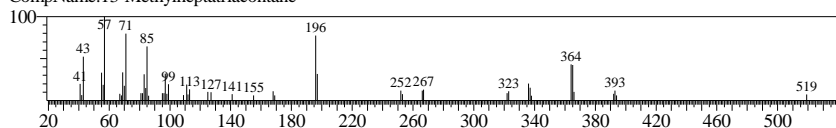
CompName:Pentatriacontane, 13-docosenylidene-



Hit#:3 Entry:186716 Library:NIST08.LIB

SI:55 Formula:C38H78 CAS:56987-87-2 MolWeight:534 RetIndex:3734

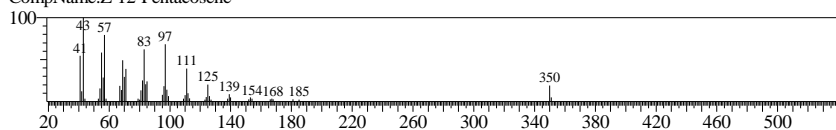
CompName:13-Methylheptatriacontane



Hit#:4 Entry:145194 Library:NIST08.LIB

SI:54 Formula:C25H50 CAS:0-00-0 MolWeight:350 RetIndex:2514

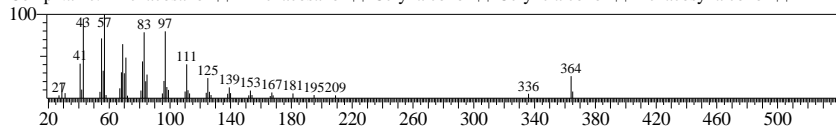
CompName:Z-12-Pentacosene



Hit#:5 Entry:161190 Library:NIST08.LIB

SI:54 Formula:C26H54O CAS:506-52-5 MolWeight:382 RetIndex:2848

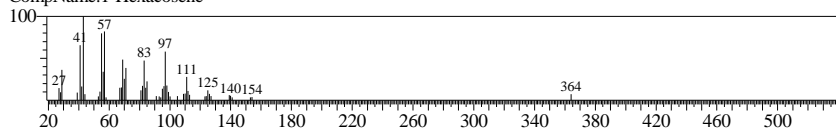
CompName:1-Hexacosanol \$n\$-Hexacosanol \$Ceryl\$ alcohol \$Cerylic\$ alcohol \$Hexacosyl\$ alcohol \$



Hit#:6 Entry:152722 Library:NIST08.LIB

SI:54 Formula:C26H52 CAS:18835-33-1 MolWeight:364 RetIndex:2596

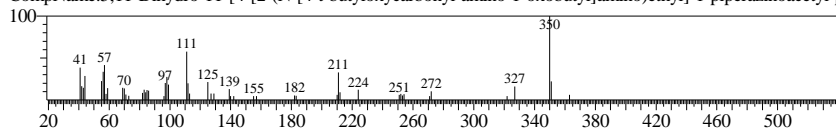
CompName:1-Hexacosene



Hit#:7 Entry:187994 Library:NIST08.LIB

SI:53 Formula:C29H39N7O5 CAS:133727-78-3 MolWeight:565 RetIndex:4739

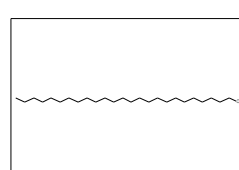
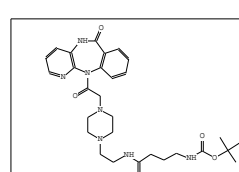
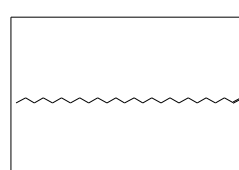
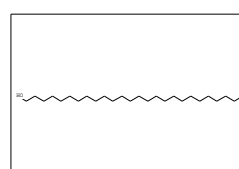
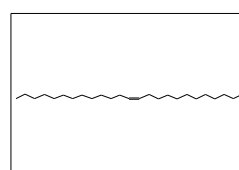
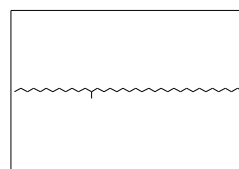
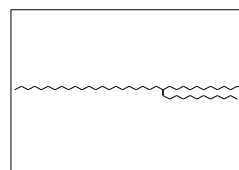
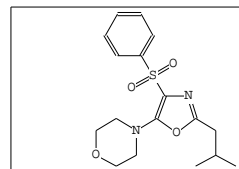
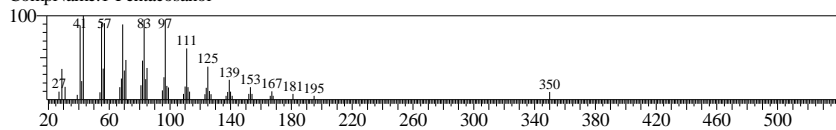
CompName:5,11-Dihydro-11-[4-[2-(N-[4-t-butylloxycarbonyl-amino-1-oxobutyl]amino)ethyl]-1-piperazinoacetyl]-



Hit#:8 Entry:154751 Library:NIST08.LIB

SI:52 Formula:C25H52O CAS:26040-98-2 MolWeight:368 RetIndex:2749

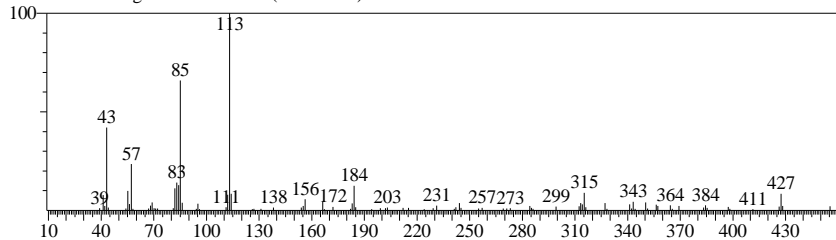
CompName:1-Pentacosanol



<< Target >>

Line#:79 R.Time:30.250(Scan#:3271) BasePeak:113.10(810457)

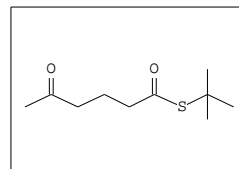
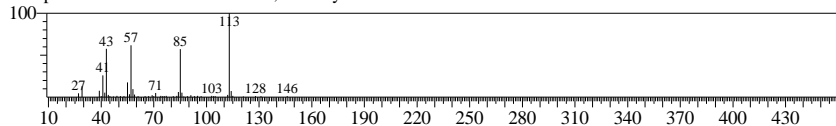
RawMode:Averaged 30.242-30.258(3270-3272) BG Mode:Calc. from Peak



Hit#:1 Entry:43725 Library:NIST08.LIB

SI:72 Formula:C10H18O2S CAS:0-00-0 MolWeight:202 RetIndex:1454

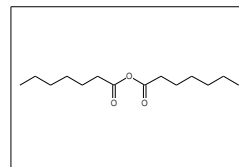
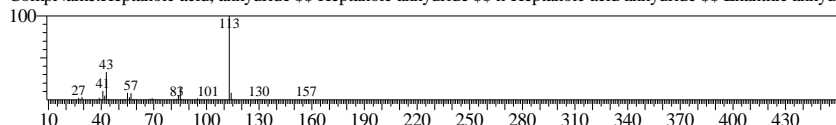
CompName:5-Oxohexanoic acid, S-t-butyl ester



Hit#:2 Entry:70799 Library:NIST08.LIB

SI:71 Formula:C14H26O3 CAS:626-27-7 MolWeight:242 RetIndex:1716

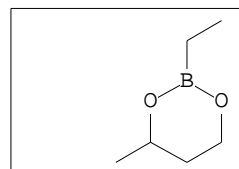
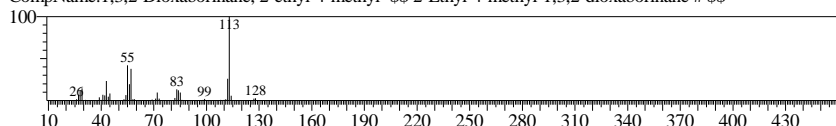
CompName:Heptanoic acid, anhydride \$\$ Heptanoic anhydride \$\$ n-Heptanoic acid anhydride \$\$ Enanthic anhyd



Hit#:3 Entry:7074 Library:NIST08.LIB

SI:69 Formula:C6H13BO2 CAS:57633-65-5 MolWeight:128 RetIndex:0

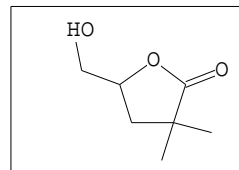
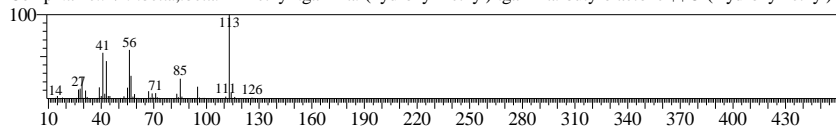
CompName:1,3,2-Dioxaborinane, 2-ethyl-4-methyl- \$\$ 2-Ethyl-4-methyl-1,3,2-dioxaborinane # \$\$



Hit#:4 Entry:12801 Library:NIST08.LIB

SI:68 Formula:C7H12O3 CAS:52398-48-8 MolWeight:144 RetIndex:1263

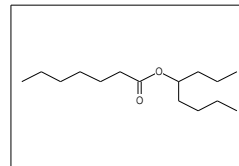
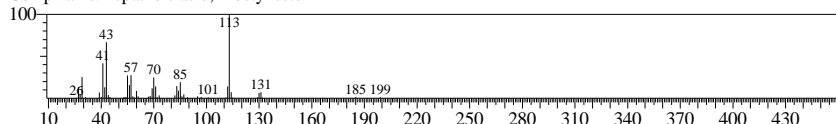
CompName: +/--.beta.,.beta.-Dimethyl-.gamma.-(hydroxy-methyl)-.gamma.-butyrolactone \$\$ 5-(Hydroxymethyl)-



Hit#:5 Entry:70969 Library:NIST08.LIB

SI:68 Formula:C15H30O2 CAS:0-00-0 MolWeight:242 RetIndex:1615

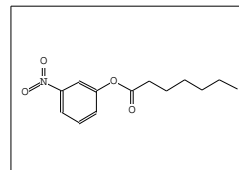
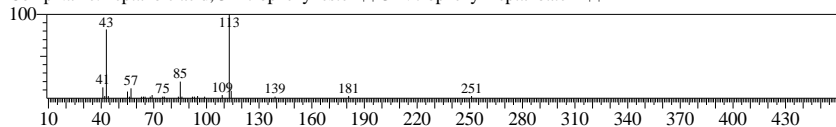
CompName:Heptanoic acid, 4-octyl ester



Hit#:6 Entry:76789 Library:NIST08.LIB

SI:68 Formula:C13H17NO4 CAS:56052-18-7 MolWeight:251 RetIndex:1953

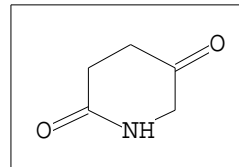
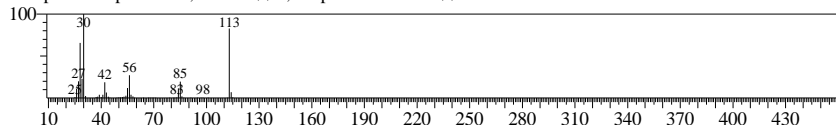
CompName:Heptanoic acid, 3-nitrophenyl ester \$\$ 3-Nitrophenyl heptanoate # \$\$



Hit#:7 Entry:3763 Library:NIST08.LIB

SI:68 Formula:C5H7NO2 CAS:52065-78-8 MolWeight:113 RetIndex:1054

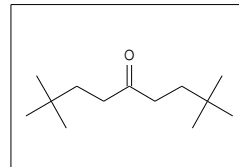
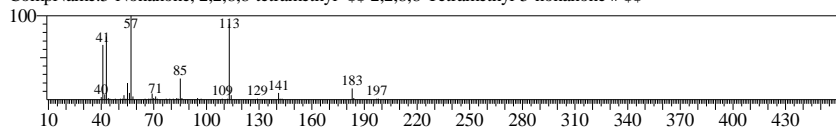
CompName:Piperidine-2,5-dione \$\$ 2,5-Piperidinedione # \$\$



Hit#:8 Entry:41876 Library:NIST08.LIB

SI:67 Formula:C13H26O CAS:5709-95-5 MolWeight:198 RetIndex:1280

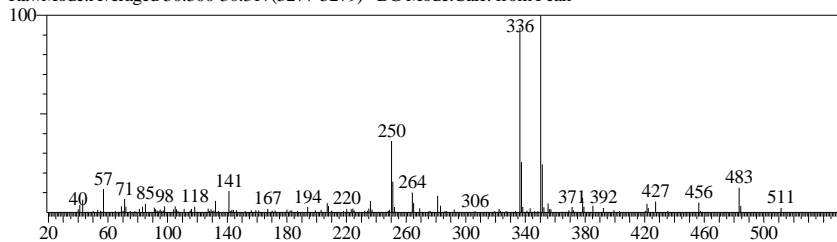
CompName:5-Nonanone, 2,2,8,8-tetramethyl- \$\$ 2,2,8,8-Tetramethyl-5-nonanone # \$\$



<< Target >>

Line#80 R.Time:30.308(Scan#:3278) BasePeak:350.30(32841)

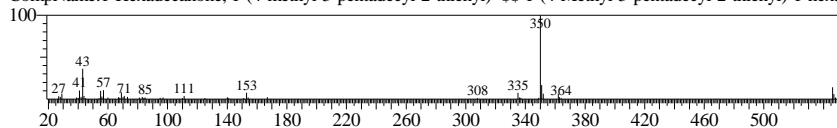
RawMode:Averaged 30.300-30.317(3277-3279) BG Mode:Calc. from Peak



Hit#1 Entry:187267 Library:NIST08.LIB

SI:51 Formula:C36H66OS CAS:59782-69-3 MolWeight:546 RetIndex:3992

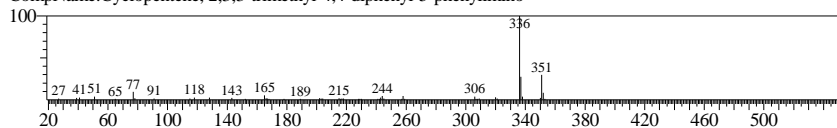
CompName:1-Hexadecanone, 1-(4-methyl-5-pentadecyl-2-thienyl)- \$- 1-(4-Methyl-5-pentadecyl-2-thienyl)-1-hexa



Hit#2 Entry:145644 Library:NIST08.LIB

SI:50 Formula:C26H25N CAS:0-00-0 MolWeight:351 RetIndex:2954

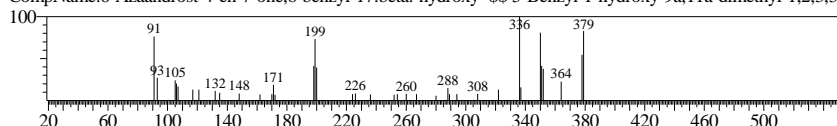
CompName:Cyclopentene, 2,3,3-trimethyl-4,4-diphenyl-5-phenylimino-



Hit#3 Entry:159856 Library:NIST08.LIB

SI:46 Formula:C25H33NO2 CAS:16373-61-8 MolWeight:379 RetIndex:2961

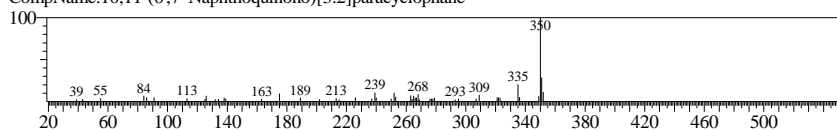
CompName:6-Azaandroster-4-en-7-one,6-benzyl-17.beta.-hydroxy-9a,11a-dimethyl-1,2,3,3-



Hit#4 Entry:145188 Library:NIST08.LIB

SI:46 Formula:C25H18O2 CAS:0-00-0 MolWeight:350 RetIndex:3108

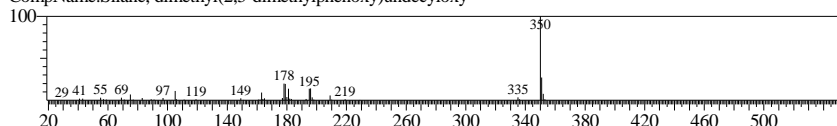
CompName:10,11-(6',7'-Naphthoquinono)[3.2]paracyclophane



Hit#5 Entry:145105 Library:NIST08.LIB

SI:45 Formula:C21H38O2Si CAS:0-00-0 MolWeight:350 RetIndex:2275

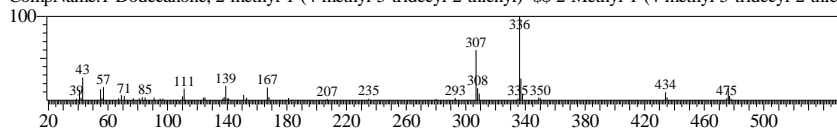
CompName:Silane, dimethyl(2,5-dimethylphenoxy)undecyloxy-



Hit#6 Entry:182423 Library:NIST08.LIB

SI:44 Formula:C31H56OS CAS:60089-65-8 MolWeight:476 RetIndex:3431

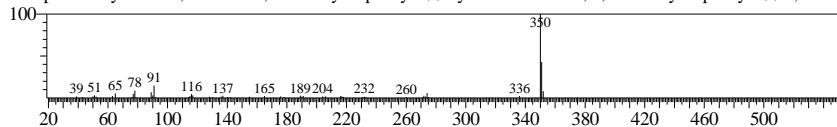
CompName:1-Dodecanone, 2-methyl-1-(4-methyl-5-tridecyl-2-thienyl)- \$- 2-Methyl-1-(4-methyl-5-tridecyl-2-thie



Hit#7 Entry:145631 Library:NIST08.LIB

SI:44 Formula:C24H21N3 CAS:6307-01-3 MolWeight:351 RetIndex:3239

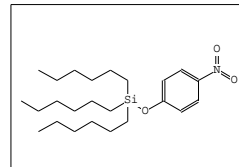
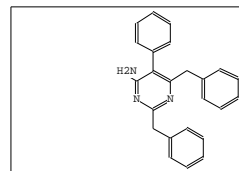
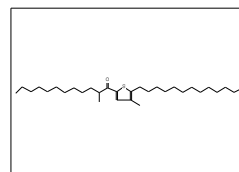
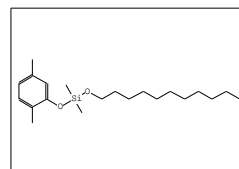
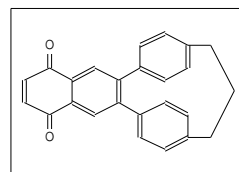
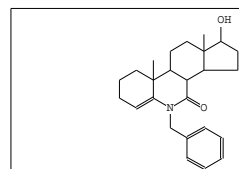
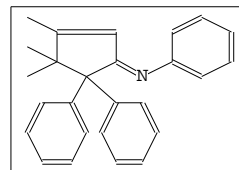
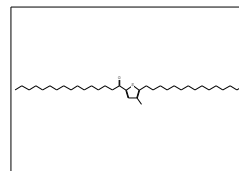
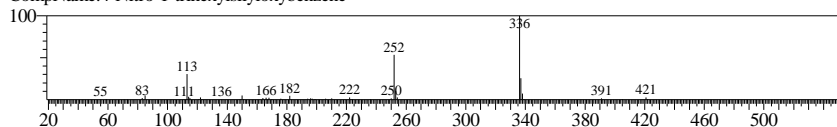
CompName:Pyrimidine, 4-amino-2,6-dibenzyl-5-phenyl- \$- Pyrimidin-4-amine, 2,6-dibenzyl-5-phenyl- \$- 2,6-Dib



Hit#8 Entry:173409 Library:NIST08.LIB

SI:43 Formula:C24H43NO3Si CAS:0-00-0 MolWeight:421 RetIndex:2865

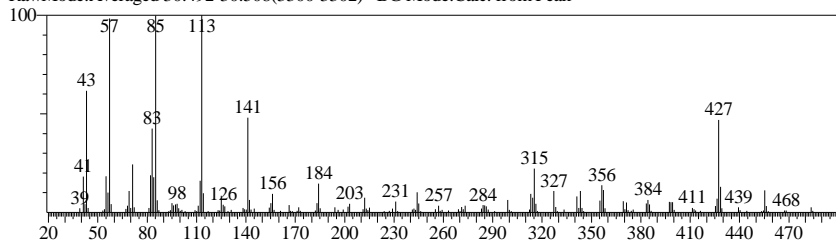
CompName:4-Nitro-1-trihexylsilyloxybenzene



<< Target >>

Line#:81 R.Time:30.500(Scan#:3301) BasePeak:113.10(1447416)

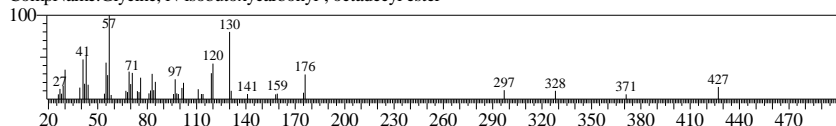
RawMode:Averaged 30.492-30.508(3300-3302) BG Mode:Calc. from Peak



Hit#:1 Entry:174750 Library:NIST08.LIB

SI:49 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975

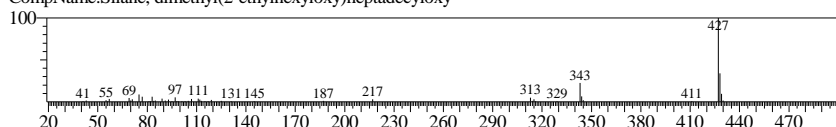
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



Hit#:2 Entry:177733 Library:NIST08.LIB

SI:46 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2703

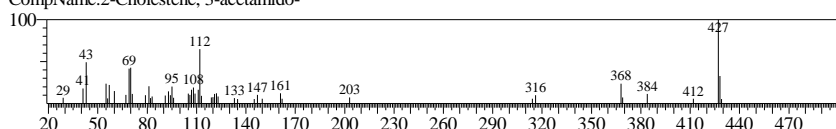
CompName:Silane, dimethyl(2-ethylhexyloxy)heptadecyloxy-



Hit#:3 Entry:174844 Library:NIST08.LIB

SI:44 Formula:C29H49NO CAS:0-00-0 MolWeight:427 RetIndex:2969

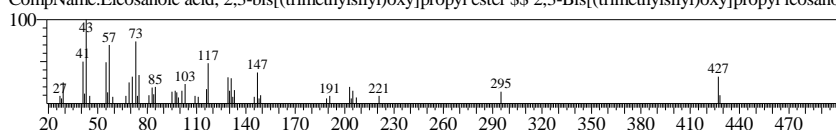
CompName:2-Cholestene, 3-acetamido-



Hit#:4 Entry:186484 Library:NIST08.LIB

SI:41 Formula:C29H62O4Si2 CAS:55517-94-7 MolWeight:530 RetIndex:2979

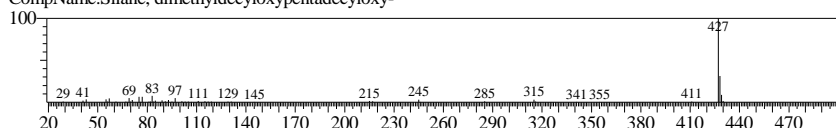
CompName:Eicosanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$\$ 2,3-Bis[(trimethylsilyl)oxy]propyl icosano



Hit#:5 Entry:177735 Library:NIST08.LIB

SI:40 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

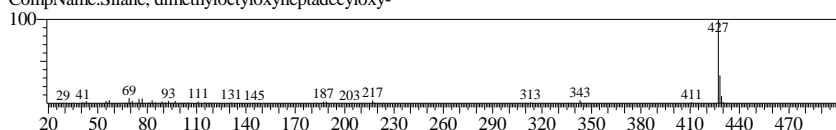
CompName:Silane, dimethyldecyloxy-pentadecyloxy-



Hit#:6 Entry:177734 Library:NIST08.LIB

SI:38 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

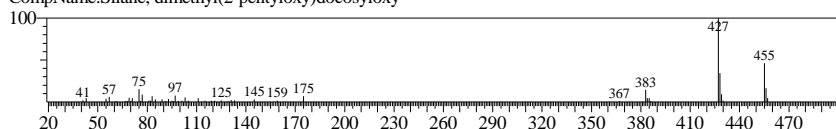
CompName:Silane, dimethyloctyloxyheptadecyloxy-



Hit#:7 Entry:181721 Library:NIST08.LIB

SI:38 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902

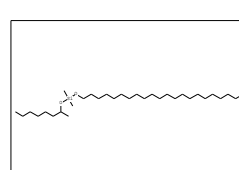
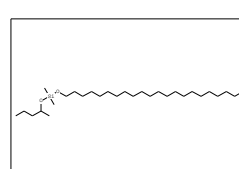
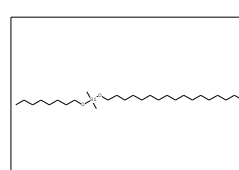
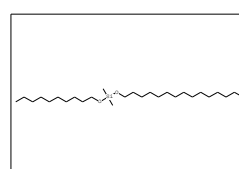
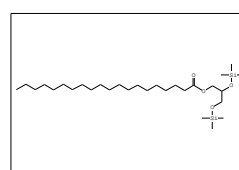
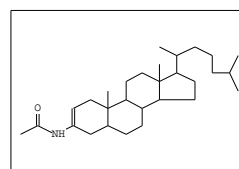
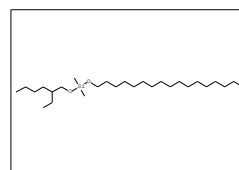
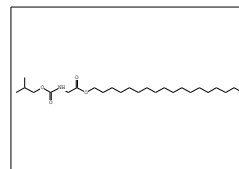
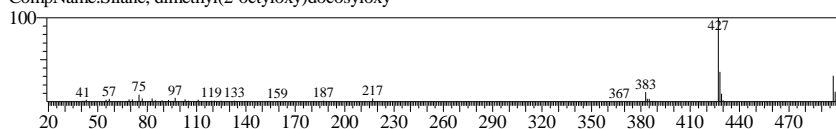
CompName:Silane, dimethyl(2-pentyloxy)docosyloxy-



Hit#:8 Entry:185523 Library:NIST08.LIB

SI:36 Formula:C32H68O2Si CAS:0-00-0 MolWeight:512 RetIndex:3200

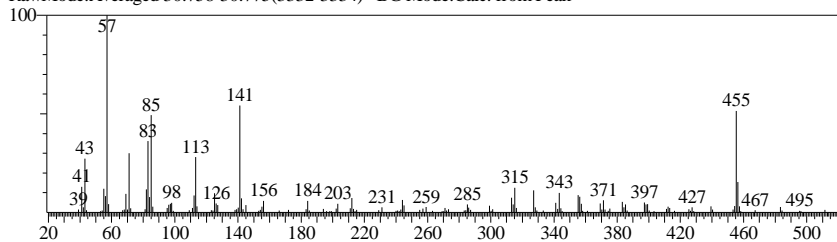
CompName:Silane, dimethyl(2-octyloxy)docosyloxy-



<< Target >>

Line#:82 R.Time:30.767(Scan#:3333) BasePeak:57.10(2852170)

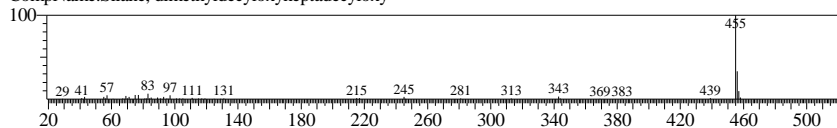
RawMode:Averaged 30.758-30.775(3332-3334) BG Mode:Calc. from Peak



Hit#:1 Entry:181724 Library:NIST08.LIB

SI:47 Formula:C₂₉H₆₂O₂Si CAS:0-00-0 MolWeight:470 RetIndex:2966

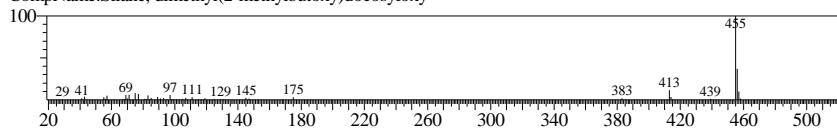
CompName:Silane, dimethyldecyloxyheptadecyloxy-



Hit#:2 Entry:181722 Library:NIST08.LIB

SI:45 Formula:C₂₉H₆₂O₂Si CAS:0-00-0 MolWeight:470 RetIndex:2902

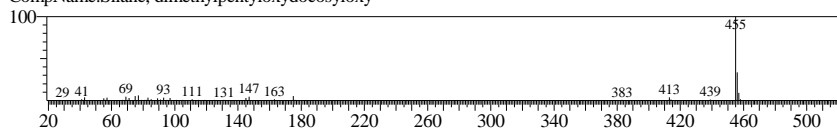
CompName:Silane, dimethyl(2-methylbutoxy)docosyloxy-



Hit#:3 Entry:181723 Library:NIST08.LIB

SI:43 Formula:C₂₉H₆₂O₂Si CAS:0-00-0 MolWeight:470 RetIndex:2966

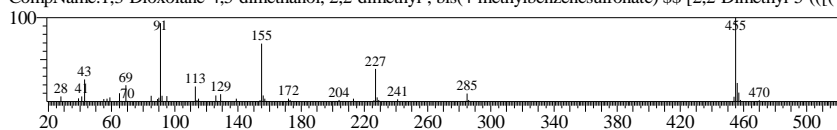
CompName:Silane, dimethylpentyloxydocosyloxy-



Hit#:4 Entry:181657 Library:NIST08.LIB

SI:37 Formula:C₂₁H₂₆O₈S₂ CAS:136881-55-5 MolWeight:470 RetIndex:3572

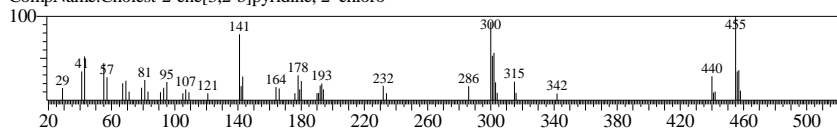
CompName:1,3-Dioxolane-4,5-dimethanol, 2,2-dimethyl-, bis(4-methylbenzenesulfonate) \$\$ [2,2-Dimethyl-5-((1,4



Hit#:5 Entry:179789 Library:NIST08.LIB

SI:36 Formula:C₃₀H₄₆ClN CAS:0-00-0 MolWeight:455 RetIndex:3018

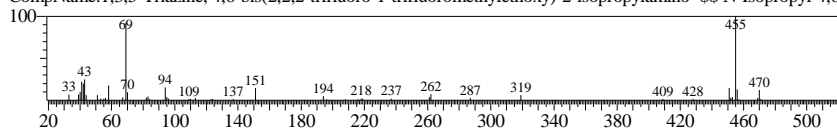
CompName:Cholest-2-ene[3,2-b]pyridine, 2'-chloro-



Hit#:6 Entry:181628 Library:NIST08.LIB

SI:35 Formula:C₁₂H₁₀F₁₂N₄O₂ CAS:304688-10-6 MolWeight:470 RetIndex:1093

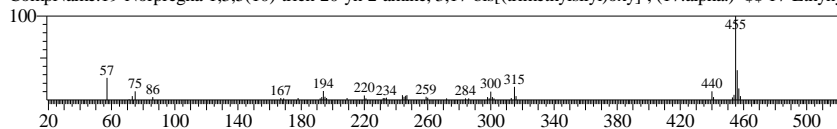
CompName:1,3,5-Triazine, 4,6-bis(2,2,2-trifluoro-1-trifluoromethylethoxy)-2-isopropylamino- \$\$ N-Isopropyl-4,6



Hit#:7 Entry:179708 Library:NIST08.LIB

SI:34 Formula:C₂₆H₄₁NO₂Si₂ CAS:77883-16-0 MolWeight:455 RetIndex:2787

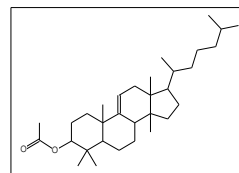
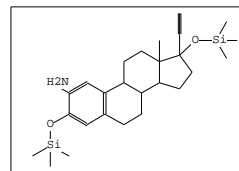
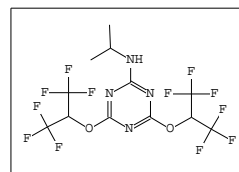
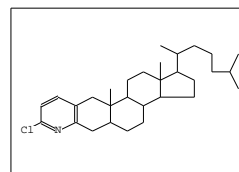
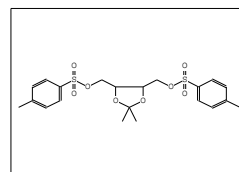
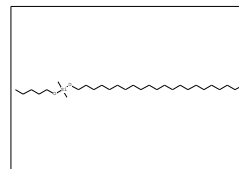
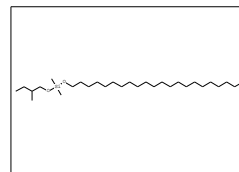
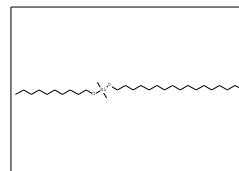
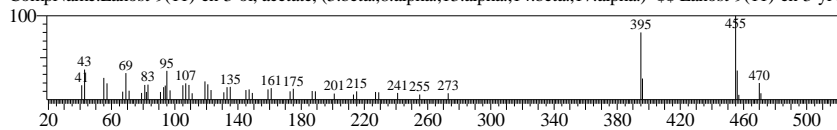
CompName:19-Norpregna-1,3,5(10)-trien-20-yn-2-amine, 3,17-bis[(trimethylsilyl)oxy]-, (17.alpha.)- \$\$ 17-Ethyn



Hit#:8 Entry:181780 Library:NIST08.LIB

SI:34 Formula:C₃₂H₅₄O₂ CAS:55515-26-9 MolWeight:470 RetIndex:2944

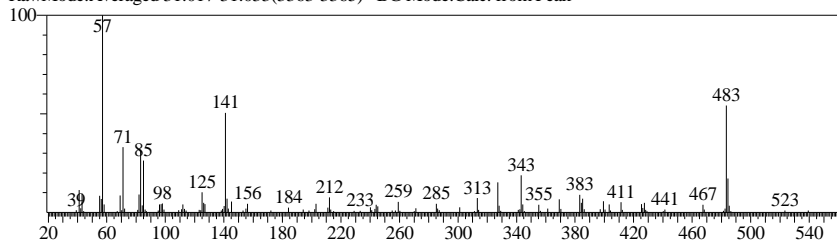
CompName:Lanost-9(11)-en-3-ol, acetate, (3.beta.,8.alpha.,13.alpha.,14.beta.,17.alpha.)- \$\$ Lanost-9(11)-en-3-yl a



<< Target >>

Line#:83 R.Time:31.025(Scan#:3364) BasePeak:57.10(2117098)

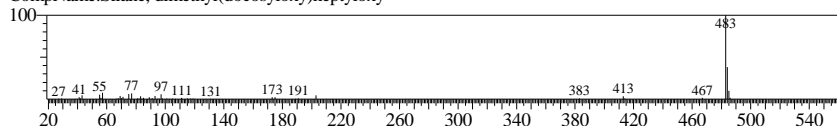
RawMode:Averaged 31.017-31.033(3363-3365) BG Mode:Calc. from Peak



Hit#:1 Entry:184491 Library:NIST08.LIB

SI:49 Formula:C31H66O2Si CAS:0-00-0 MolWeight:498 RetIndex:3164

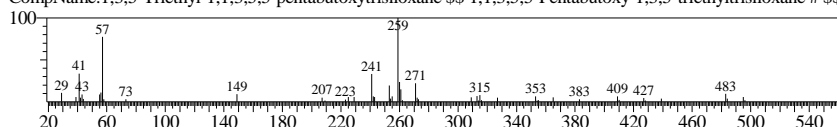
CompName:Silane, dimethyl(docosyloxy)heptyloxy-



Hit#:2 Entry:188092 Library:NIST08.LIB

SI:39 Formula:C26H60O7Si3 CAS:110991-10-1 MolWeight:568 RetIndex:2867

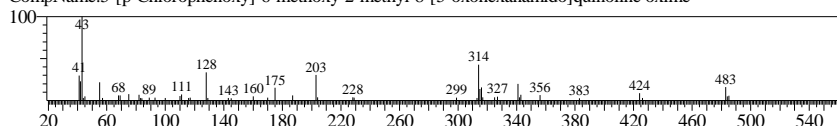
CompName:1,3,5-Triethyl-1,1,3,5,5-pentabutoxytrisiloxane \$\$ 1,1,3,5,5-Pentabutoxy-1,3,5-triethyltrisiloxane # \$\$



Hit#:3 Entry:183114 Library:NIST08.LIB

SI:35 Formula:C25H26ClN3O5 CAS:0-00-0 MolWeight:483 RetIndex:4040

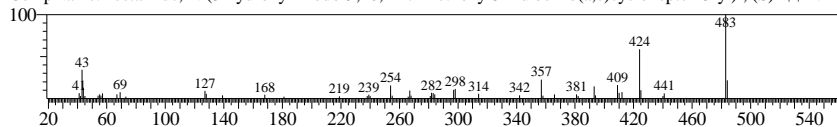
CompName:5-[p-Chlorophenoxyl]-6-methoxy-2-methyl-8-[5-oxohexanamido]quinoline oxime



Hit#:4 Entry:183100 Library:NIST08.LIB

SI:31 Formula:C20H22INO5 CAS:38838-27-6 MolWeight:483 RetIndex:3486

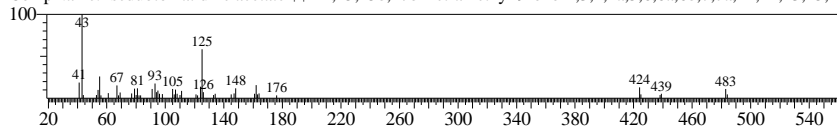
CompName:Acetamide, N-(3-hydroxy-2-iodo-9,10,11-trimethoxy-5H-dibenzo(a,c)cyclohepten-5-yl)-, (S)- \$\$ N-A



Hit#:5 Entry:183154 Library:NIST08.LIB

SI:29 Formula:C30H45NO4 CAS:34638-83-0 MolWeight:483 RetIndex:3391

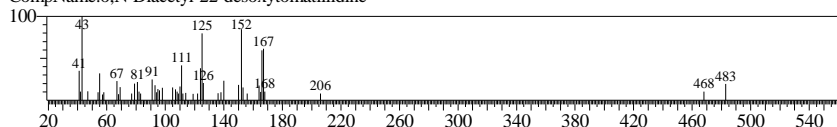
CompName:Pseudotomatidine acetate \$\$ 12,15,15b,17b-Tetramethyl-9-oxo-2,3,4,4a,5,6,6a,6b,7,7a,11,12,13,15,1:



Hit#:6 Entry:183159 Library:NIST08.LIB

SI:28 Formula:C31H49NO3 CAS:0-00-0 MolWeight:483 RetIndex:3253

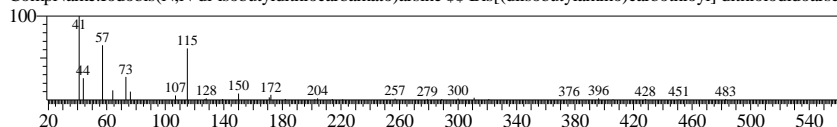
CompName:o,N-Diacetyl-22-desoxytomatillidine



Hit#:7 Entry:189212 Library:NIST08.LIB

SI:27 Formula:C18H36AsIN2S4 CAS:59196-55-3 MolWeight:610 RetIndex:0

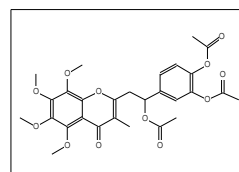
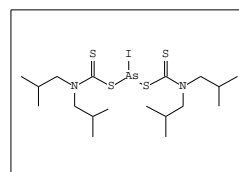
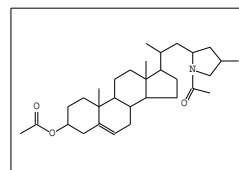
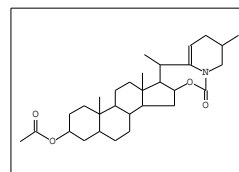
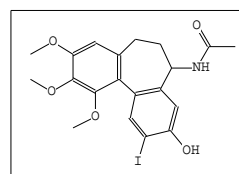
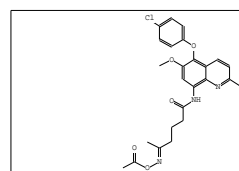
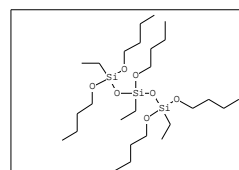
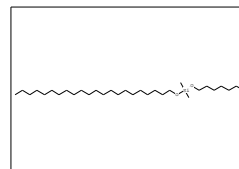
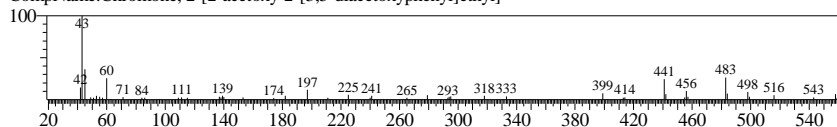
CompName:Iodobis(N,N-di-isobutylidithiocarbamato)arsine \$\$ Bis[(diisobutylamino)carbothioyl] dithioiodidoarse



Hit#:8 Entry:187733 Library:NIST08.LIB

SI:26 Formula:C28H30O12 CAS:0-00-0 MolWeight:558 RetIndex:4079

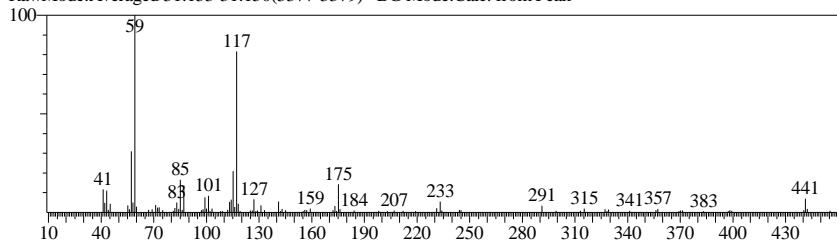
CompName:Chromone, 2-[2-acetoxy-2-[3,5-diacetoxyphenyl]ethyl]-



<< Target >>

Line#:84 R.Time:31.142(Scan#:3378) BasePeak:59.10(376962)

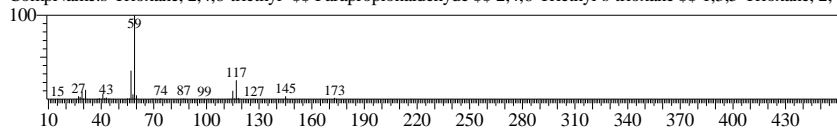
RawMode:Averaged 31.133-31.150(3377-3379) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:74 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

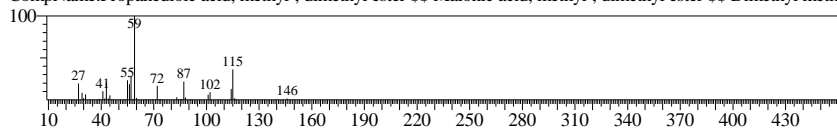
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$ 2,4,6-Triethyl-s-trioxane \$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:13447 Library:NIST08.LIB

SI:72 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

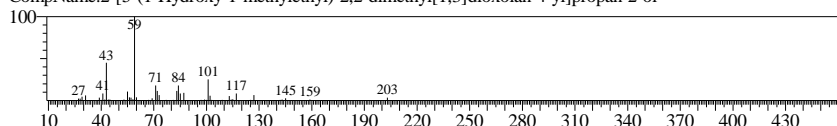
CompName:Propanedioic acid, methyl-, dimethyl ester \$ Malonic acid, methyl-, dimethyl ester \$ Dimethyl meth



Hit#:3 Entry:54251 Library:NIST08.LIB

SI:71 Formula:C11H22O4 CAS:0-00-0 MolWeight:218 RetIndex:1415

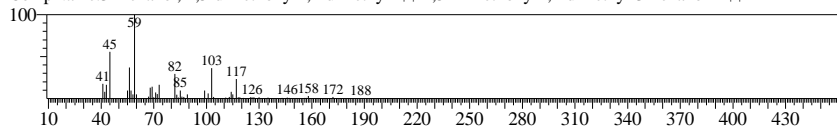
CompName:2-[5-(1-Hydroxy-1-methylethyl)-2,2-dimethyl[1,3]dioxolan-4-yl]propan-2-ol



Hit#:4 Entry:36580 Library:NIST08.LIB

SI:70 Formula:C10H22O3 CAS:13897-22-8 MolWeight:190 RetIndex:1137

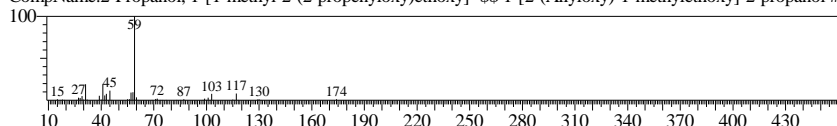
CompName:3-Hexanol, 1,5-dimethoxy-2,4-dimethyl- \$ 1,5-Dimethoxy-2,4-dimethyl-3-hexanol # \$ \$



Hit#:5 Entry:27550 Library:NIST08.LIB

SI:70 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

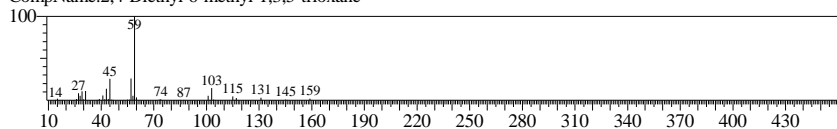
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:6 Entry:19968 Library:NIST08.LIB

SI:70 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

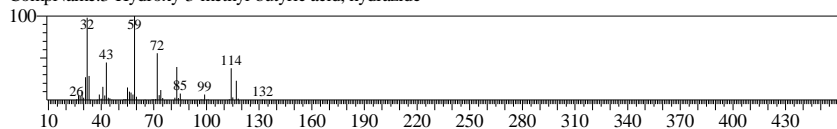
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#:7 Entry:8329 Library:NIST08.LIB

SI:69 Formula:C5H12N2O2 CAS:0-00-0 MolWeight:132 RetIndex:1253

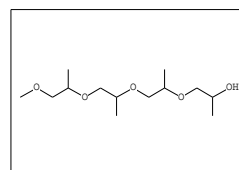
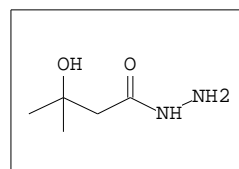
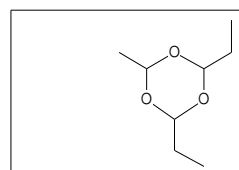
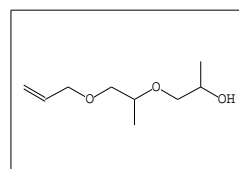
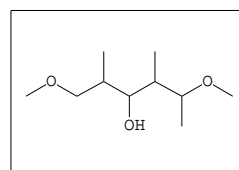
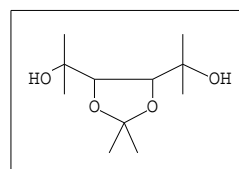
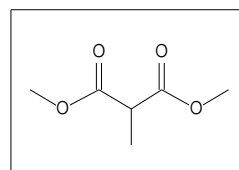
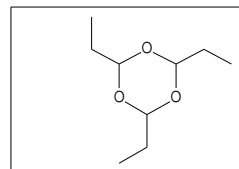
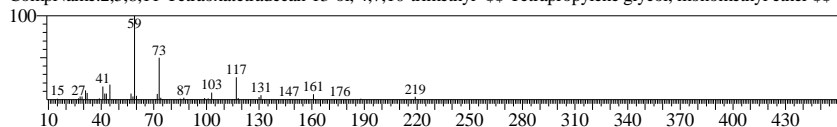
CompName:3-Hydroxy-3-methyl-butyric acid, hydrazide



Hit#:8 Entry:85908 Library:NIST08.LIB

SI:69 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

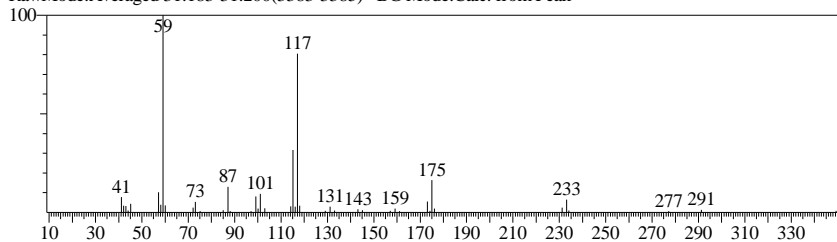
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$ Tetrapropylene glycol, monomethyl ether \$



<< Target >>

Line#85 R.Time:31.192(Scan#:3384) BasePeak:59.10(339640)

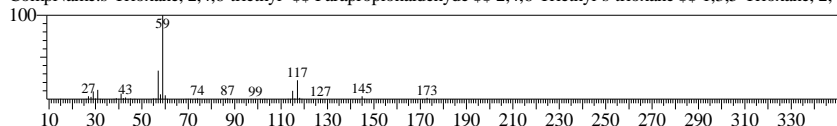
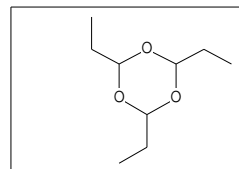
RawMode:Averaged 31.183-31.200(3383-3385) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:76 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

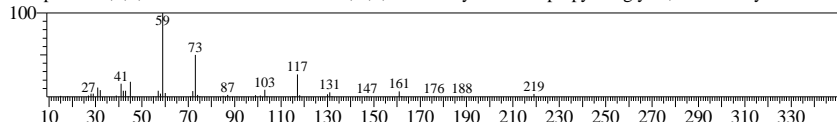
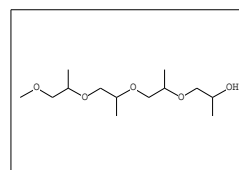
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:74 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

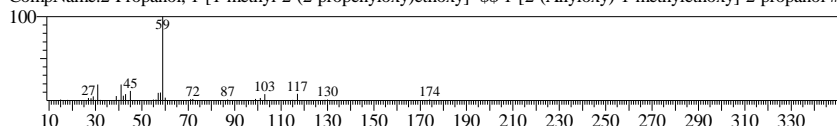
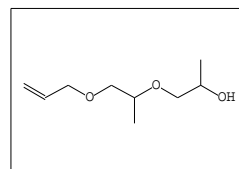
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#3 Entry:27550 Library:NIST08.LIB

SI:73 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

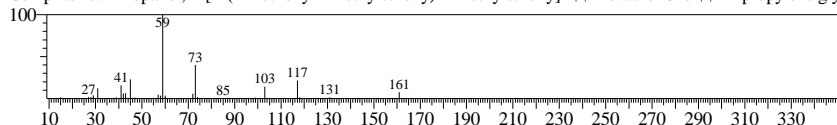
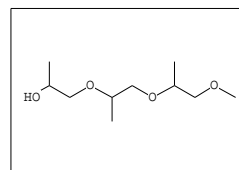
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:17307 Library:NIST08s.LIB

SI:73 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

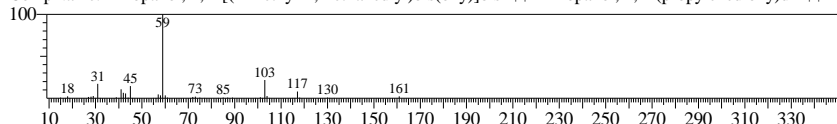
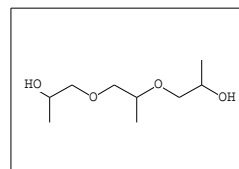
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#5 Entry:37499 Library:NIST08.LIB

SI:72 Formula:C9H20O4 CAS:1638-16-0 MolWeight:192 RetIndex:1328

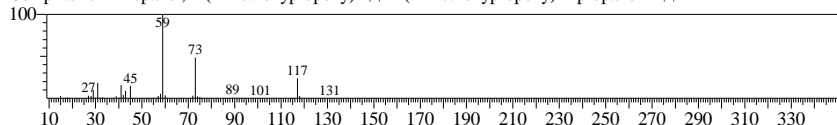
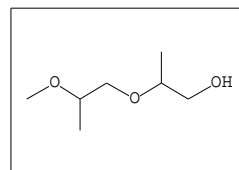
CompName:2-Propanol, 1,1'-[(1-methyl-1,2-ethanediyl)bis(oxy)]bis- \$\$ 2-Propanol, 1,1'-(propylenedioxy)di- \$\$ T



Hit#6 Entry:14297 Library:NIST08.LIB

SI:72 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

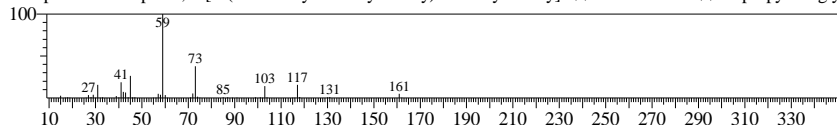
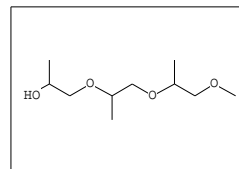
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#7 Entry:46317 Library:NIST08.LIB

SI:72 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

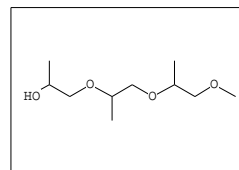
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#8 Entry:17306 Library:NIST08s.LIB

SI:72 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

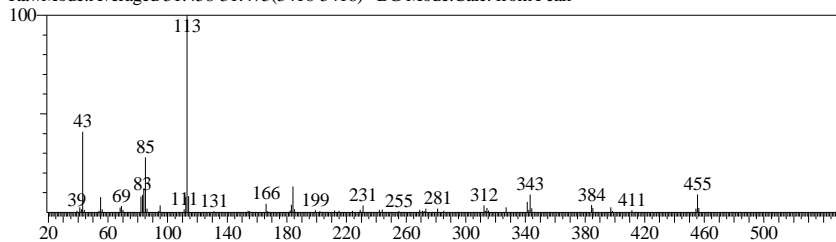
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



<< Target >>

Line#:86 R.Time:31.467(Scan#:3417) BasePeak:113.10(188825)

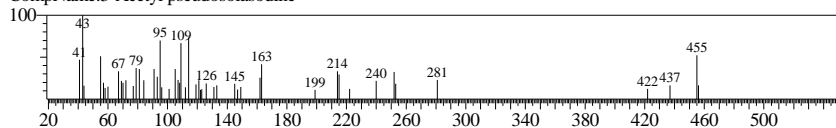
RawMode:Averaged 31.458-31.475(3416-3418) BG Mode:Calc. from Peak



Hit#:1 Entry:179782 Library:NIST08.LIB

SI:26 Formula:C₂₉H₄₅NO₃ CAS:0-00-0 MolWeight:455 RetIndex:3161

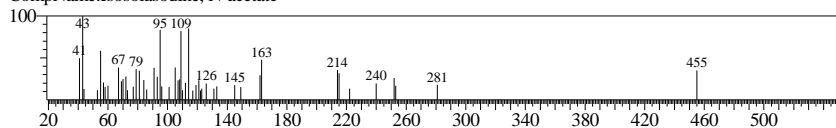
CompName:3-Acetyl pseudosolasodine



Hit#:2 Entry:179781 Library:NIST08.LIB

SI:25 Formula:C₂₉H₄₅NO₃ CAS:0-00-0 MolWeight:455 RetIndex:3312

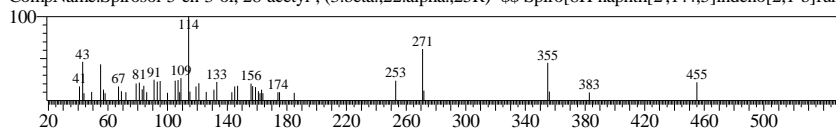
CompName:Isosolasodine, N-acetate



Hit#:3 Entry:179785 Library:NIST08.LIB

SI:24 Formula:C₂₉H₄₅NO₃ CAS:58822-29-0 MolWeight:455 RetIndex:3203

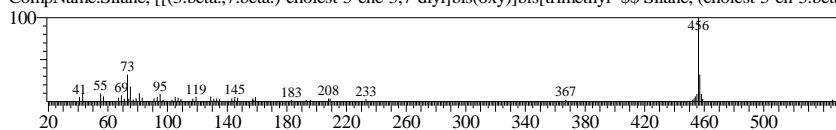
CompName:Spinosol-5-en-3-ol, 28-acetyl-, (3.beta.,22.alpha.,25R)- \$\$ Spiro[8H-naphth[2',1':4,5]indeno[2,1-b]fur:



Hit#:4 Entry:28125 Library:NIST08s.LIB

SI:23 Formula:C₃₃H₆₂O₂Si₂ CAS:33287-26-2 MolWeight:546 RetIndex:2900

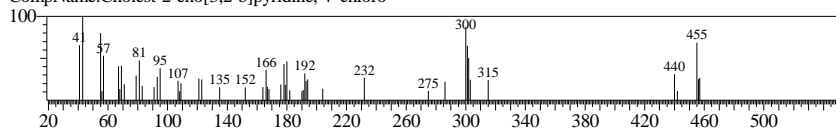
CompName:Silane, [[(3.beta.,7.beta.)-cholest-5-ene-3,7-diyl]bis(oxy)]bis[trimethyl-



Hit#:5 Entry:179788 Library:NIST08.LIB

SI:21 Formula:C₃₀H₄₆ClN CAS:0-00-0 MolWeight:455 RetIndex:3018

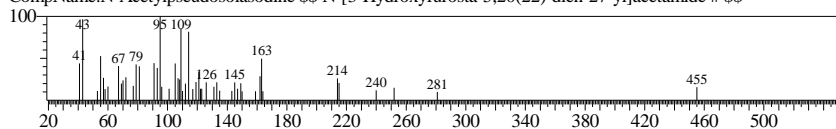
CompName:Cholest-2-eno[3,2-b]pyridine, 4'-chloro-



Hit#:6 Entry:179783 Library:NIST08.LIB

SI:21 Formula:C₂₉H₄₅NO₃ CAS:24261-59-4 MolWeight:455 RetIndex:3312

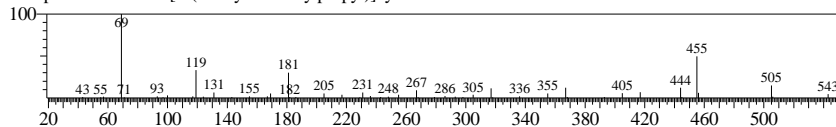
CompName:N-Acetyl pseudosolasodine \$\$ N-[3-Hydroxyfurosta-5,20(22)-dien-27-yl]acetamide # \$\$



Hit#:7 Entry:187858 Library:NIST08.LIB

SI:20 Formula:C₁₂F₂₂ CAS:0-00-0 MolWeight:562 RetIndex:0

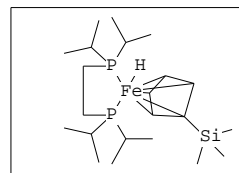
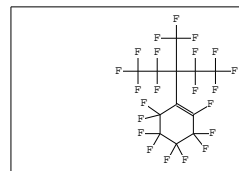
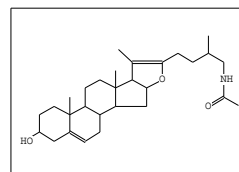
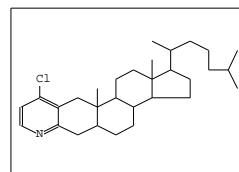
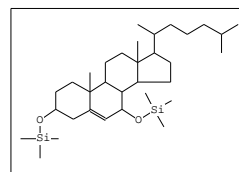
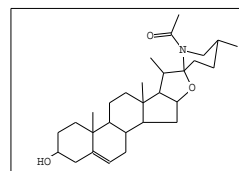
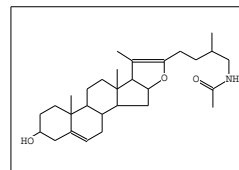
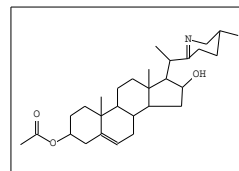
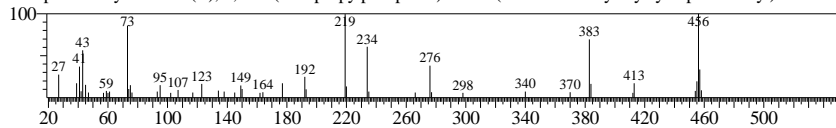
CompName:Perfluoro[1-(1-ethyl-1-methylpropyl)]cyclohexene



Hit#:8 Entry:179843 Library:NIST08.LIB

SI:18 Formula:C₂₂H₄₆FeP₂Si CAS:0-00-0 MolWeight:456 RetIndex:0

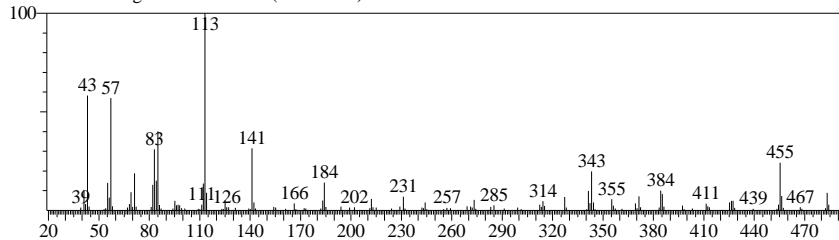
CompName:Hydridoiron(II), 1,2-bis(diisopropylphosphino)ethane-(eta.-5-trimethylsilylcyclopentadienyl)-



<< Target >>

Line#:87 R.Time:31.775(Scan#:3454) BasePeak:113.10(807600)

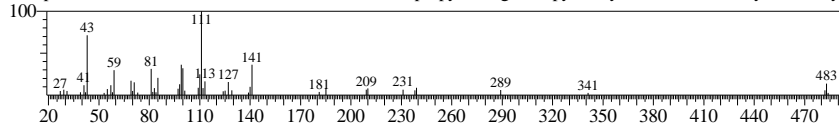
RawMode:Averaged 31.767-31.783(3453-3455) BG Mode:Calc. from Peak



Hit#:1 Entry:184448 Library:NIST08.LIB

SI:50 Formula:C22H36B2O11 CAS:93524-14-2 MolWeight:498 RetIndex:0

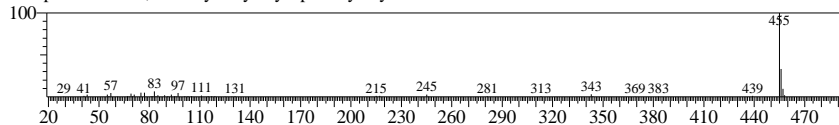
CompName:.beta.-d-Mannofuranoside, 1,2-3,4-di-O-isopropylidene-galactopyranosyl-2,3-5,6-bis-O-ethylboranedi-



Hit#:2 Entry:181724 Library:NIST08.LIB

SI:39 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966

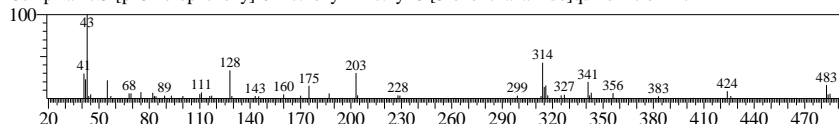
CompName:Silane, dimethyldecyloxyheptadecyloxy-



Hit#:3 Entry:183114 Library:NIST08.LIB

SI:38 Formula:C25H26ClN3O5 CAS:0-00-0 MolWeight:483 RetIndex:4040

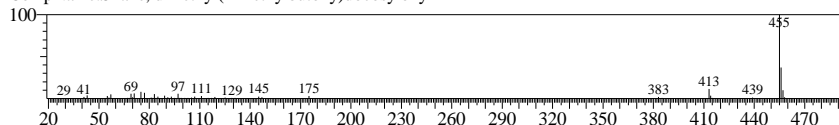
CompName:5-[p-Chlorophenoxyl]-6-methoxy-2-methyl-8-[5-oxohexanamido]quinoline oxime



Hit#:4 Entry:181722 Library:NIST08.LIB

SI:37 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902

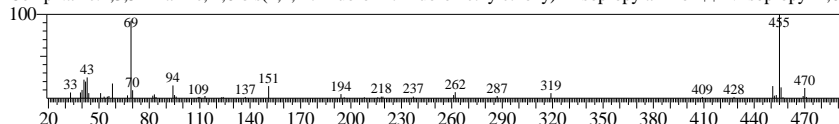
CompName:Silane, dimethyl(2-methylbutoxy)docosyloxy-



Hit#:5 Entry:181628 Library:NIST08.LIB

SI:36 Formula:C12H10F12N4O2 CAS:304688-10-6 MolWeight:470 RetIndex:1093

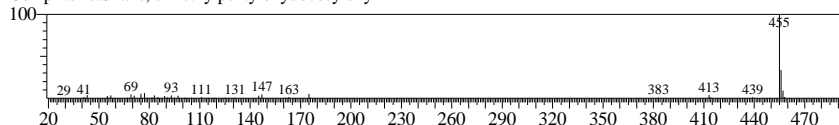
CompName:1,3,5-Triazine, 4,6-bis(2,2,2-trifluoro-1-trifluoromethylethoxy)-2-isopropylamino- \$N\$-Isopropyl-4,6



Hit#:6 Entry:181723 Library:NIST08.LIB

SI:35 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966

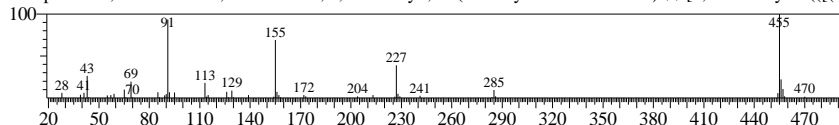
CompName:Silane, dimethylpentyloxydocosyloxy-



Hit#:7 Entry:181657 Library:NIST08.LIB

SI:33 Formula:C21H26O8S2 CAS:136881-55-5 MolWeight:470 RetIndex:3572

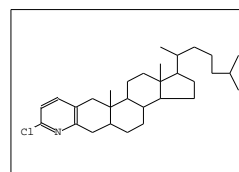
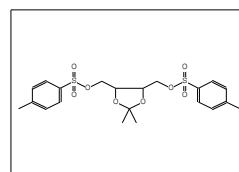
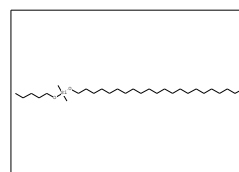
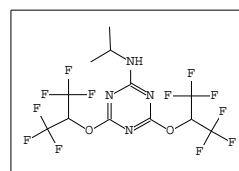
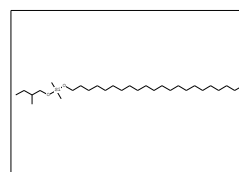
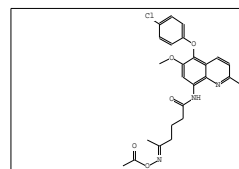
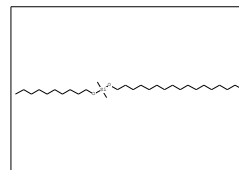
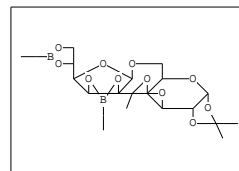
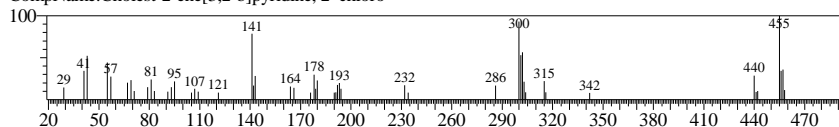
CompName:1,3-Dioxolane-4,5-dimethanol, 2,2-dimethyl-, bis(4-methylbenzenesulfonate) \$[2,2\$-Dimethyl-5-((1,4



Hit#:8 Entry:179789 Library:NIST08.LIB

SI:32 Formula:C30H46ClN CAS:0-00-0 MolWeight:455 RetIndex:3018

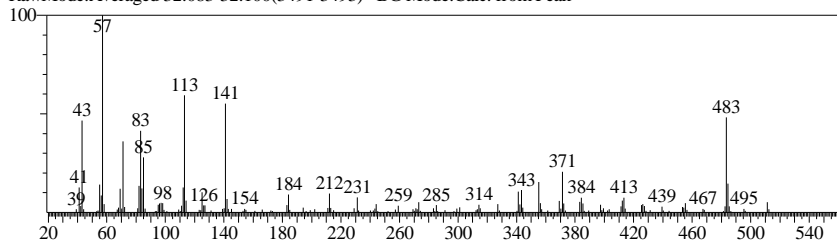
CompName:Cholest-2-ene[3,2-b]pyridine, 2'-chloro-



<< Target >>

Line#:88 R.Time:32.092(Scan#:3492) BasePeak:57.10(1430188)

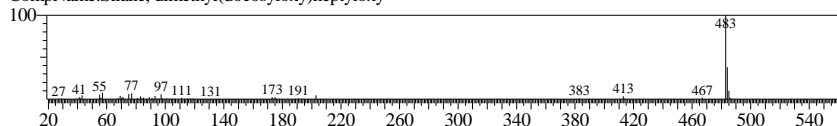
RawMode:Averaged 32.083-32.100(3491-3493) BG Mode:Calc. from Peak



Hit#:1 Entry:184491 Library:NIST08.LIB

SI:42 Formula:C31H66O2Si CAS:0-00-0 MolWeight:498 RetIndex:3164

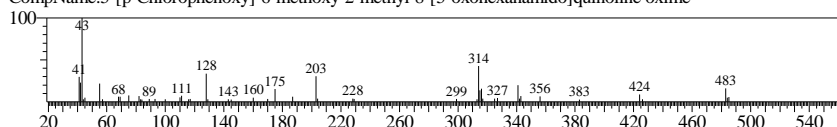
CompName:Silane, dimethyl(docosyloxy)heptyloxy-



Hit#:2 Entry:183114 Library:NIST08.LIB

SI:38 Formula:C25H26ClN3O5 CAS:0-00-0 MolWeight:483 RetIndex:4040

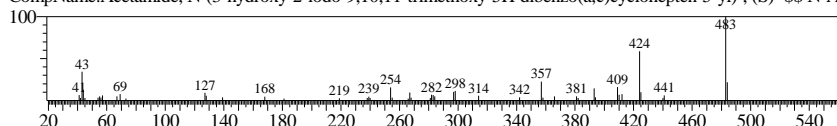
CompName:5-[p-Chlorophenoxyl]-6-methoxy-2-methyl-8-[5-oxohexanamido]quinoline oxime



Hit#:3 Entry:183100 Library:NIST08.LIB

SI:31 Formula:C20H22INO5 CAS:38838-27-6 MolWeight:483 RetIndex:3486

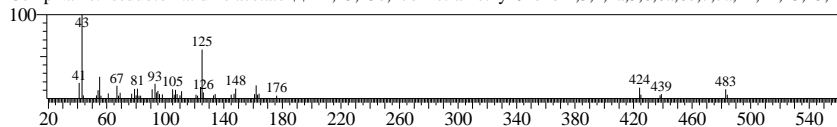
CompName:Acetamide, N-(3-hydroxy-2-iodo-9,10,11-trimethoxy-5H-dibenzo(a,c)cyclohepten-5-yl)-, (S)-



Hit#:4 Entry:183154 Library:NIST08.LIB

SI:30 Formula:C30H45NO4 CAS:34638-83-0 MolWeight:483 RetIndex:3391

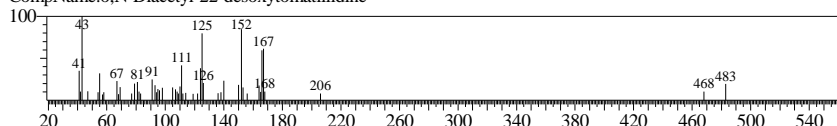
CompName:Pseudoatomitidine acetate



Hit#:5 Entry:183159 Library:NIST08.LIB

SI:29 Formula:C31H49NO3 CAS:0-00-0 MolWeight:483 RetIndex:3253

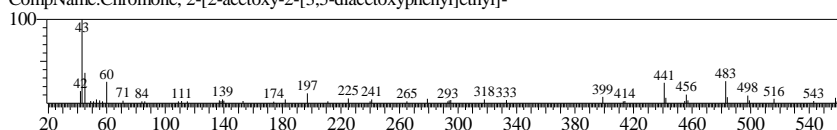
CompName:o,N-Diacetyl-22-desoxytomatillidine



Hit#:6 Entry:187733 Library:NIST08.LIB

SI:28 Formula:C28H30O12 CAS:0-00-0 MolWeight:558 RetIndex:4079

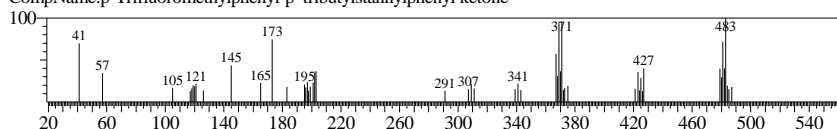
CompName:Chromone, 2-[2-acetoxy-2-[3,5-diacetoxyphenyl]ethyl]-



Hit#:7 Entry:186971 Library:NIST08.LIB

SI:24 Formula:C26H35F3OSn CAS:0-00-0 MolWeight:540 RetIndex:0

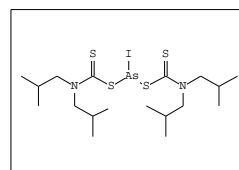
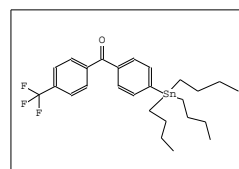
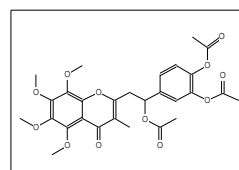
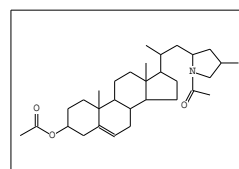
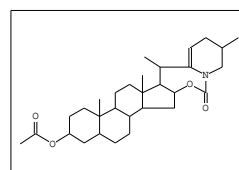
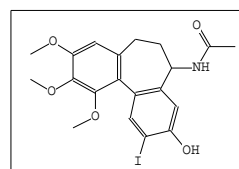
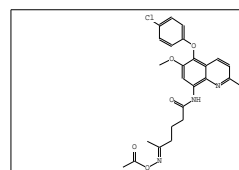
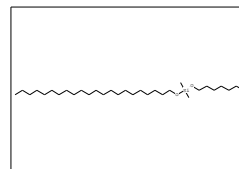
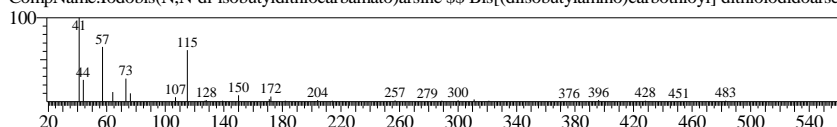
CompName:p-Trifluoromethylphenyl-p'-tributylstannylphenyl ketone



Hit#:8 Entry:189212 Library:NIST08.LIB

SI:22 Formula:C18H36AsIN2S4 CAS:59196-55-3 MolWeight:610 RetIndex:0

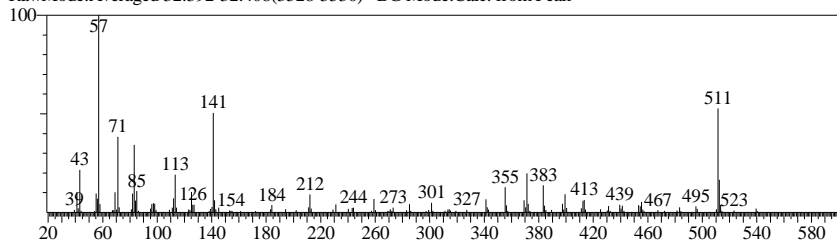
CompName:Iodobis(N,N-di-isobutylthiocarbamato)arsine



<< Target >>

Line#:89 R.Time:32.400(Scan#:3529) BasePeak:57.10(1911178)

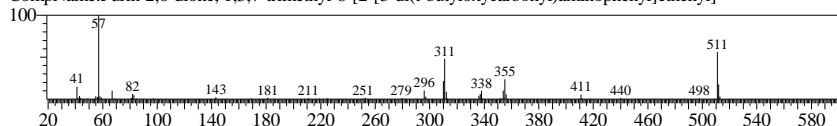
RawMode:Averaged 32.392-32.408(3528-3530) BG Mode:Calc. from Peak



Hit#:1 Entry:185431 Library:NIST08.LIB

SI:50 Formula:C26H33N5O6 CAS:0-00-0 MolWeight:511 RetIndex:3851

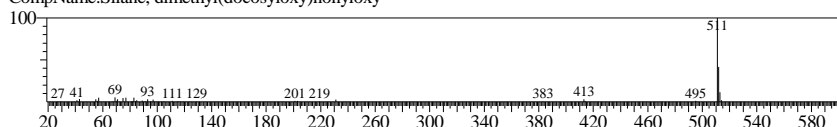
CompName:Purin-2,6-dione, 1,3,7-trimethyl-8-[2-[3-di(t-butyloxycarbonyl)aminophenyl]ethenyl]-



Hit#:2 Entry:186328 Library:NIST08.LIB

SI:49 Formula:C33H70O2Si CAS:0-00-0 MolWeight:526 RetIndex:3363

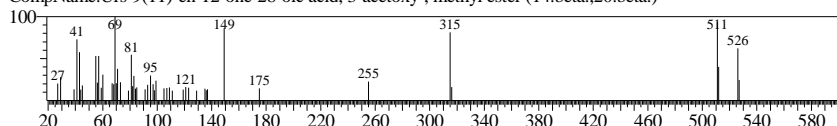
CompName:Silane, dimethyl(docosyloxy)nonyloxy-



Hit#:3 Entry:186325 Library:NIST08.LIB

SI:38 Formula:C33H50O5 CAS:0-00-0 MolWeight:526 RetIndex:3451

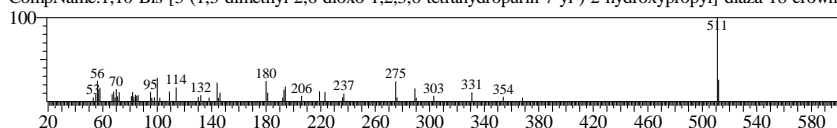
CompName:Urs-9(11)-en-12-one-28-oic acid, 3-acetoxy-, methyl ester (14.beta.,20.beta.)



Hit#:4 Entry:190782 Library:NIST08.LIB

SI:34 Formula:C32H50N10O10 CAS:171524-70-2 MolWeight:734 RetIndex:6411

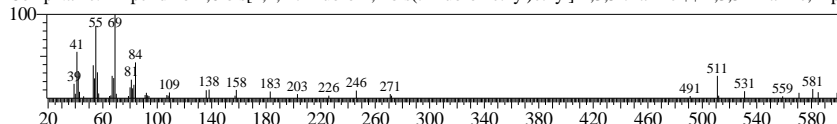
CompName:1,10-Bis-[3-(1,3-dimethyl-2,6-dioxo-1,2,3,6-tetrahydropurin-7-yl)-2-hydroxypropyl]-diazia-18-crown-



Hit#:5 Entry:189009 Library:NIST08.LIB

SI:33 Formula:C16H10F18N4 CAS:254118-24-6 MolWeight:600 RetIndex:1047

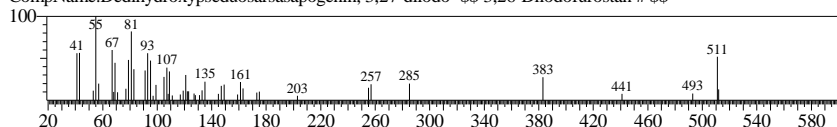
CompName:2-Piperidino-4,6-bis[2,2,2-trifluoro-1,1-bis(trifluoromethyl)ethyl]-1,3,5-triazine



Hit#:6 Entry:189748 Library:NIST08.LIB

SI:33 Formula:C27H44I2O CAS:111824-22-7 MolWeight:638 RetIndex:3266

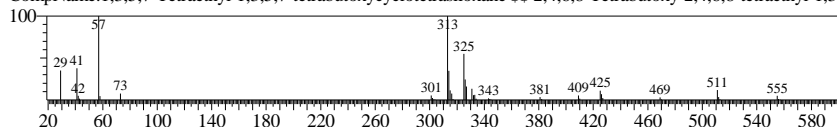
CompName:Dedihydroxypseduosarsapogenin, 3,27-diiodo- 3,26-Diiodofurostan



Hit#:7 Entry:188586 Library:NIST08.LIB

SI:30 Formula:C24H56O8Si4 CAS:110991-13-4 MolWeight:584 RetIndex:2721

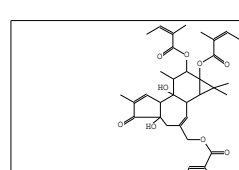
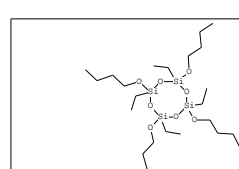
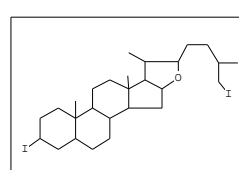
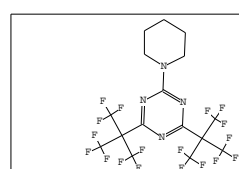
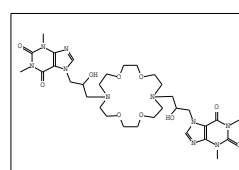
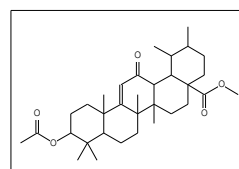
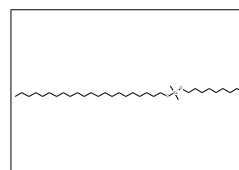
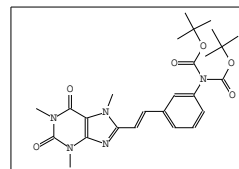
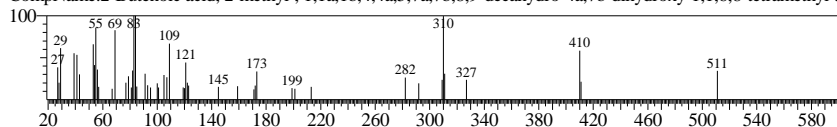
CompName:1,3,5,7-Tetraethyl-1,3,5,7-tetrabutoxycyclotetrasiloxane



Hit#:8 Entry:189229 Library:NIST08.LIB

SI:28 Formula:C35H46O9 CAS:77573-12-7 MolWeight:610 RetIndex:4077

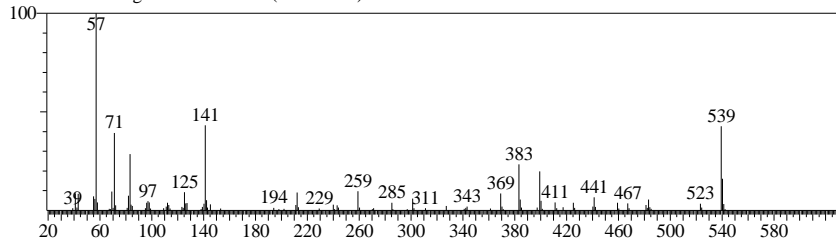
CompName:2-Butenoic acid, 2-methyl-, 1,1a,1b,4,4a,5,7a,7b,8,9-decahydro-4a,7b-dihydroxy-1,1,6,8-tetramethyl-



<< Target >>

Line#90 R.Time:32.717(Scan#:3567) BasePeak:57.10(1194047)

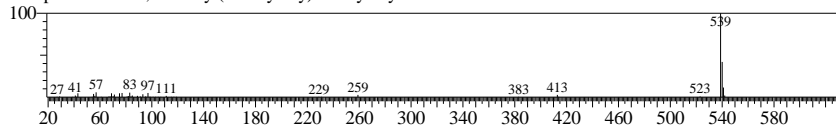
RawMode:Averaged 32.708-32.725(3566-3568) BG Mode:Calc. from Peak



Hit#1 Entry:187600 Library:NIST08.LIB

SI:50 Formula:C35H74O2Si CAS:0-00-0 MolWeight:554 RetIndex:3562

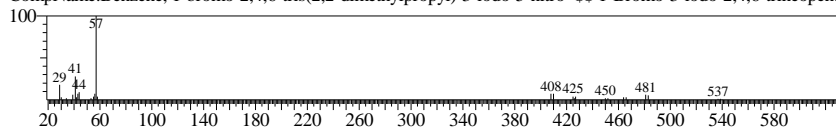
CompName:Silane, dimethyl(docosyloxy)undecyloxy-



Hit#2 Entry:186809 Library:NIST08.LIB

SI:45 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

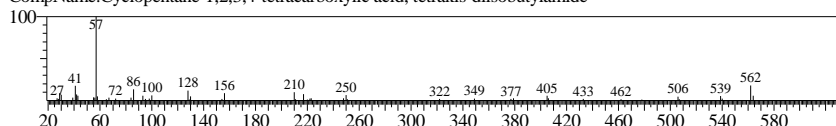
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- \$S\$ 1-Bromo-3-iodo-2,4,6-trineopent



Hit#3 Entry:190443 Library:NIST08.LIB

SI:39 Formula:C41H78N4O4 CAS:0-00-0 MolWeight:690 RetIndex:4316

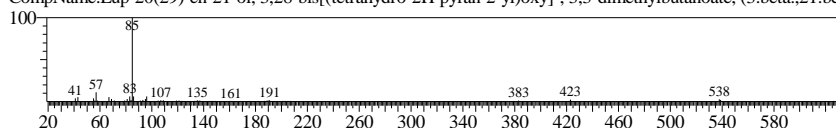
CompName:Cyclopentane-1,2,3,4-tetracarboxylic acid, tetrakis-diisobutylamide



Hit#4 Entry:190726 Library:NIST08.LIB

SI:37 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

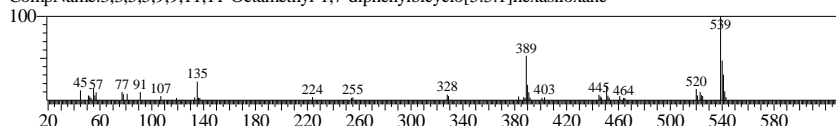
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



Hit#5 Entry:187547 Library:NIST08.LIB

SI:29 Formula:C20H34O7Si6 CAS:49538-51-4 MolWeight:554 RetIndex:2276

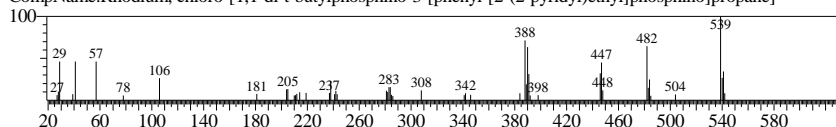
CompName:3,3,5,5,9,9,11,11-Octamethyl-1,7-diphenylbicyclo[5.5.1]hexasiloxane



Hit#6 Entry:186909 Library:NIST08.LIB

SI:27 Formula:C24H37ClNP2Rh CAS:0-00-0 MolWeight:539 RetIndex:0

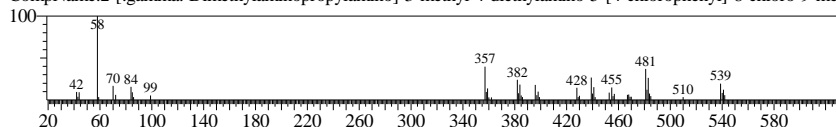
CompName:Rhodium, chloro-[1,1-di-t-butylphosphino-3-[phenyl-[2-(2-pyridyl)ethyl]phosphino]propane]



Hit#7 Entry:186919 Library:NIST08.LIB

SI:27 Formula:C29H35Cl2N5O CAS:0-00-0 MolWeight:539 RetIndex:4338

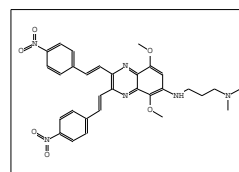
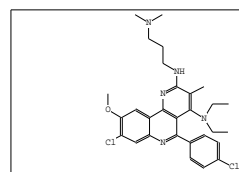
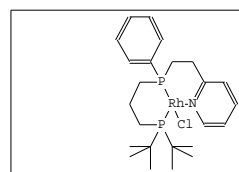
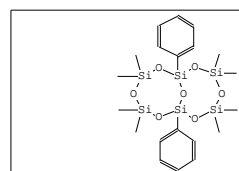
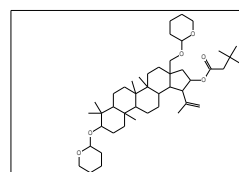
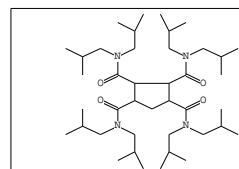
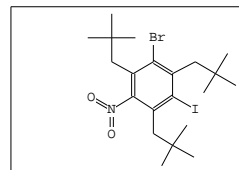
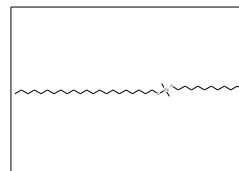
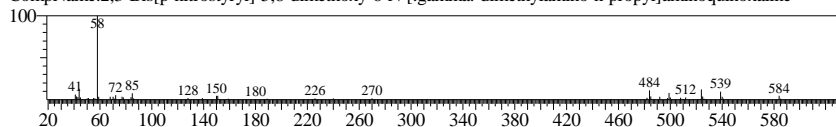
CompName:2-[gamma-Dimethylaminopropylamino]-3-methyl-4-diethylamino-5-[4-chlorophenyl]-8-chloro-9-me



Hit#8 Entry:188600 Library:NIST08.LIB

SI:24 Formula:C31H32N6O6 CAS:56393-52-3 MolWeight:584 RetIndex:4934

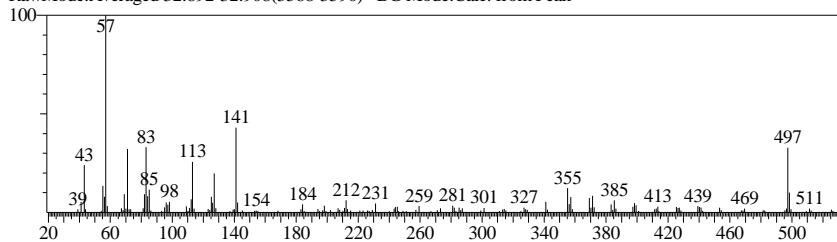
CompName:2,3-Bis[p-nitrotyrlyl]-5,8-dimethoxy-6-N-[gamma-dimethylamino-n-propyl]aminoquinoxaline



<< Target >>

Line#:91 R.Time:32.900(Scan#:3589) BasePeak:57.10(76131)

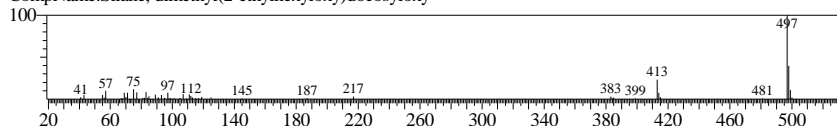
RawMode:Averaged 32.892-32.908(3588-3590) BG Mode:Calc. from Peak



Hit#:1 Entry:185524 Library:NIST08.LIB

SI:49 Formula:C32H68O2Si CAS:0-00-0 MolWeight:512 RetIndex:3200

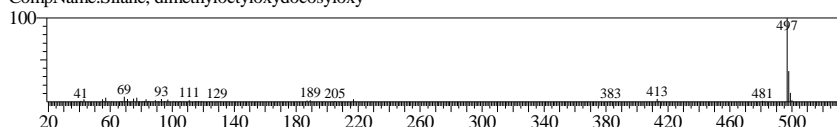
CompName:Silane, dimethyl(2-ethylhexyloxy)docosyloxy-



Hit#:2 Entry:185525 Library:NIST08.LIB

SI:42 Formula:C32H68O2Si CAS:0-00-0 MolWeight:512 RetIndex:3264

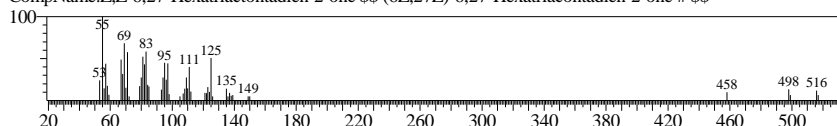
CompName:Silane, dimethyloctyloxydocosyloxy-



Hit#:3 Entry:185782 Library:NIST08.LIB

SI:41 Formula:C36H68O CAS:133530-19-5 MolWeight:516 RetIndex:3752

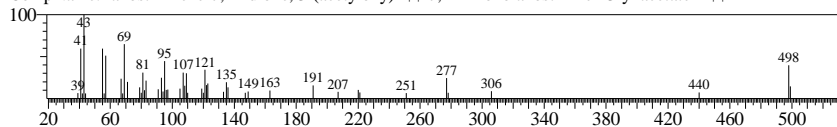
CompName:Z,Z-6,27-Hexatriacontadien-2-one (6Z,27Z)-6,27-Hexatriacontadien-2-one #



Hit#:4 Entry:184495 Library:NIST08.LIB

SI:38 Formula:C32H50O4 CAS:55700-81-7 MolWeight:498 RetIndex:3325

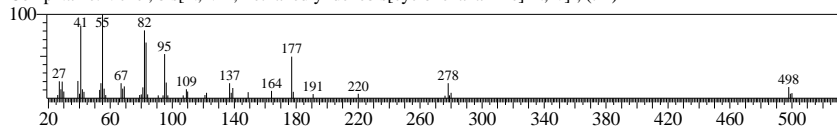
CompName:Lanost-24-ene-7,11-dione, 3-(acetyloxy)- 7,11-Dioxolanost-24-en-3-yl acetate #



Hit#:5 Entry:184469 Library:NIST08.LIB

SI:36 Formula:C28H48N4Ni CAS:63576-86-3 MolWeight:498 RetIndex:0

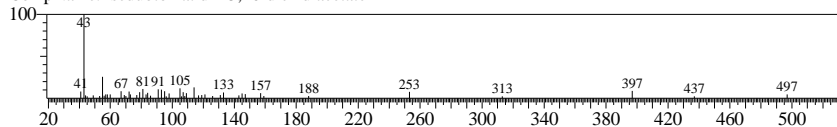
CompName:Nickel, bis[N,N'-1,2-ethanediyldenebis(cyclohexanamine)-N,N']-, (1-4)-



Hit#:6 Entry:184412 Library:NIST08.LIB

SI:35 Formula:C31H47NO4 CAS:0-00-0 MolWeight:497 RetIndex:3452

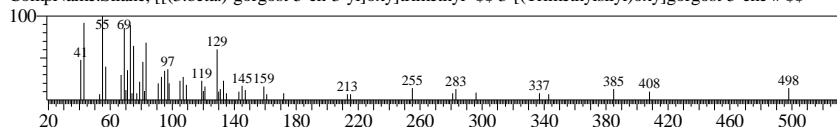
CompName:Pseudoatomatin-5,20-dien diacetate



Hit#:7 Entry:184517 Library:NIST08.LIB

SI:35 Formula:C33H58OSi CAS:55103-85-0 MolWeight:498 RetIndex:2826

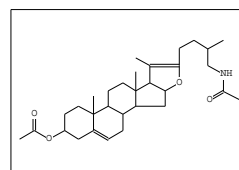
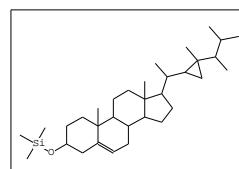
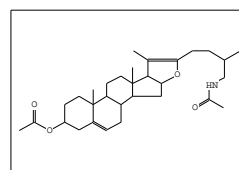
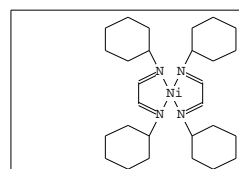
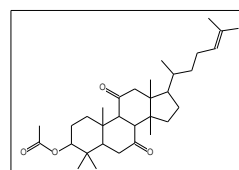
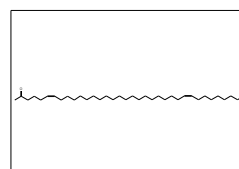
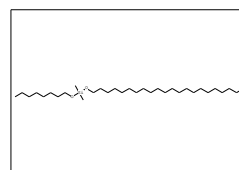
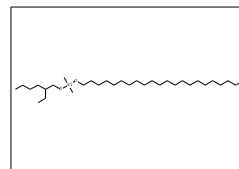
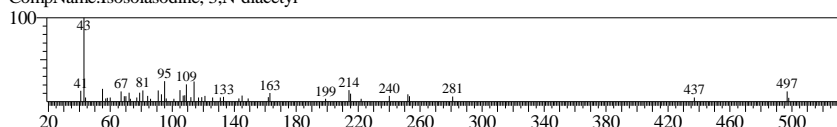
CompName:Silane, [(3.beta.)-gorgost-5-en-3-yl]oxy]trimethyl- 3-[(Trimethylsilyl)oxy]gorgost-5-ene #



Hit#:8 Entry:184413 Library:NIST08.LIB

SI:32 Formula:C31H47NO4 CAS:0-00-0 MolWeight:497 RetIndex:3452

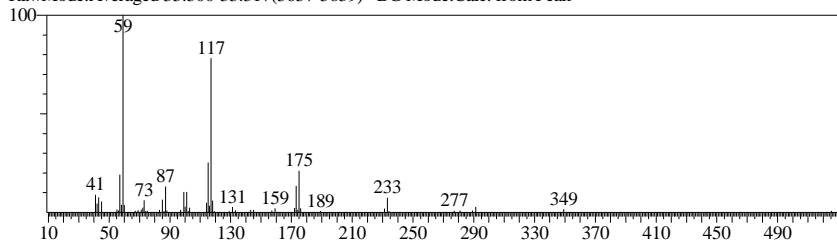
CompName:Isosolasodine, 3,N-diacetyl-



<< Target >>

Line# 92 R.Time:33.308(Scan#:3638) BasePeak:59.10(339267)

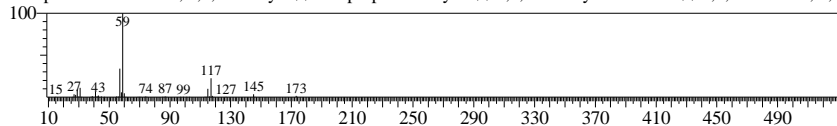
RawMode:Averaged 33.300-33.317(3637-3639) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:75 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

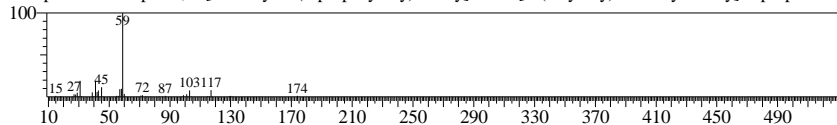
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:27550 Library:NIST08.LIB

SI:71 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

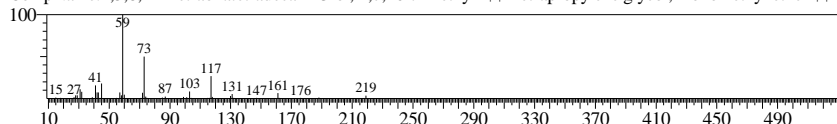
CompName:1-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#3 Entry:85908 Library:NIST08.LIB

SI:71 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

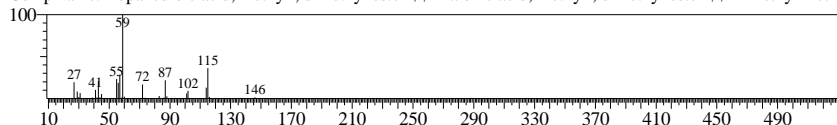
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#4 Entry:13447 Library:NIST08.LIB

SI:71 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

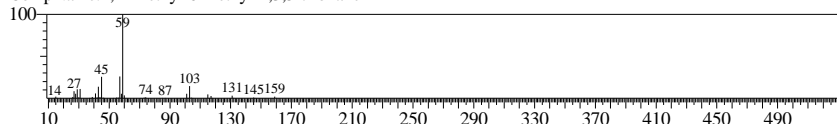
CompName:Propanedioic acid, methyl-, dimethyl ester \$\$ Malonic acid, methyl-, dimethyl ester \$\$ Dimethyl meth



Hit#5 Entry:19968 Library:NIST08.LIB

SI:70 Formula:C8H16O3 CAS:117888-04-7 MolWeight:160 RetIndex:1069

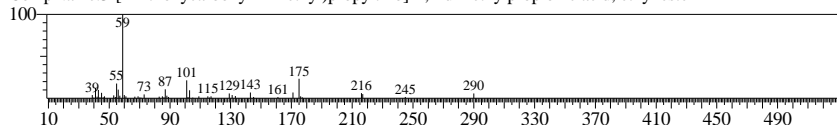
CompName:2,4-Diethyl-6-methyl-1,3,5-trioxane



Hit#6 Entry:104652 Library:NIST08.LIB

SI:70 Formula:C14H26O4S CAS:21153-32-2 MolWeight:290 RetIndex:1829

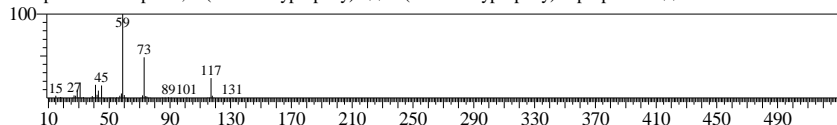
CompName:3-[2-Ethoxycarbonyl-2-methylpropylthio]-2,2-dimethylpropionic acid, ethyl ester



Hit#7 Entry:14297 Library:NIST08.LIB

SI:70 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

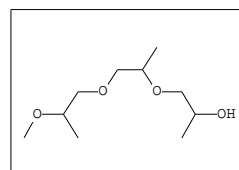
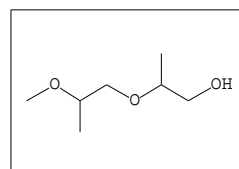
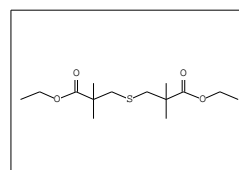
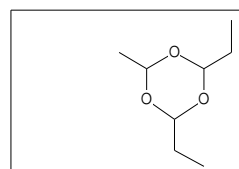
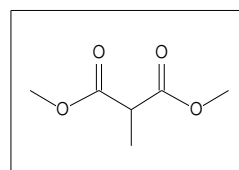
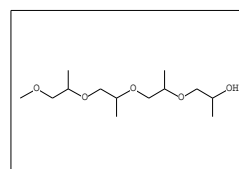
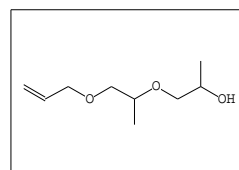
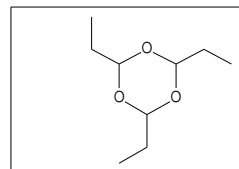
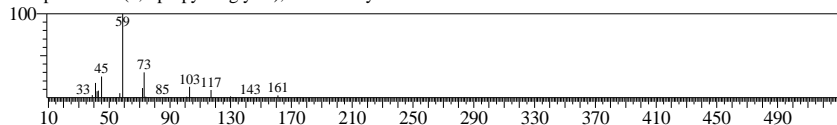
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



Hit#8 Entry:46316 Library:NIST08.LIB

SI:69 Formula:C10H22O4 CAS:0-00-0 MolWeight:206 RetIndex:1277

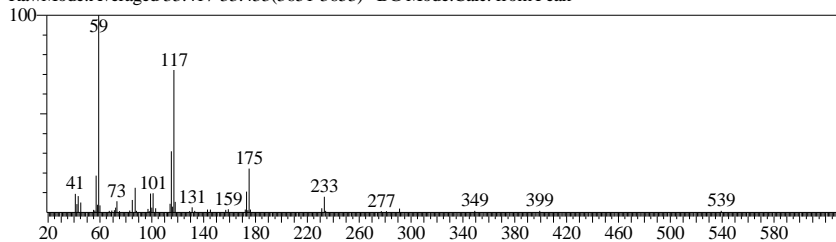
CompName:Tri(1,2-propyleneglycol), monomethyl ether



<< Target >>

Line# 93 R.Time:33.425(Scan#:3652) BasePeak:59.10(703695)

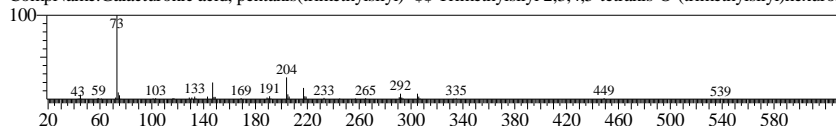
RawMode:Averaged 33.417-33.433(3651-3653) BG Mode:Calc. from Peak



Hit#:1 Entry:187550 Library:NIST08.LIB

SI:32 Formula:C21H50O7Si5 CAS:65337-27-1 MolWeight:554 RetIndex:2062

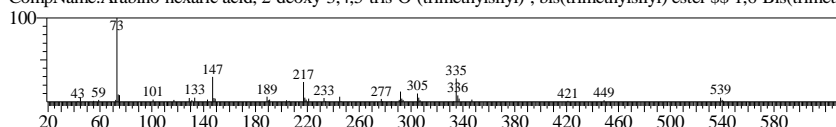
CompName:Galacturonic acid, pentakis(trimethylsilyl)- \$ \$ Trimethylsilyl 2,3,4,5-tetrakis-O-(trimethylsilyl)hexuror



Hit#:2 Entry:187549 Library:NIST08.LIB

SI:30 Formula:C21H50O7Si5 CAS:38165-98-9 MolWeight:554 RetIndex:2029

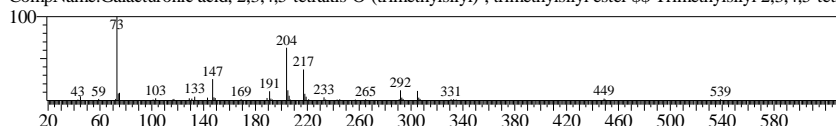
CompName:Arabino-hexaric acid, 2-deoxy-3,4,5-tris-O-(trimethylsilyl)-, bis(trimethylsilyl) ester \$ \$ 1,6-Bis(trimet



Hit#:3 Entry:187552 Library:NIST08.LIB

SI:29 Formula:C21H50O7Si5 CAS:56192-87-1 MolWeight:554 RetIndex:2062

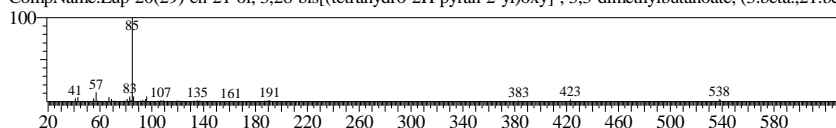
CompName:Galacturonic acid, 2,3,4,5-tetrakis-O-(trimethylsilyl)-, trimethylsilyl ester \$ \$ Trimethylsilyl 2,3,4,5-tetr



Hit#:4 Entry:190726 Library:NIST08.LIB

SI:25 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

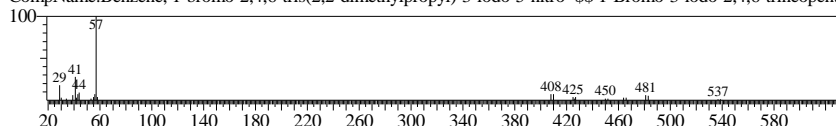
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



Hit#:5 Entry:186809 Library:NIST08.LIB

SI:24 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

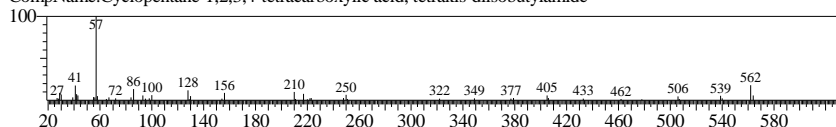
CompName:Benzene, 1-bromo-2,4,6-tris[(2,2-dimethylpropyl)-3-iodo-5-nitro-\$ \$ 1-Bromo-3-iodo-2,4,6-trineopent



Hit#:6 Entry:190443 Library:NIST08.LIB

SI:23 Formula:C41H78N4O4 CAS:0-00-0 MolWeight:690 RetIndex:4316

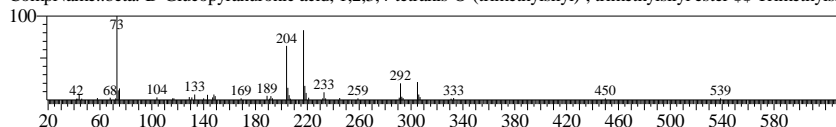
CompName:Cyclopentane-1,2,3,4-tetracarboxylic acid, tetrakis-diisobutylamide



Hit#:7 Entry:28136 Library:NIST08s.LIB

SI:20 Formula:C21H50O7Si5 CAS:52842-24-7 MolWeight:554 RetIndex:2128

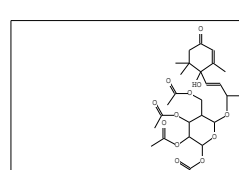
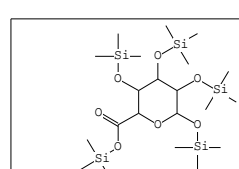
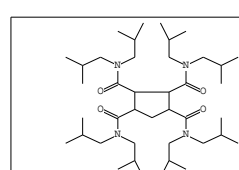
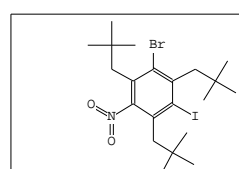
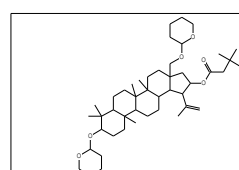
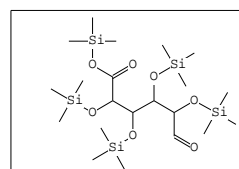
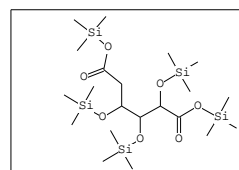
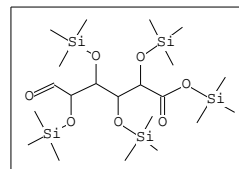
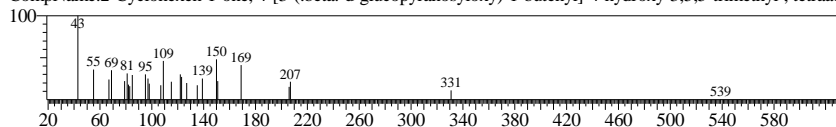
CompName:.beta.-D-Glucopyranuronic acid, 1,2,3,4-tetrakis-O-(trimethylsilyl)-, trimethylsilyl ester \$ \$ Trimethylsi



Hit#:8 Entry:187574 Library:NIST08.LIB

SI:20 Formula:C27H38O12 CAS:0-00-0 MolWeight:554 RetIndex:3686

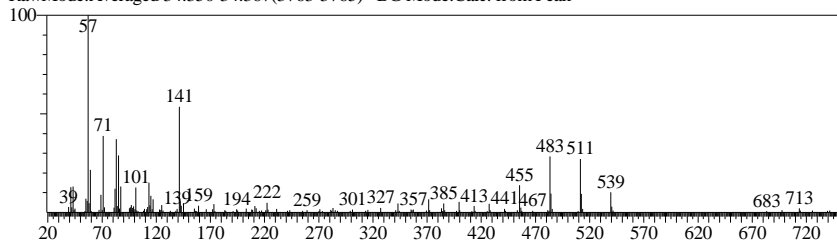
CompName:2-Cyclohexen-1-one, 4-[3-(.beta.-d-glucopyranosyloxy)-1-butenyl]-4-hydroxy-3,5,5-trimethyl-, tetraac



<< Target >>

Line# 94 R.Time:34.358(Scan#:3764) BasePeak:57.10(101210)

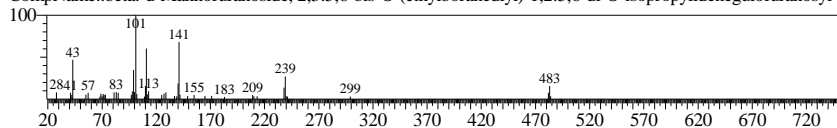
RawMode:Averaged 34.350-34.367(3763-3765) BG Mode:Calc. from Peak



Hit#:1 Entry:184447 Library:NIST08.LIB

SI:47 Formula:C22H36B2O11 CAS:0-00-0 MolWeight:498 RetIndex:0

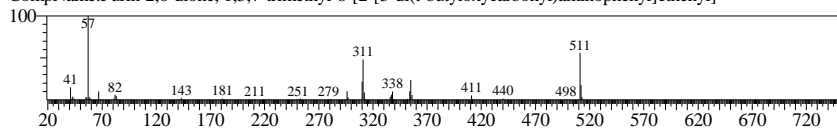
CompName:.beta.-d-Mannofuranoside, 2,3:5,6-bis-O-(ethylboranediyl)-1,2:5,6-di-O-isopropylidenegulofuranosyl-



Hit#:2 Entry:185431 Library:NIST08.LIB

SI:45 Formula:C26H33N5O6 CAS:0-00-0 MolWeight:511 RetIndex:3851

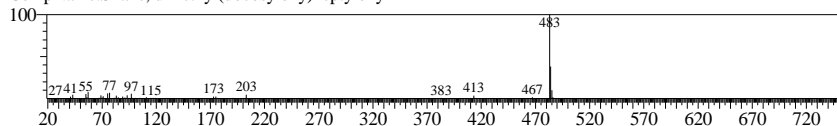
CompName:Purin-2,6-dione, 1,3,7-trimethyl-8-[2-[3-di(t-butyloxycarbonyl)aminophenyl]ethenyl]-



Hit#:3 Entry:184491 Library:NIST08.LIB

SI:43 Formula:C31H66O2Si CAS:0-00-0 MolWeight:498 RetIndex:3164

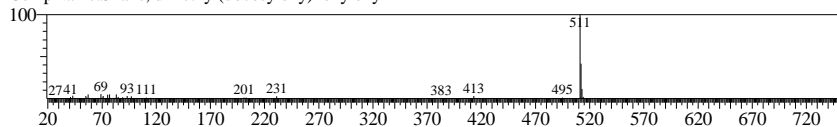
CompName:Silane, dimethyl(docosyloxy)heptyloxy-



Hit#:4 Entry:186328 Library:NIST08.LIB

SI:42 Formula:C33H70O2Si CAS:0-00-0 MolWeight:526 RetIndex:3363

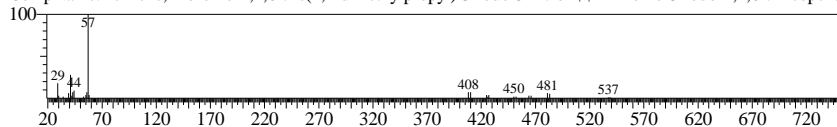
CompName:Silane, dimethyl(docosyloxy)nonyloxy-



Hit#:5 Entry:186809 Library:NIST08.LIB

SI:41 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

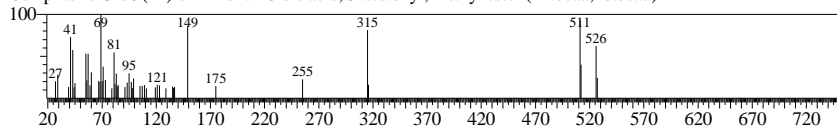
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- β -1-Bromo-3-iodo-2,4,6-trineopent



Hit#:6 Entry:186325 Library:NIST08.LIB

SI:40 Formula:C33H50O5 CAS:0-00-0 MolWeight:526 RetIndex:3451

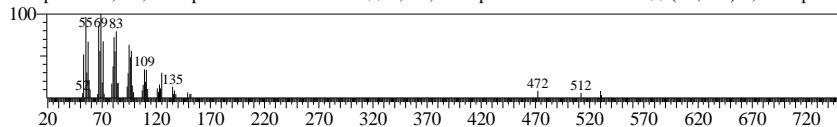
CompName:Urs-9(11)-en-12-one-28-oic acid, 3-acetoxy-, methyl ester (14.beta.,20.beta.)



Hit#:7 Entry:186505 Library:NIST08.LIB

SI:39 Formula:C37H70O CAS:133530-21-9 MolWeight:530 RetIndex:3851

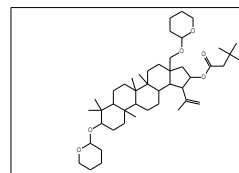
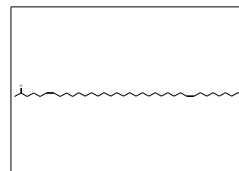
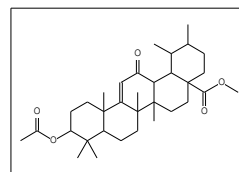
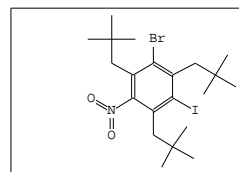
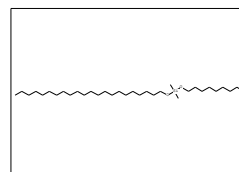
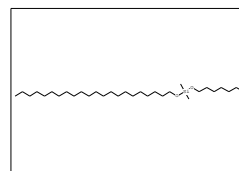
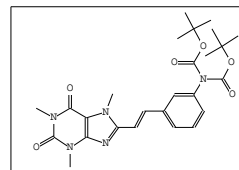
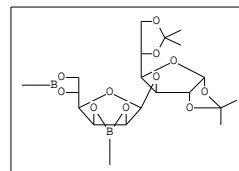
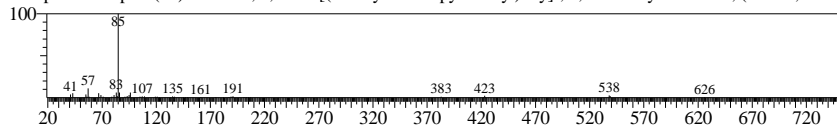
CompName:Z,Z-6,28-Heptatriacontadien-2-one β -Z,Z-6,28-Heptatriacontadien-2-one β (6Z,28Z)-6,28-Heptatri



Hit#:8 Entry:190726 Library:NIST08.LIB

SI:37 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

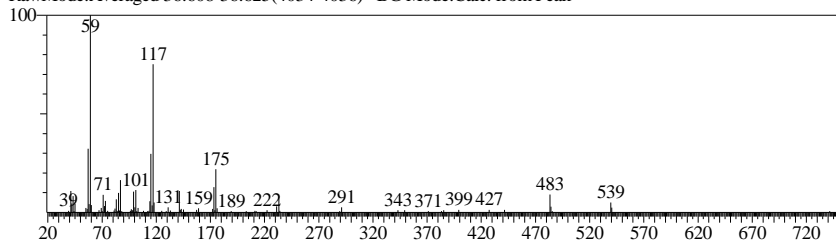
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



<< Target >>

Line#95 R.Time:36.617(Scan#:4035) BasePeak:59.10(514126)

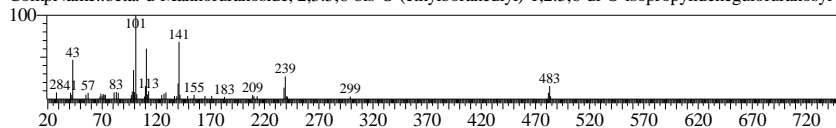
RawMode:Averaged 36.608-36.625(4034-4036) BG Mode:Calc. from Peak



Hit#:1 Entry:184447 Library:NIST08.LIB

SI:33 Formula:C22H36B2O11 CAS:0-00-0 MolWeight:498 RetIndex:0

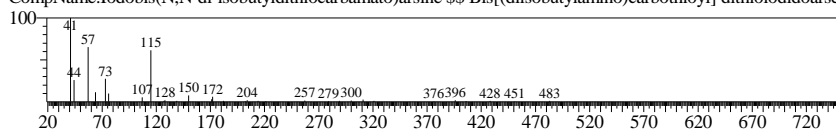
CompName:.beta.-d-Mannofuranoside, 2,3:5,6-bis-O-(ethylboranediyl)-1,2:5,6-di-O-isopropylidenegulofuranosyl-



Hit#:2 Entry:189212 Library:NIST08.LIB

SI:31 Formula:C18H36AsIN2S4 CAS:59196-55-3 MolWeight:610 RetIndex:0

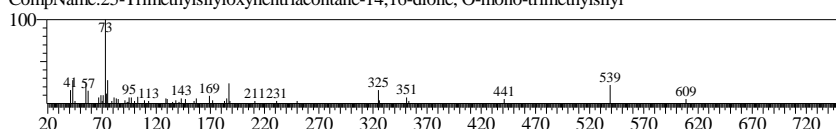
CompName:Iodobis(N,N-di-isobutylthiocarbamato)arsine \$\$ Bis[(diisobutylamino)carbothioyl] dithioiodidoarse



Hit#:3 Entry:189513 Library:NIST08.LIB

SI:31 Formula:C37H76O3Si2 CAS:0-00-0 MolWeight:624 RetIndex:3728

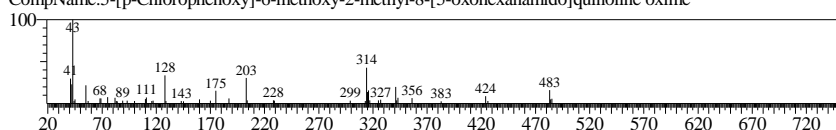
CompName:25-Trimethylsilyloxyhentriacontane-14,16-dione, O-mono-trimethylsilyl



Hit#:4 Entry:183114 Library:NIST08.LIB

SI:30 Formula:C25H26ClN3O5 CAS:0-00-0 MolWeight:483 RetIndex:4040

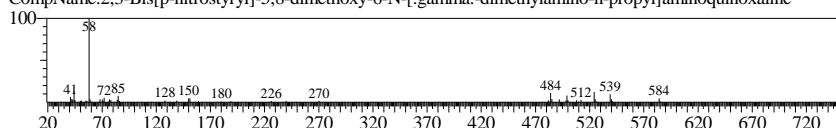
CompName:5-[p-Chlorophenoxyl]-6-methoxy-2-methyl-8-[5-oxohexanamido]quinoline oxime



Hit#:5 Entry:188600 Library:NIST08.LIB

SI:30 Formula:C31H32N6O6 CAS:56393-52-3 MolWeight:584 RetIndex:4934

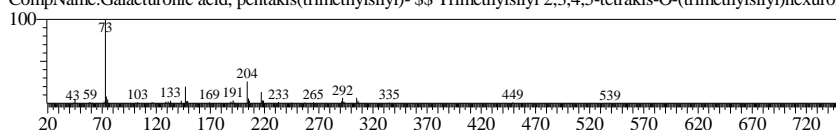
CompName:2,3-Bis[p-nitrostyryl]-5,8-dimethoxy-6-N-[gamma.-dimethylamino-n-propyl]aminoquinoxaline



Hit#:6 Entry:187550 Library:NIST08.LIB

SI:28 Formula:C21H50O7Si5 CAS:65337-27-1 MolWeight:554 RetIndex:2062

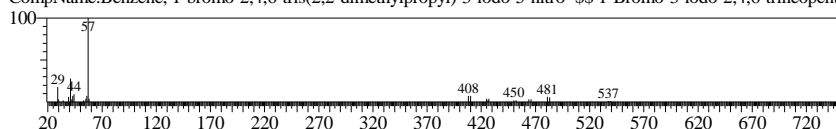
CompName:Galacturonic acid, pentakis(trimethylsilyl)- \$\$ Trimethylsilyl 2,3,4,5-tetrakis-O-(trimethylsilyl)hexuror



Hit#:7 Entry:186809 Library:NIST08.LIB

SI:28 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

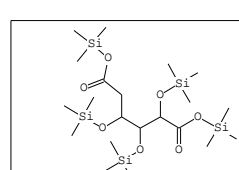
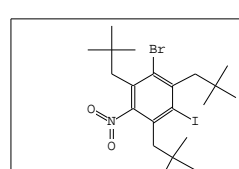
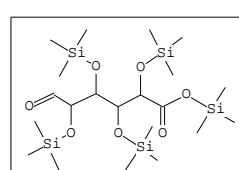
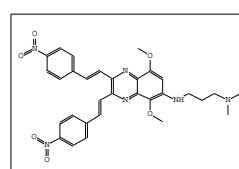
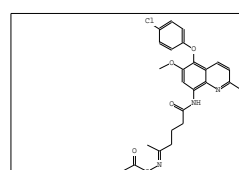
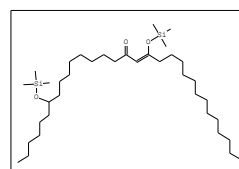
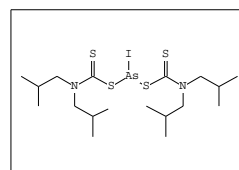
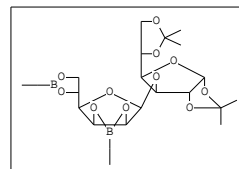
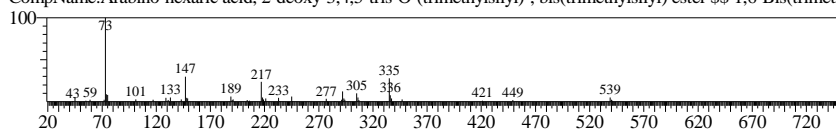
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- \$\$ 1-Bromo-3-iodo-2,4,6-trineopent



Hit#:8 Entry:187549 Library:NIST08.LIB

SI:27 Formula:C21H50O7Si5 CAS:38165-98-9 MolWeight:554 RetIndex:2029

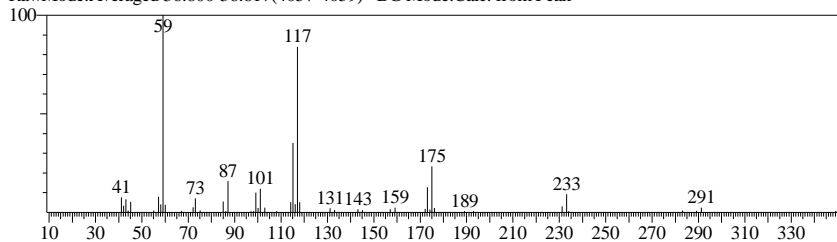
CompName:Arabino-hexaric acid, 2-deoxy-3,4,5-tris-O-(trimethylsilyl)-, bis(trimethylsilyl) ester \$\$ 1,6-Bis(trimet



<< Target >>

Line#96 R.Time:36.808(Scan#:4058) BasePeak:59.10(306816)

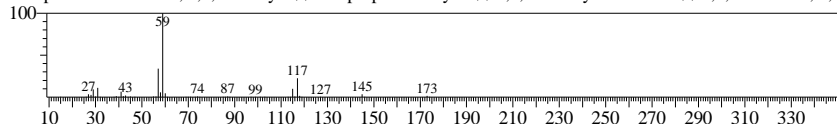
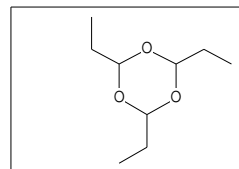
RawMode:Averaged 36.800-36.817(4057-4059) BG Mode:Calc. from Peak



Hit#1 Entry:27552 Library:NIST08.LIB

SI:72 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

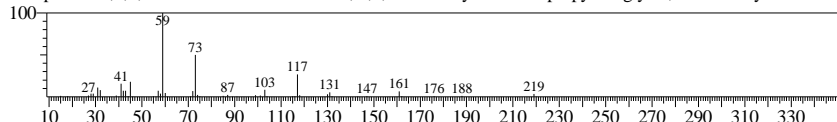
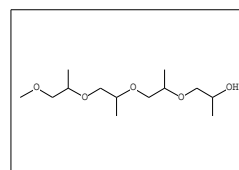
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#2 Entry:85908 Library:NIST08.LIB

SI:70 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

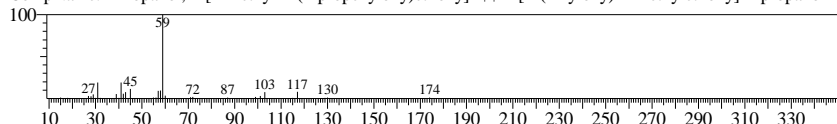
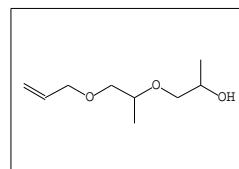
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$



Hit#3 Entry:27550 Library:NIST08.LIB

SI:69 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

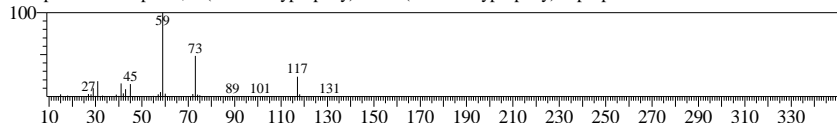
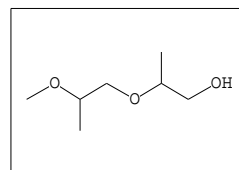
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#4 Entry:14297 Library:NIST08.LIB

SI:69 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

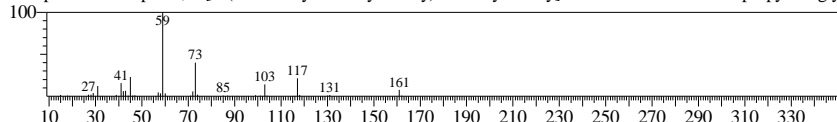
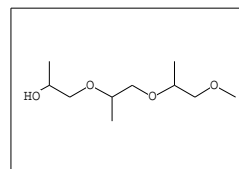
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol #



Hit#5 Entry:17307 Library:NIST08.LIB

SI:69 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

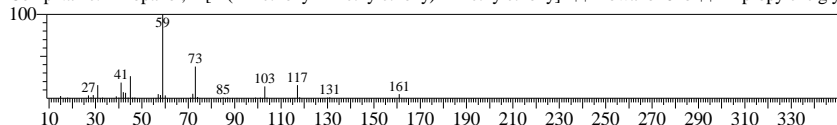
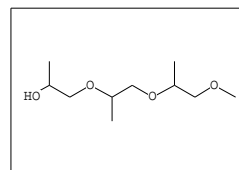
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#6 Entry:46317 Library:NIST08.LIB

SI:69 Formula:C10H22O4 CAS:20324-33-8 MolWeight:206 RetIndex:1277

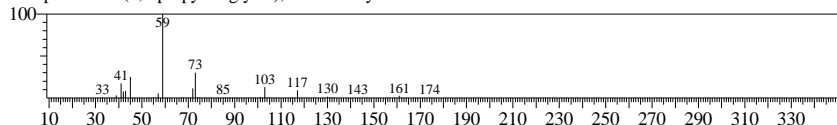
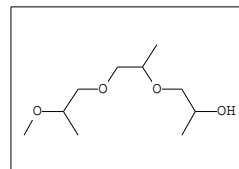
CompName:2-Propanol, 1-[2-(2-methoxy-1-methylethoxy)-1-methylethoxy]- \$\$ Dowanol 62b \$\$ Tripropylene gly



Hit#7 Entry:46316 Library:NIST08.LIB

SI:69 Formula:C10H22O4 CAS:0-00-0 MolWeight:206 RetIndex:1277

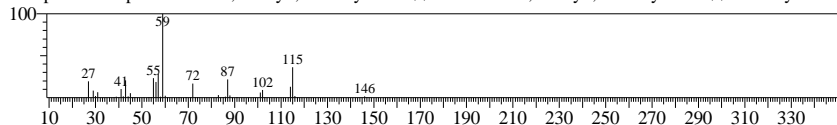
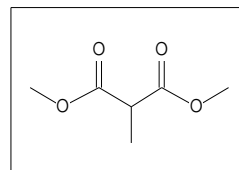
CompName:Tri(1,2-propyleneglycol), monomethyl ether



Hit#8 Entry:13447 Library:NIST08.LIB

SI:69 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

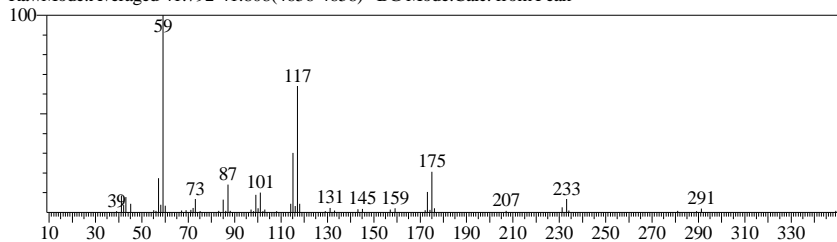
CompName:Propanedioic acid, methyl-, dimethyl ester \$\$ Malonic acid, methyl-, dimethyl ester \$\$ Dimethyl meth



<< Target >>

Line#:97 R.Time:41.800(Scan#:4657) BasePeak:59.10(1171434)

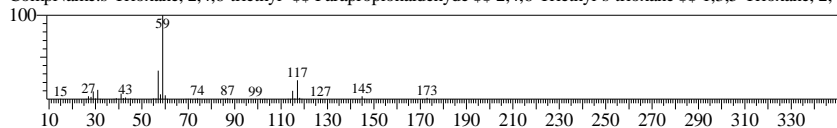
RawMode:Averaged 41.792-41.808(4656-4658) BG Mode:Calc. from Peak



Hit#:1 Entry:27552 Library:NIST08.LIB

SI:76 Formula:C9H18O3 CAS:2396-42-1 MolWeight:174 RetIndex:1168

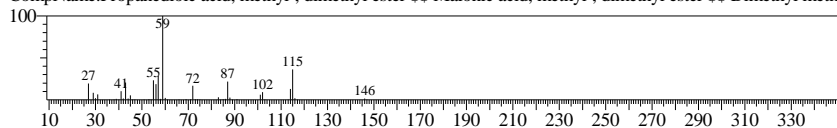
CompName:s-Trioxane, 2,4,6-triethyl- \$\$ Parapropionaldehyde \$\$ 2,4,6-Triethyl-s-trioxane \$\$ 1,3,5-Trioxane, 2,4



Hit#:2 Entry:13447 Library:NIST08.LIB

SI:73 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

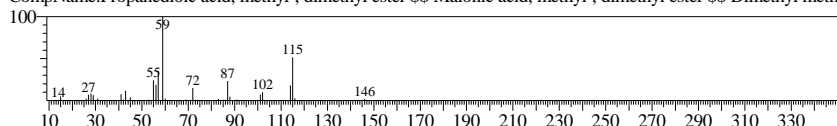
CompName:Propanedioic acid, methyl-, dimethyl ester \$\$ Malonic acid, methyl-, dimethyl ester \$\$ Dimethyl meth



Hit#:3 Entry:7890 Library:NIST08.LIB

SI:73 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

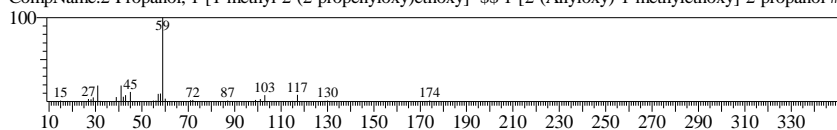
CompName:Propanedioic acid, methyl-, dimethyl ester \$\$ Malonic acid, methyl-, dimethyl ester \$\$ Dimethyl meth



Hit#:4 Entry:27550 Library:NIST08.LIB

SI:72 Formula:C9H18O3 CAS:55956-25-7 MolWeight:174 RetIndex:1156

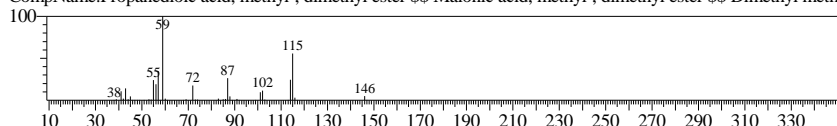
CompName:2-Propanol, 1-[1-methyl-2-(2-propenyloxy)ethoxy]- \$\$ 1-[2-(Allyloxy)-1-methylethoxy]-2-propanol #



Hit#:5 Entry:7889 Library:NIST08.LIB

SI:72 Formula:C6H10O4 CAS:609-02-9 MolWeight:146 RetIndex:888

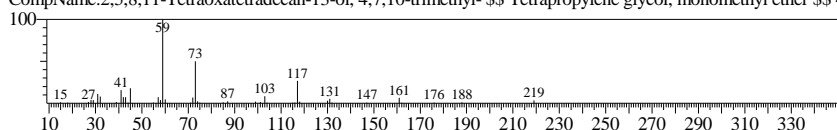
CompName:Propanedioic acid, methyl-, dimethyl ester \$\$ Malonic acid, methyl-, dimethyl ester \$\$ Dimethyl meth



Hit#:6 Entry:85908 Library:NIST08.LIB

SI:72 Formula:C13H28O5 CAS:20324-34-9 MolWeight:264 RetIndex:1587

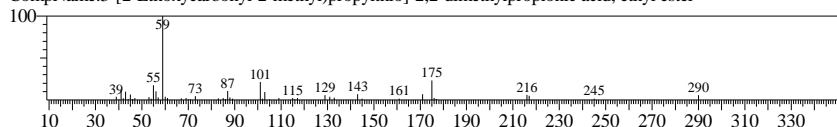
CompName:2,5,8,11-Tetraoxatetradecan-13-ol, 4,7,10-trimethyl- \$\$ Tetrapropylene glycol, monomethyl ether \$\$ 4



Hit#:7 Entry:104652 Library:NIST08.LIB

SI:71 Formula:C14H26O4S CAS:21153-32-2 MolWeight:290 RetIndex:1829

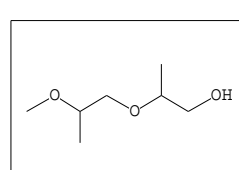
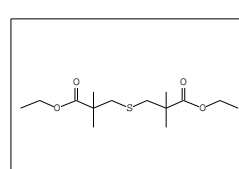
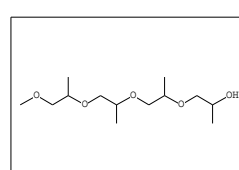
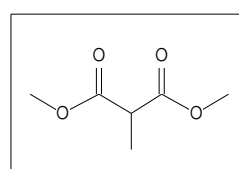
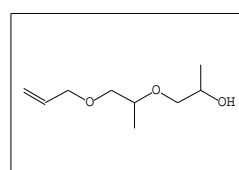
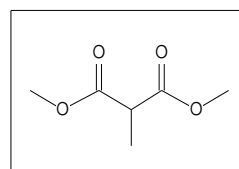
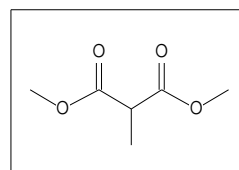
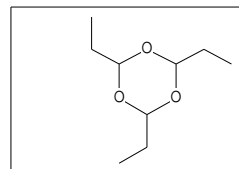
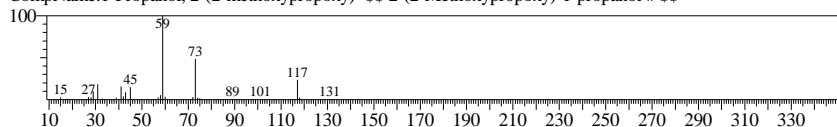
CompName:3-[2-Ethoxycarbonyl-2-methyl)propylthio]-2,2-dimethylpropionic acid, ethyl ester



Hit#:8 Entry:14297 Library:NIST08.LIB

SI:71 Formula:C7H16O3 CAS:13588-28-8 MolWeight:148 RetIndex:983

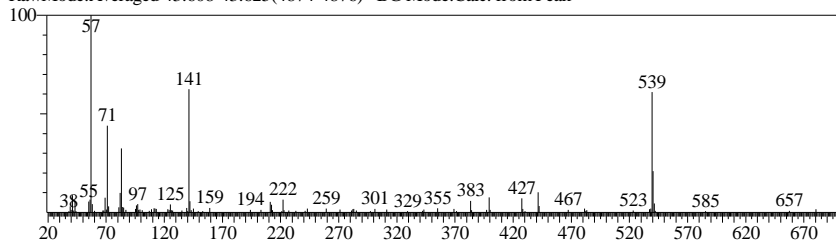
CompName:1-Propanol, 2-(2-methoxypropoxy)- \$\$ 2-(2-Methoxypropoxy)-1-propanol # \$\$



<< Target >>

Line#98 R.Time:43.617(Scan#:4875) BasePeak:57.10(83741)

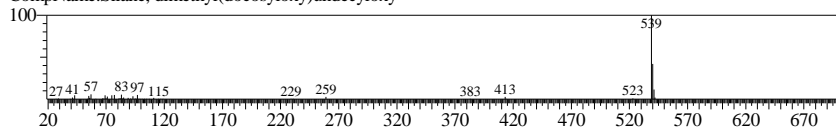
RawMode:Averaged 43.608-43.625(4874-4876) BG Mode:Calc. from Peak



Hit#1 Entry:187600 Library:NIST08.LIB

SI:56 Formula:C35H74O2Si CAS:0-00-0 MolWeight:554 RetIndex:3562

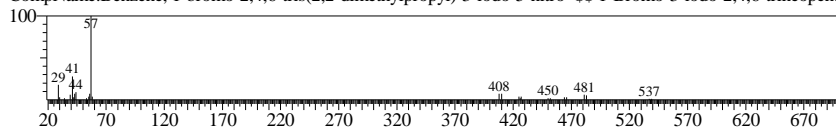
CompName:Silane, dimethyl(docosyloxy)undecyloxy-



Hit#2 Entry:186809 Library:NIST08.LIB

SI:47 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

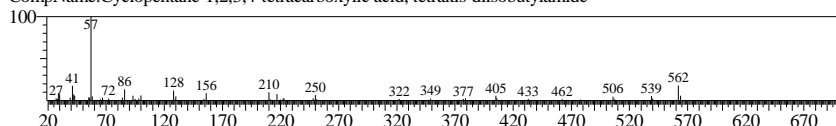
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- \$S\$ 1-Bromo-3-iodo-2,4,6-trineopent



Hit#3 Entry:190443 Library:NIST08.LIB

SI:43 Formula:C41H78N4O4 CAS:0-00-0 MolWeight:690 RetIndex:4316

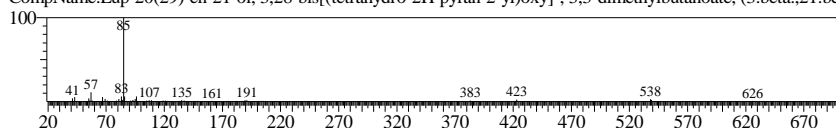
CompName:Cyclopentane-1,2,3,4-tetracarboxylic acid, tetrakis-diisobutylamide



Hit#4 Entry:190726 Library:NIST08.LIB

SI:42 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

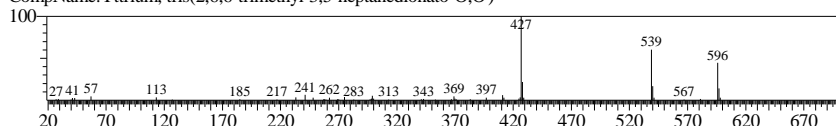
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



Hit#5 Entry:188910 Library:NIST08.LIB

SI:39 Formula:C30H51O6Y CAS:36502-83-7 MolWeight:596 RetIndex:0

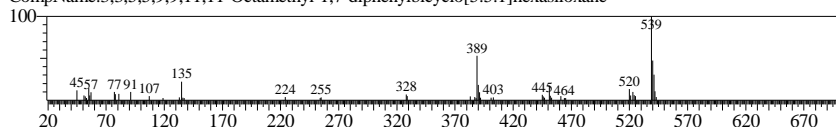
CompName:Yttrium, tris(2,6,6-trimethyl-3,5-heptanedionato-O,O')-



Hit#6 Entry:187547 Library:NIST08.LIB

SI:32 Formula:C20H34O7Si6 CAS:49538-51-4 MolWeight:554 RetIndex:2276

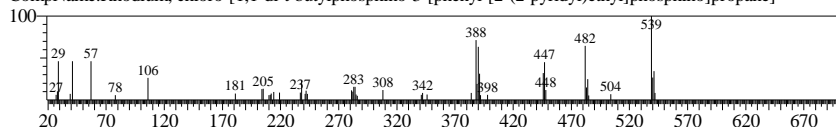
CompName:3,3,5,5,9,9,11,11-Octamethyl-1,7-diphenylbicyclo[5.5.1]hexasiloxane



Hit#7 Entry:186909 Library:NIST08.LIB

SI:29 Formula:C24H37ClNP2Rh CAS:0-00-0 MolWeight:539 RetIndex:0

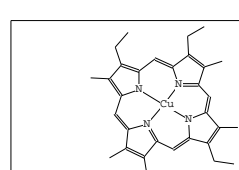
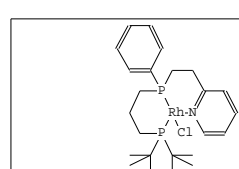
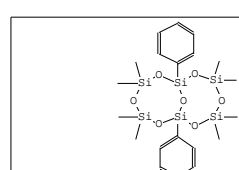
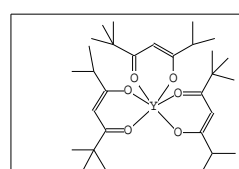
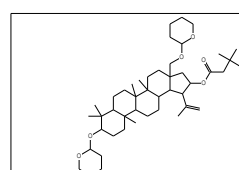
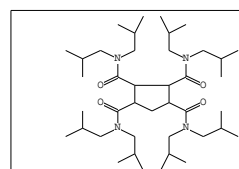
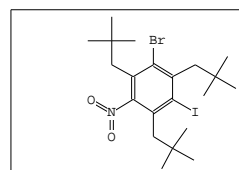
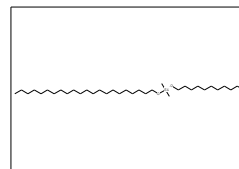
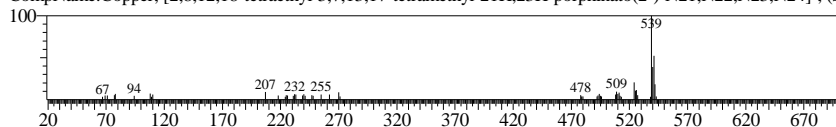
CompName:Rhodium, chloro-[1,1-di-t-butylphosphino-3-[phenyl-[2-(2-pyridyl)ethyl]phosphino]propane]



Hit#8 Entry:186923 Library:NIST08.LIB

SI:28 Formula:C32H36CuN4 CAS:14055-18-6 MolWeight:539 RetIndex:0

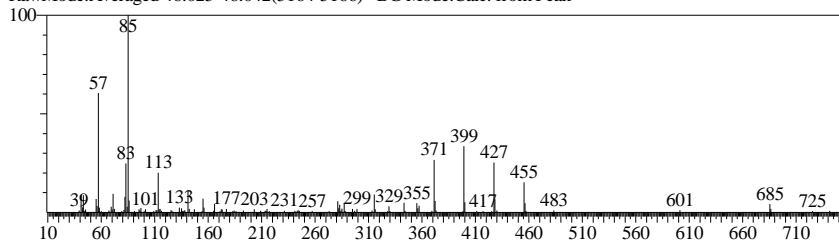
CompName:Copper, [2,8,12,18-tetraethyl-3,7,13,17-tetramethyl-21H,23H-porphinato(2-)-N21,N22,N23,N24]-, (S



<< Target >>

Line#:99 R.Time:46.033(Scan#:5165) BasePeak:85.10(53699)

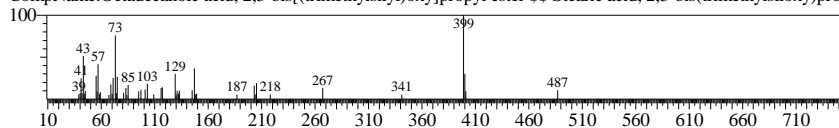
RawMode:Averaged 46.025-46.042(5164-5166) BG Mode:Calc. from Peak



Hit#:1 Entry:28029 Library:NIST08s.LIB

SI:38 Formula:C27H58O4Si2 CAS:1188-75-6 MolWeight:502 RetIndex:2780

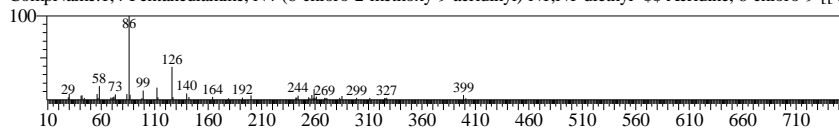
CompName:Octadecanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$\$ Stearic acid, 2,3-bis(trimethylsiloxy)pro



Hit#:2 Entry:27290 Library:NIST08s.LIB

SI:26 Formula:C23H30ClN3O CAS:83-89-6 MolWeight:399 RetIndex:3237

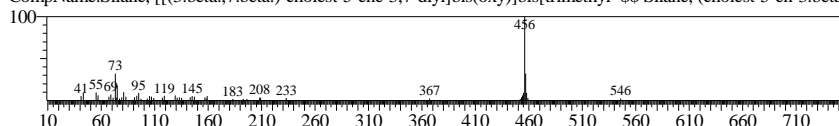
CompName:1,4-Pentanediamine, N4-(6-chloro-2-methoxy-9-acridinyl)-N1,N1-diethyl- \$\$ Acridine, 6-chloro-9-[[4



Hit#:3 Entry:28125 Library:NIST08s.LIB

SI:24 Formula:C33H62O2Si2 CAS:33287-26-2 MolWeight:546 RetIndex:2900

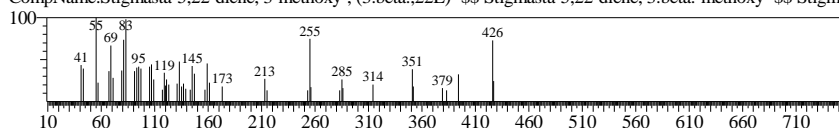
CompName:Silane, [[(3.beta.,7.beta.)-cholest-5-ene-3,7-diyl]bis(oxy)]bis(trimethyl- \$\$ Silane, (cholest-5-en-3.beta



Hit#:4 Entry:27615 Library:NIST08s.LIB

SI:23 Formula:C30H50O CAS:10453-25-5 MolWeight:426 RetIndex:2688

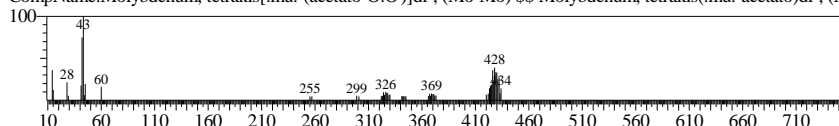
CompName:Stigmasta-5,22-diene, 3-methoxy-, (3.beta.,22E)- \$\$ Stigmasta-5,22-diene, 3.beta.-methoxy- \$\$ Stigm



Hit#:5 Entry:27668 Library:NIST08s.LIB

SI:22 Formula:C8H12Mo2O8 CAS:14221-06-8 MolWeight:432 RetIndex:0

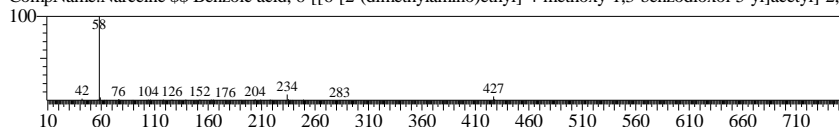
CompName:Molybdenum, tetrakis[.mu.-(acetato-O:O')]di-, (Mo-Mo) \$\$ Molybdenum, tetrakis(.mu.-acetato)di-, (



Hit#:6 Entry:27757 Library:NIST08s.LIB

SI:20 Formula:C23H27NO8 CAS:131-28-2 MolWeight:445 RetIndex:3544

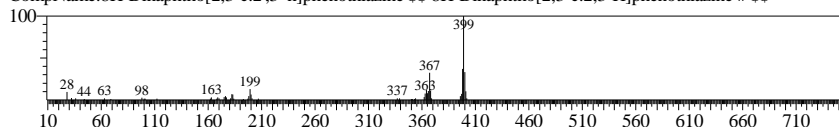
CompName:Narceine \$\$ Benzoic acid, 6-[[6-[2-(dimethylamino)ethyl]-4-methoxy-1,3-benzodioxol-5-yl]acetyl]-2,



Hit#:7 Entry:27291 Library:NIST08s.LIB

SI:15 Formula:C28H17NS CAS:222-06-0 MolWeight:399 RetIndex:3981

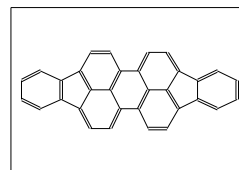
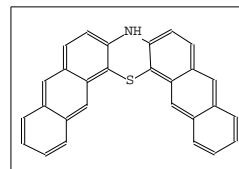
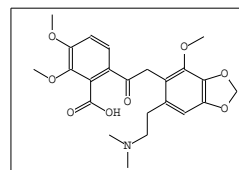
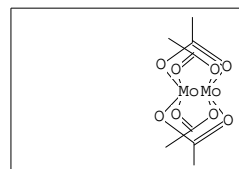
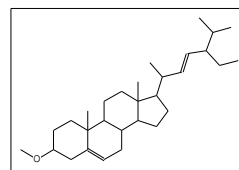
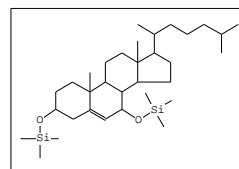
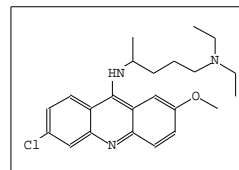
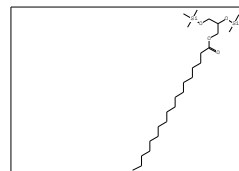
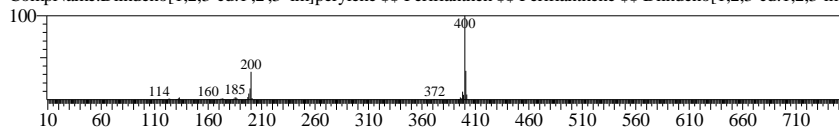
CompName:8H-Dinaphtho[2,3-c:2',3'-h]phenothiazine \$\$ 8H-Dinaphtho[2,3-c:2,3-h]phenothiazine # \$\$



Hit#:8 Entry:27311 Library:NIST08s.LIB

SI:15 Formula:C32H16 CAS:188-94-3 MolWeight:400 RetIndex:3739

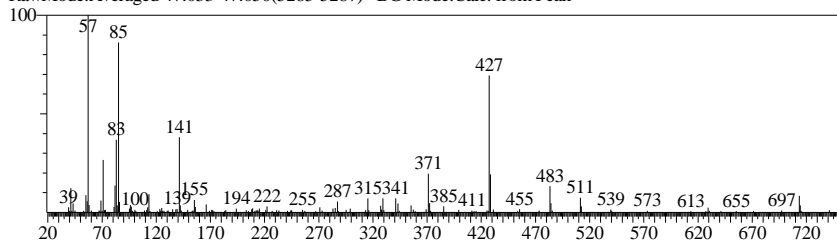
CompName:Diindenol[1,2,3-cd:1',2',3'-lm]perylene \$\$ Periflanthen \$\$ Diindenol[1,2,3-cd:1,2,3-lr



<< Target >>

Line#:100 R.Time:47.042(Scan#:5286) BasePeak:57.10(53776)

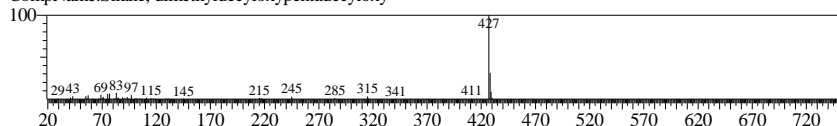
RawMode:Averaged 47.033-47.050(5285-5287) BG Mode:Calc. from Peak



Hit#:1 Entry:177735 Library:NIST08.LIB

SI:52 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

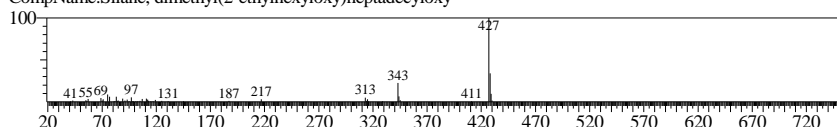
CompName:Silane, dimethyldecyloxypentadecyloxy-



Hit#:2 Entry:177733 Library:NIST08.LIB

SI:51 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2703

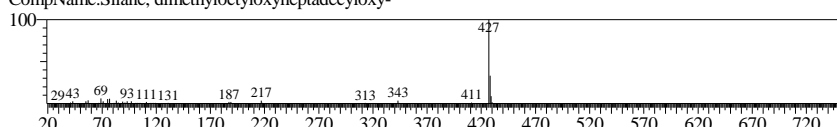
CompName:Silane, dimethyl(2-ethylhexyloxy)heptadecyloxy-



Hit#:3 Entry:177734 Library:NIST08.LIB

SI:50 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

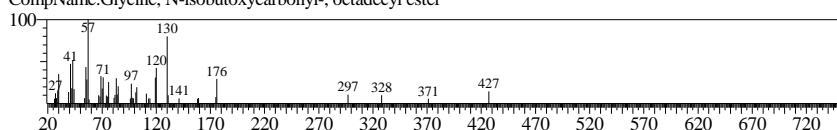
CompName:Silane, dimethyloctyloxyheptadecyloxy-



Hit#:4 Entry:174750 Library:NIST08.LIB

SI:50 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975

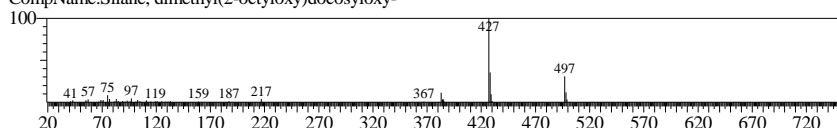
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



Hit#:5 Entry:185523 Library:NIST08.LIB

SI:46 Formula:C32H68O2Si CAS:0-00-0 MolWeight:512 RetIndex:3200

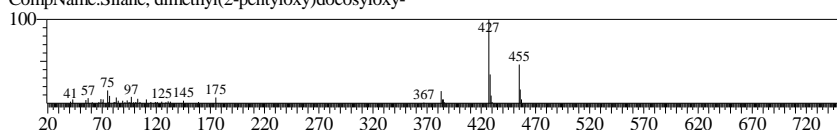
CompName:Silane, dimethyl(2-octyloxy)docosyloxy-



Hit#:6 Entry:181721 Library:NIST08.LIB

SI:45 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902

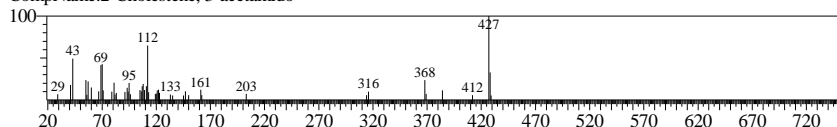
CompName:Silane, dimethyl(2-pentyloxy)docosyloxy-



Hit#:7 Entry:174844 Library:NIST08.LIB

SI:41 Formula:C29H49NO CAS:0-00-0 MolWeight:427 RetIndex:2969

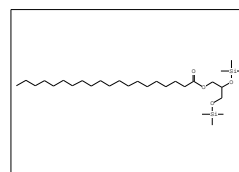
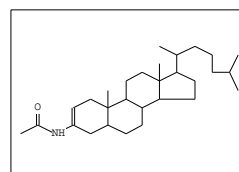
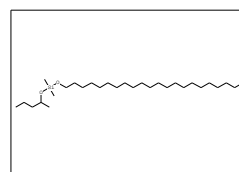
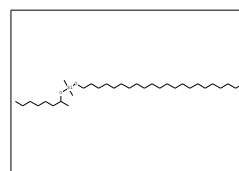
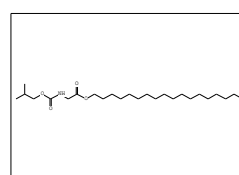
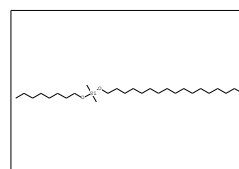
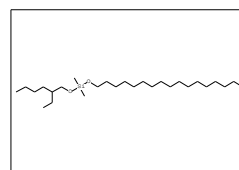
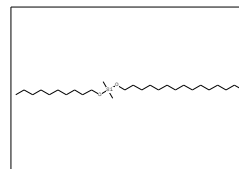
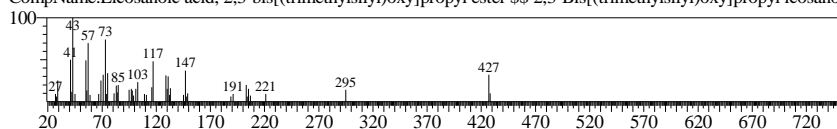
CompName:2-Cholestene, 3-acetamido-



Hit#:8 Entry:186484 Library:NIST08.LIB

SI:40 Formula:C29H62O4Si2 CAS:55517-94-7 MolWeight:530 RetIndex:2979

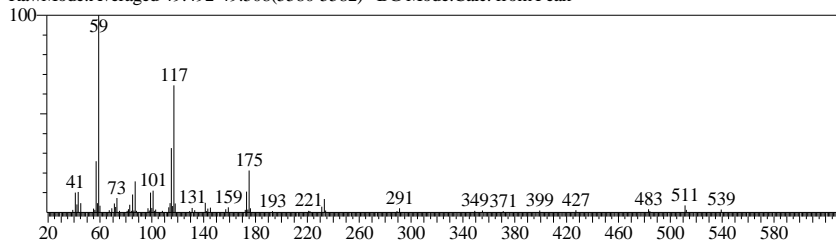
CompName:Eicosanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$\$ 2,3-Bis[(trimethylsilyl)oxy]propyl icosano



<< Target >>

Line#:101 R.Time:49.500(Scan#:5581) BasePeak:59.10(363316)

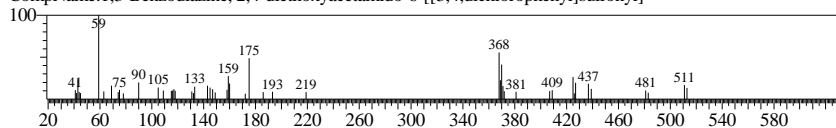
RawMode:Averaged 49.492-49.508(5580-5582) BG Mode:Calc. from Peak



Hit#:1 Entry:186963 Library:NIST08.LIB

SI:42 Formula:C22H22Cl2N4O6S CAS:0-00-0 MolWeight:540 RetIndex:4547

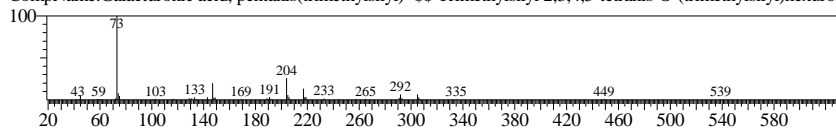
CompName:1,3-Benzodiazine, 2,4-diethoxyacetamido-6-[[3,4-dichlorophenyl]sulfonyl]



Hit#:2 Entry:187550 Library:NIST08.LIB

SI:30 Formula:C21H50O7Si5 CAS:65337-27-1 MolWeight:554 RetIndex:2062

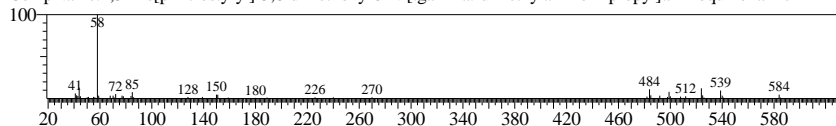
CompName:Galacturonic acid, pentakis(trimethylsilyl)- \$ \$ Trimethylsilyl 2,3,4,5-tetrakis-O-(trimethylsilyl)hexuror



Hit#:3 Entry:188600 Library:NIST08.LIB

SI:29 Formula:C31H32N6O6 CAS:56393-52-3 MolWeight:584 RetIndex:4934

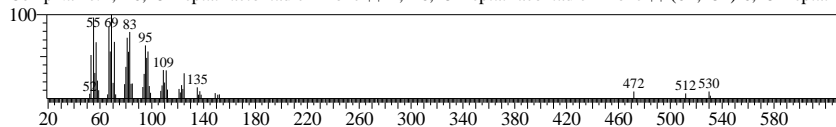
CompName:2,3-Bis[p-nitrotyrilyl]-5,8-dimethoxy-6-N-[gamma-dimethylamino-n-propyl]aminoquinoxaline



Hit#:4 Entry:186505 Library:NIST08.LIB

SI:28 Formula:C37H70O CAS:133530-21-9 MolWeight:530 RetIndex:3851

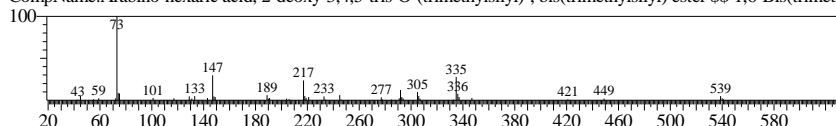
CompName:Z,Z-6,28-Heptatriacontadien-2-one \$ \$ Z,Z-6,28-Heptatriacontadien-2-one \$ \$ (6Z,28Z)-6,28-Heptatri



Hit#:5 Entry:187549 Library:NIST08.LIB

SI:27 Formula:C21H50O7Si5 CAS:38165-98-9 MolWeight:554 RetIndex:2029

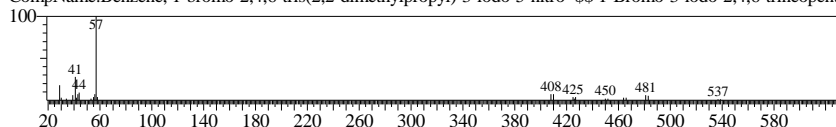
CompName:Arabino-hexaric acid, 2-deoxy-3,4,5-tris-O-(trimethylsilyl)-, bis(trimethylsilyl) ester \$ \$ 1,6-Bis(trimet



Hit#:6 Entry:186809 Library:NIST08.LIB

SI:27 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

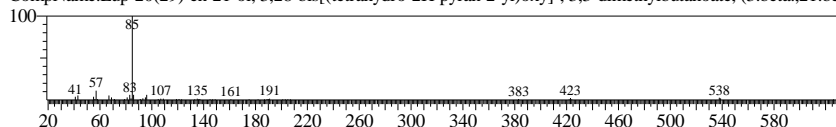
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- \$ \$ 1-Bromo-3-iodo-2,4,6-trineopent



Hit#:7 Entry:190726 Library:NIST08.LIB

SI:27 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

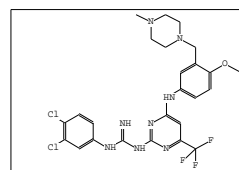
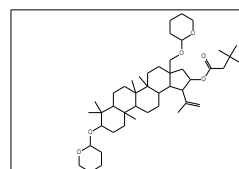
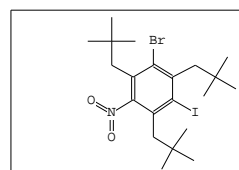
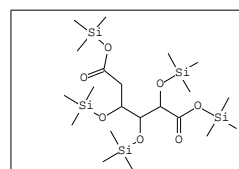
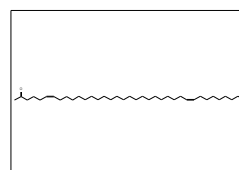
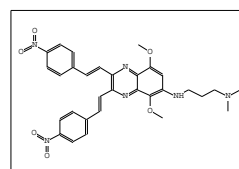
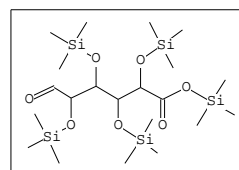
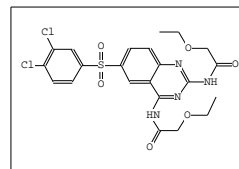
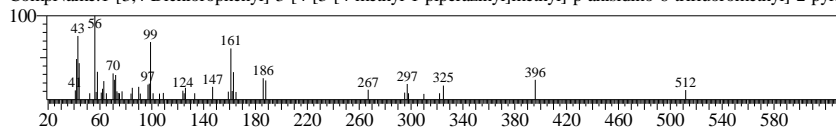
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



Hit#:8 Entry:188524 Library:NIST08.LIB

SI:26 Formula:C25H27Cl2F3N8O CAS:0-00-0 MolWeight:582 RetIndex:4130

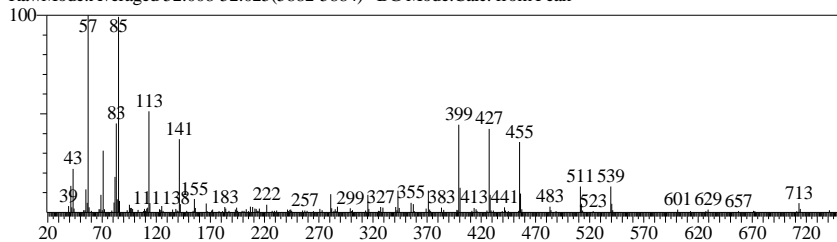
CompName:1-[3,4-Dichlorophenyl]-3-[4-[3-[4-methyl-1-piperazinyl)methyl]-p-anisidino-6-trifluoromethyl]-2-pyri



<< Target >>

Line#:102 R.Time:52.017(Scan#:5883) BasePeak:57.10(53373)

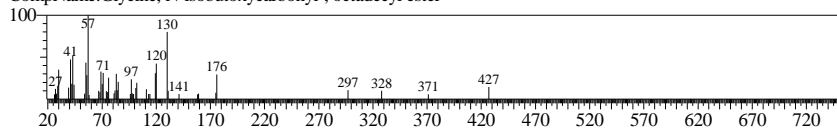
RawMode:Averaged 52.008-52.025(5882-5884) BG Mode:Calc. from Peak



Hit#:1 Entry:174750 Library:NIST08.LIB

SI:49 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975

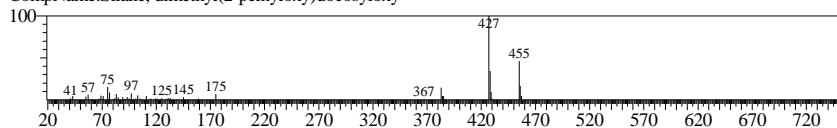
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



Hit#:2 Entry:181721 Library:NIST08.LIB

SI:43 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902

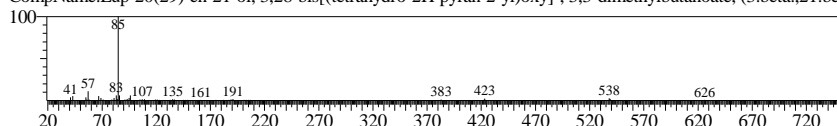
CompName:Silane, dimethyl(2-pentyloxy)docosyloxy-



Hit#:3 Entry:190726 Library:NIST08.LIB

SI:43 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

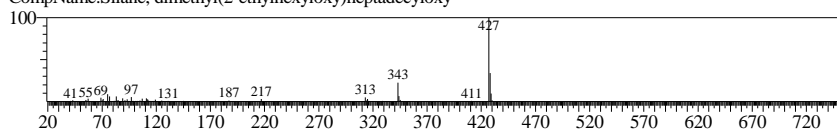
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



Hit#:4 Entry:177733 Library:NIST08.LIB

SI:42 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2703

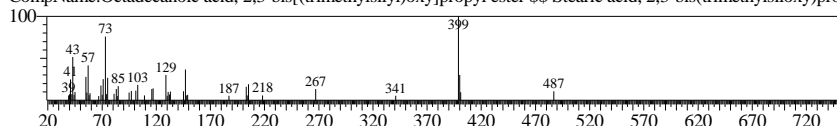
CompName:Silane, dimethyl(2-ethylhexyloxy)heptadecyloxy-



Hit#:5 Entry:28029 Library:NIST08s.LIB

SI:41 Formula:C27H58O4Si2 CAS:1188-75-6 MolWeight:502 RetIndex:2780

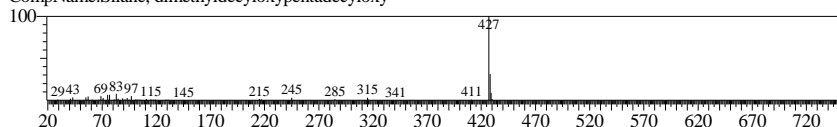
CompName:Octadecanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$\$ Stearic acid, 2,3-bis(trimethylsiloxy)pro



Hit#:6 Entry:177735 Library:NIST08.LIB

SI:40 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767

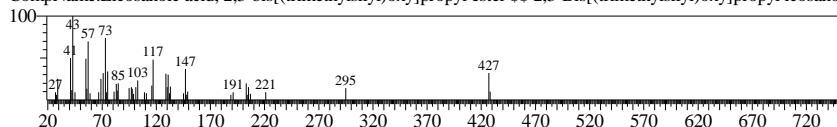
CompName:Silane, dimethyldecyloxyptadecyloxy-



Hit#:7 Entry:186484 Library:NIST08.LIB

SI:40 Formula:C29H62O4Si2 CAS:55517-94-7 MolWeight:530 RetIndex:2979

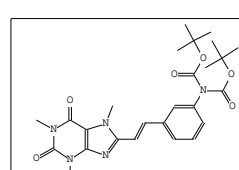
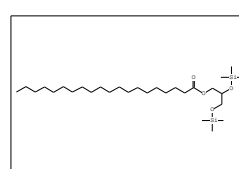
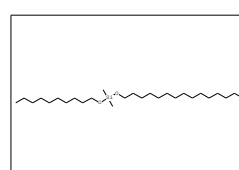
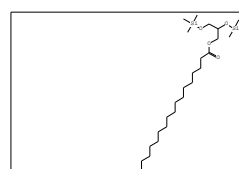
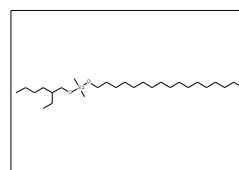
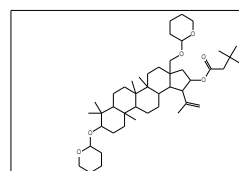
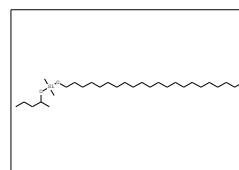
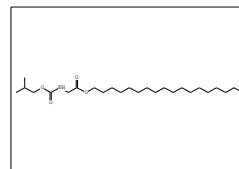
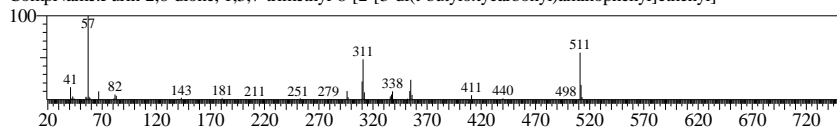
CompName:Eicosanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$\$ 2,3-Bis[(trimethylsilyl)oxy]propyl icosano



Hit#:8 Entry:185431 Library:NIST08.LIB

SI:39 Formula:C26H33N5O6 CAS:0-00-0 MolWeight:511 RetIndex:3851

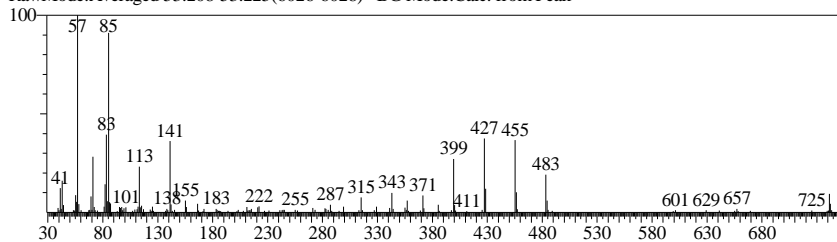
CompName:Purin-2,6-dione, 1,3,7-trimethyl-8-[2-[3-di(t-butyloxycarbonyl)aminophenyl]ethenyl]-



<< Target >>

Line#:103 R.Time:53.217(Scan#:6027) BasePeak:57.10(74299)

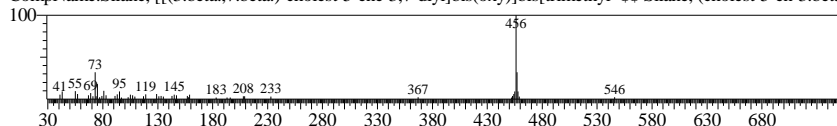
RawMode:Averaged 53.208-53.225(6026-6028) BG Mode:Calc. from Peak



Hit#:1 Entry:28125 Library:NIST08s.LIB

SI:28 Formula:C33H62O2Si2 CAS:33287-26-2 MolWeight:546 RetIndex:2900

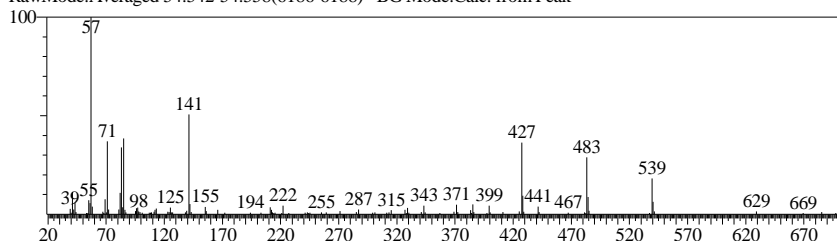
CompName:Silane, [(3.beta.,7.beta.)-cholest-5-ene-3,7-diyl]bis[trimethyl-]bis[trimethyl-] Silane, (cholest-5-en-3.beta.



<< Target >>

Line#:104 R.Time:54.550(Scan#:6187) BasePeak:57.10(195614)

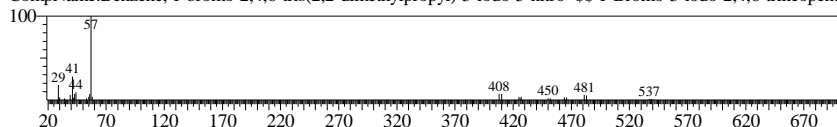
RawMode:Averaged 54.542-54.558(6186-6188) BG Mode:Calc. from Peak



Hit#:1 Entry:186809 Library:NIST08.LIB

SI:47 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079

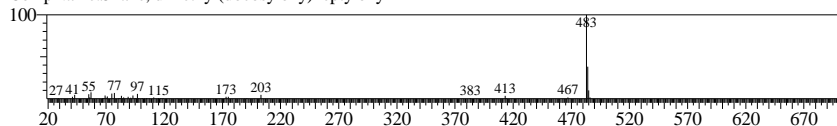
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- 1-Bromo-3-iodo-2,4,6-trineopent



Hit#:2 Entry:184491 Library:NIST08.LIB

SI:45 Formula:C31H66O2Si CAS:0-00-0 MolWeight:498 RetIndex:3164

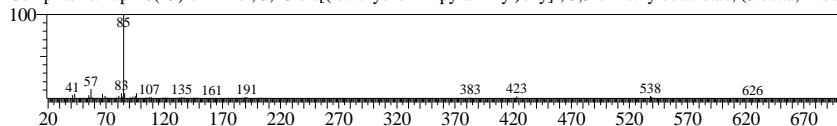
CompName:Silane, dimethyl(docosyloxy)heptyloxy-



Hit#:3 Entry:190726 Library:NIST08.LIB

SI:44 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712

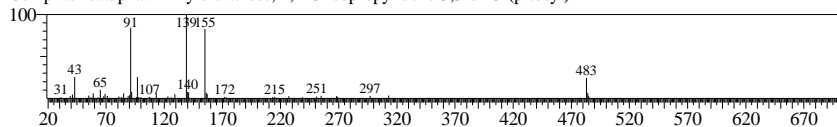
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



Hit#:4 Entry:184445 Library:NIST08.LIB

SI:42 Formula:C22H26O9S2 CAS:0-00-0 MolWeight:498 RetIndex:3723

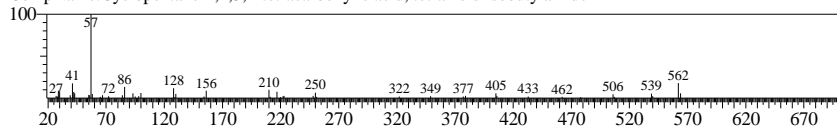
CompName:.alpha.-D-Xylofuranose, 1,2-O-isopropylidene-3,5-di-O-(p-tosyl)-



Hit#:5 Entry:190443 Library:NIST08.LIB

SI:41 Formula:C41H78N4O4 CAS:0-00-0 MolWeight:690 RetIndex:4316

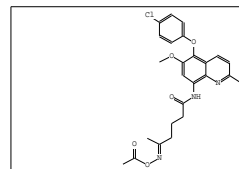
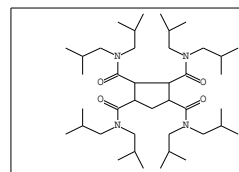
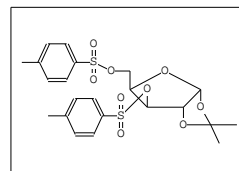
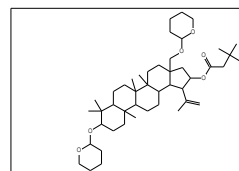
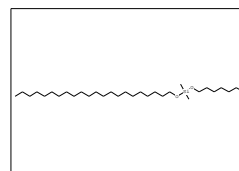
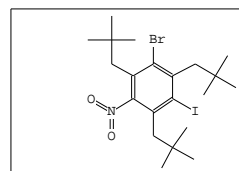
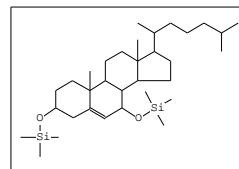
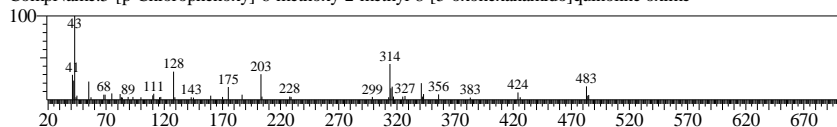
CompName:Cyclopentane-1,2,3,4-tetracarboxylic acid, tetrakis-diisobutylamide



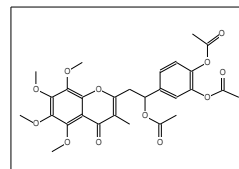
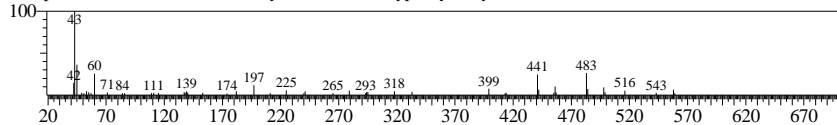
Hit#:6 Entry:183114 Library:NIST08.LIB

SI:31 Formula:C25H26ClN3O5 CAS:0-00-0 MolWeight:483 RetIndex:4040

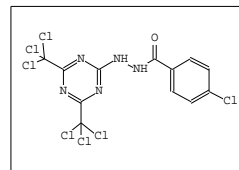
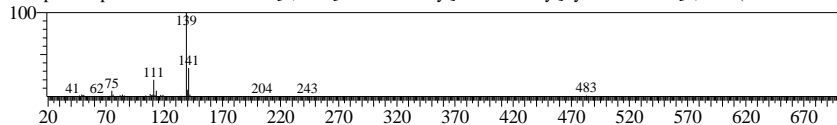
CompName:5-[p-Chlorophenoxy]-6-methoxy-2-methyl-8-[5-oxohexanamido]quinoline oxime



Hit#:7 Entry:187733 Library:NIST08.LIB
SI:30 Formula:C28H30O12 CAS:0-00-0 MolWeight:558 RetIndex:4079
CompName:Chromone, 2-[2-acetoxy-2-[3,5-diacetoxyphenyl]ethyl]-

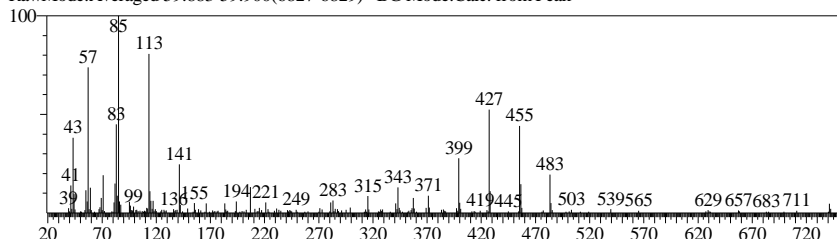


Hit#:8 Entry:182847 Library:NIST08.LIB
SI:28 Formula:C12H6Cl7N5O CAS:110045-61-9 MolWeight:481 RetIndex:3228
CompName:p-Chlorobenzoic acid 2-[4,6-bis(trichloromethyl)-S-triazin-2-yl]hydrazide \$ 2-[4,6-Bis(trichloromethyl)-

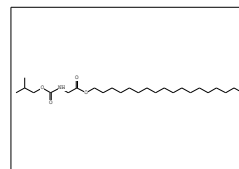
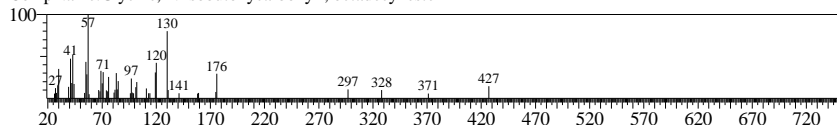


<< Target >>

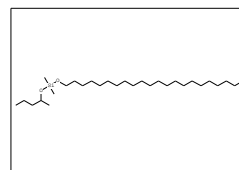
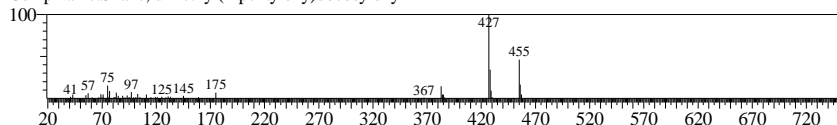
Line#:105 R.Time:59.892(Scan#:6828) BasePeak:85.10(27011)
RawMode:Averaged 59.883-59.900(6827-6829) BG Mode:Calc. from Peak



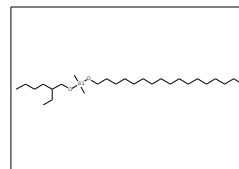
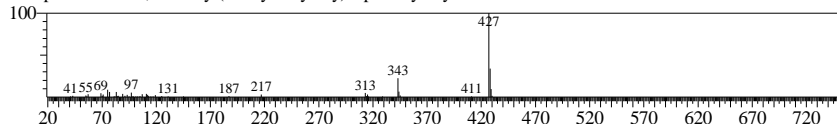
Hit#:1 Entry:174750 Library:NIST08.LIB
SI:48 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



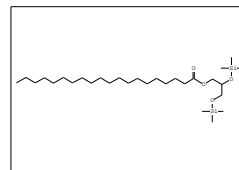
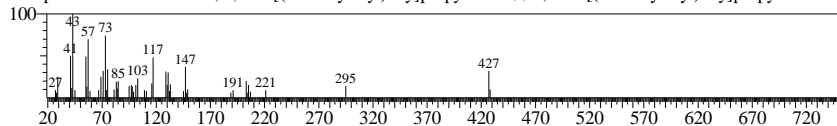
Hit#:2 Entry:181721 Library:NIST08.LIB
SI:45 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902
CompName:Silane, dimethyl(2-pentyloxy)docosyloxy-



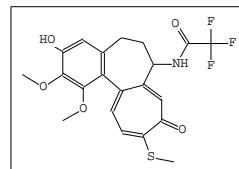
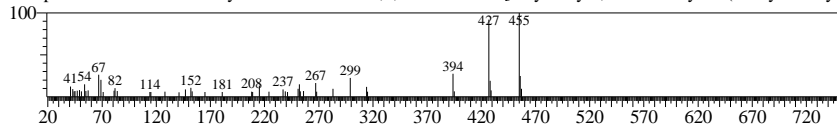
Hit#:3 Entry:177733 Library:NIST08.LIB
SI:44 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2703
CompName:Silane, dimethyl(2-ethylhexyloxy)heptadecyloxy-



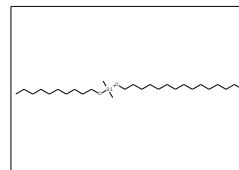
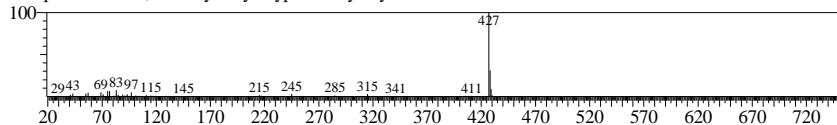
Hit#:4 Entry:186484 Library:NIST08.LIB
SI:43 Formula:C29H62O4Si2 CAS:55517-94-7 MolWeight:530 RetIndex:2979
CompName:Eicosanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$ 2,3-Bis[(trimethylsilyl)oxy]propyl icosano



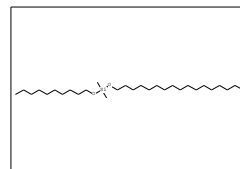
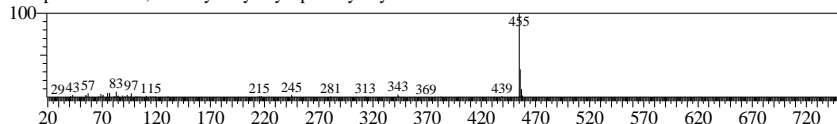
Hit#:5 Entry:179685 Library:NIST08.LIB
SI:42 Formula:C21H20F3NO5S CAS:123643-51-6 MolWeight:455 RetIndex:3297
CompName:N-TFA-3-demethylthicolchicine \$ 2,2,2-Trifluoro-N-[3-hydroxy-1,2-dimethoxy-10-(methylsulfanyl



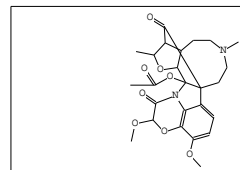
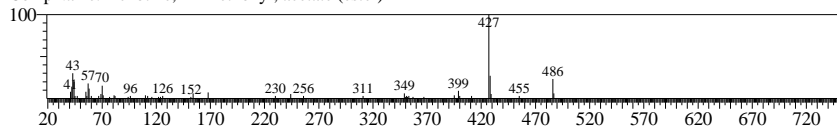
Hit#:6 Entry:177735 Library:NIST08.LIB
SI:40 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2767
CompName:Silane, dimethyldecyloxy pentadecyloxy-



Hit#:7 Entry:181724 Library:NIST08.LIB
SI:39 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966
CompName:Silane, dimethyldecyloxyheptadecyloxy-

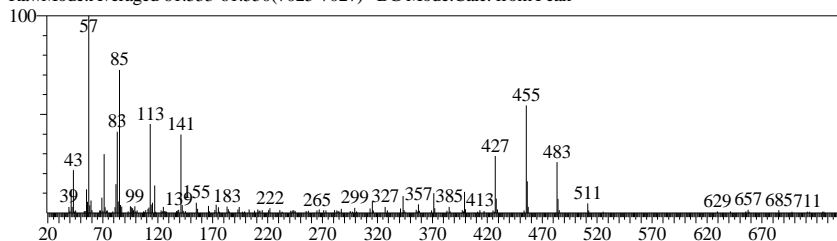


Hit#:8 Entry:183457 Library:NIST08.LIB
SI:39 Formula:C25H30N2O8 CAS:29474-86-0 MolWeight:486 RetIndex:3525
CompName:Dichotine, 11-methoxy-, acetate (ester)

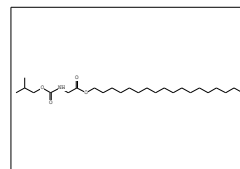
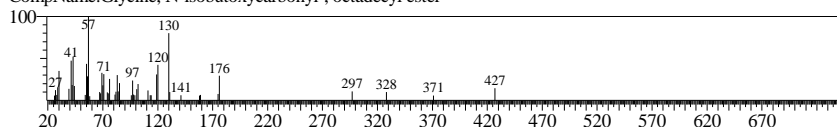


<< Target >>

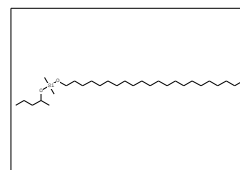
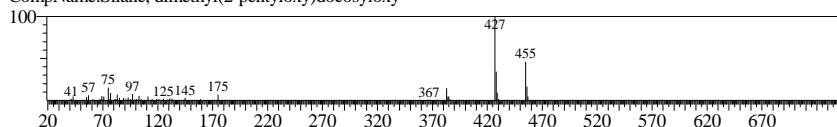
Line#:106 R.Time:61.542(Scan#:7026) BasePeak:57.10(77262)
RawMode:Averaged 61.533-61.550(7025-7027) BG Mode:Calc. from Peak



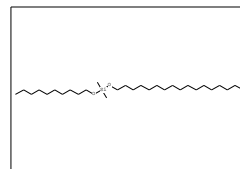
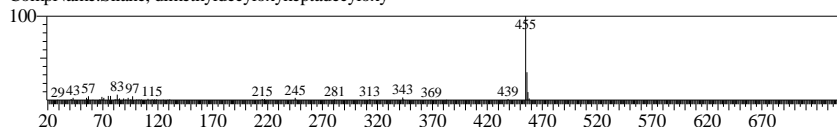
Hit#:1 Entry:174750 Library:NIST08.LIB
SI:50 Formula:C25H49NO4 CAS:0-00-0 MolWeight:427 RetIndex:2975
CompName:Glycine, N-isobutoxycarbonyl-, octadecyl ester



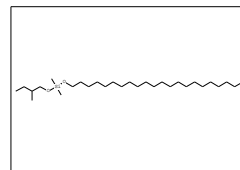
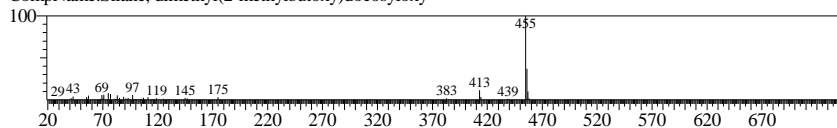
Hit#:2 Entry:181721 Library:NIST08.LIB
SI:48 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902
CompName:Silane, dimethyl(2-pentyloxy)docosyloxy-



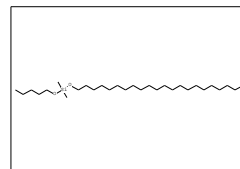
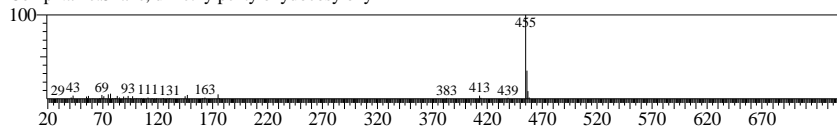
Hit#:3 Entry:181724 Library:NIST08.LIB
SI:46 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966
CompName:Silane, dimethyldecyloxyheptadecyloxy-



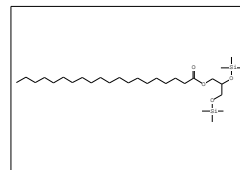
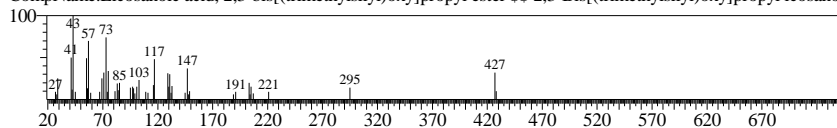
Hit#:4 Entry:181722 Library:NIST08.LIB
SI:45 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902
CompName:Silane, dimethyl(2-methylbutoxy)docosyloxy-



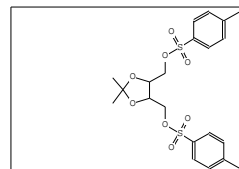
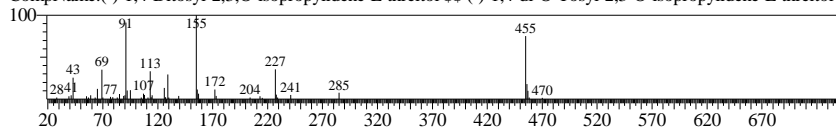
Hit#:5 Entry:181723 Library:NIST08.LIB
SI:44 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966
CompName:Silane, dimethylpentyloxydocosyloxy-



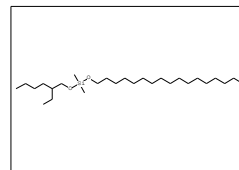
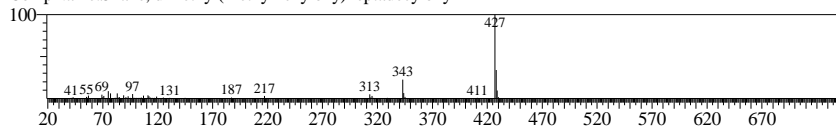
Hit#:6 Entry:186484 Library:NIST08.LIB
SI:43 Formula:C29H62O4Si2 CAS:55517-94-7 MolWeight:530 RetIndex:2979
CompName:Eicosanoic acid, 2,3-bis[(trimethylsilyl)oxy]propyl ester \$ 2,3-Bis[(trimethylsilyl)oxy]propyl icosano



Hit#:7 Entry:181656 Library:NIST08.LIB
SI:42 Formula:C21H26O8S2 CAS:37002-45-2 MolWeight:470 RetIndex:3572
CompName:(-)-1,4-Ditosyl-2,3-O-isopropylidene-L-threitol

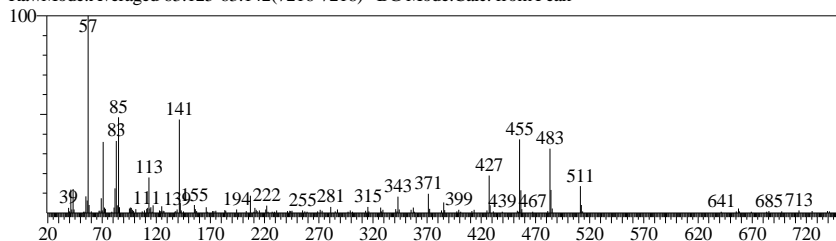


Hit#:8 Entry:177733 Library:NIST08.LIB
SI:42 Formula:C27H58O2Si CAS:0-00-0 MolWeight:442 RetIndex:2703
CompName:Silane, dimethyl(2-ethylhexyloxy)heptadecyloxy-

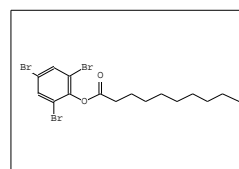
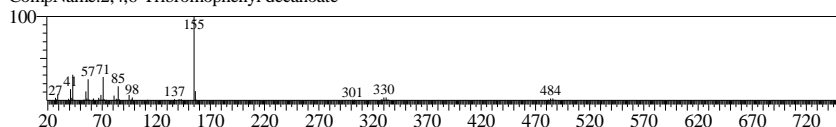


<< Target >>

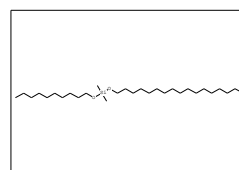
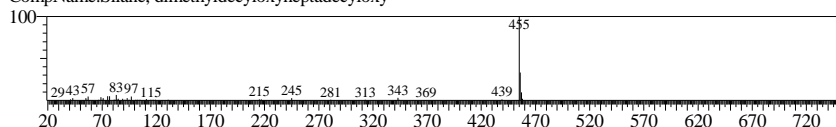
Line#:107 R.Time:63.133(Scan#:7217) BasePeak:57.10(117629)
RawMode:Averaged 63.125-63.142(7216-7218) BG Mode:Calc. from Peak



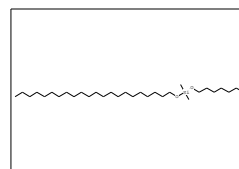
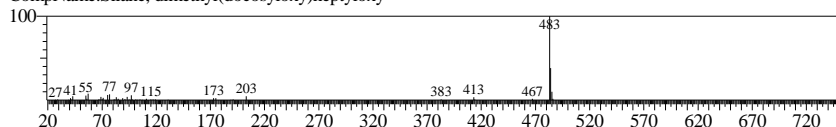
Hit#:1 Entry:182936 Library:NIST08.LIB
SI:49 Formula:C16H21Br3O2 CAS:0-00-0 MolWeight:482 RetIndex:2815
CompName:2,4,6-Tribromophenyl decanoate



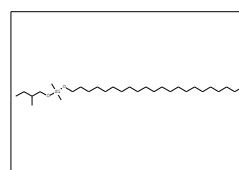
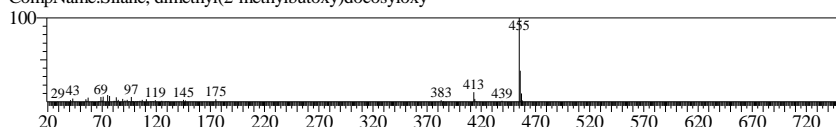
Hit#:2 Entry:181724 Library:NIST08.LIB
SI:44 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966
CompName:Silane, dimethyldecyloxyheptadecyloxy-



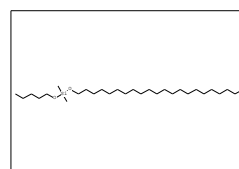
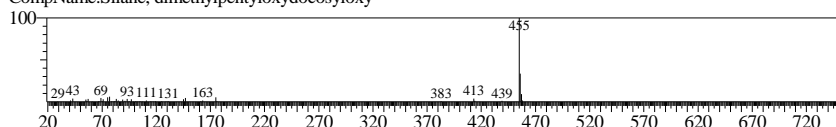
Hit#:3 Entry:184491 Library:NIST08.LIB
SI:43 Formula:C31H66O2Si CAS:0-00-0 MolWeight:498 RetIndex:3164
CompName:Silane, dimethyl(docosyloxy)heptyloxy-



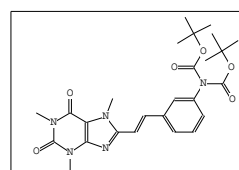
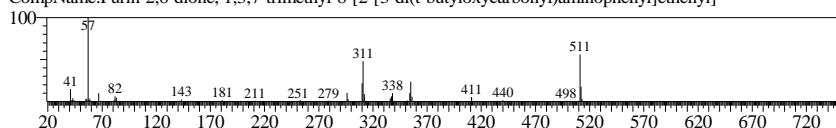
Hit#:4 Entry:181722 Library:NIST08.LIB
SI:43 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2902
CompName:Silane, dimethyl(2-methylbutoxy)docosyloxy-



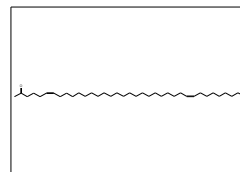
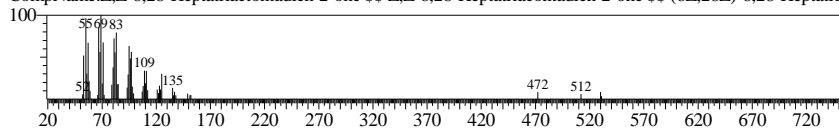
Hit#:5 Entry:181723 Library:NIST08.LIB
SI:41 Formula:C29H62O2Si CAS:0-00-0 MolWeight:470 RetIndex:2966
CompName:Silane, dimethylpentyloxydocosyloxy-



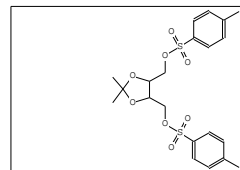
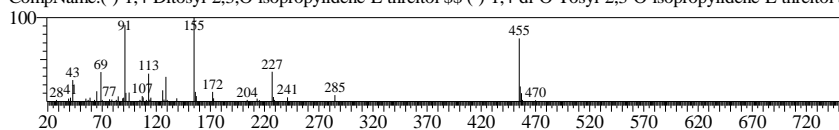
Hit#:6 Entry:185431 Library:NIST08.LIB
SI:41 Formula:C26H33N5O6 CAS:0-00-0 MolWeight:511 RetIndex:3851
CompName:Purin-2,6-dione, 1,3,7-trimethyl-8-[2-[3-di(t-butyloxycarbonyl)aminophenyl]ethenyl]-



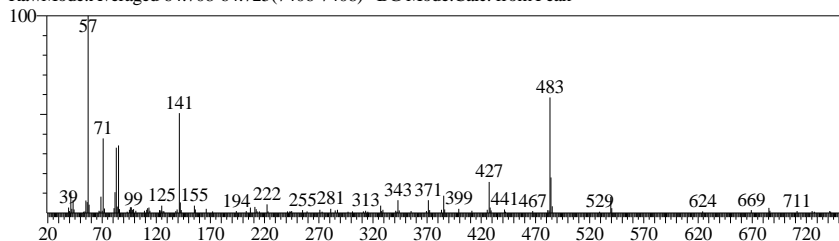
Hit#:7 Entry:186505 Library:NIST08.LIB
SI:39 Formula:C37H70O CAS:133530-21-9 MolWeight:530 RetIndex:3851
CompName:Z,Z-6,28-Heptatriacontadien-2-one \$Z,Z-6,28-Heptatriacontadien-2-one \$ (6Z,28Z)-6,28-Heptatri



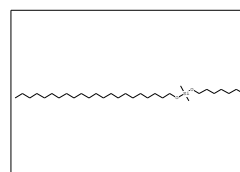
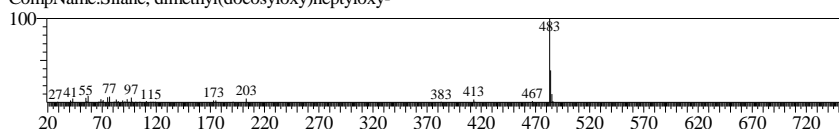
Hit#:8 Entry:181656 Library:NIST08.LIB
SI:36 Formula:C21H26O8S2 CAS:37002-45-2 MolWeight:470 RetIndex:3572
CompName:(-)-1,4-Ditosyl-2,3-O-isopropylidene-L-threitol \$(-)-1,4-di-O-Tosyl-2,3-O-isopropylidene-L-threitol :



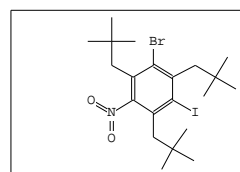
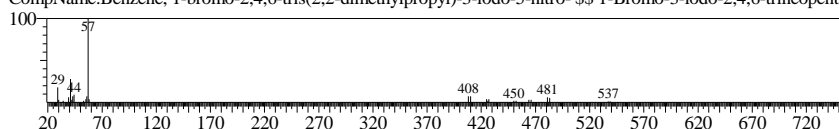
<< Target >>
Line#:108 R.Time:64.717(Scan#:7407) BasePeak:57.10(188963)
RawMode:Averaged 64.708-64.725(7406-7408) BG Mode:Calc. from Peak



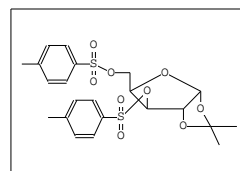
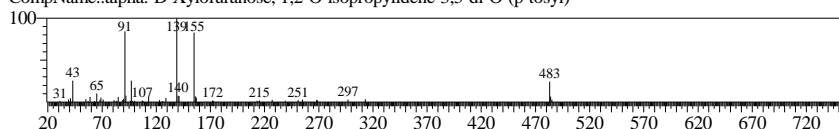
Hit#:1 Entry:184491 Library:NIST08.LIB
SI:52 Formula:C31H66O2Si CAS:0-00-0 MolWeight:498 RetIndex:3164
CompName:Silane, dimethyl(docosyloxy)heptyloxy-



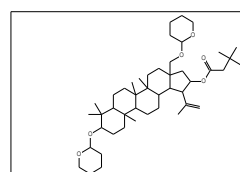
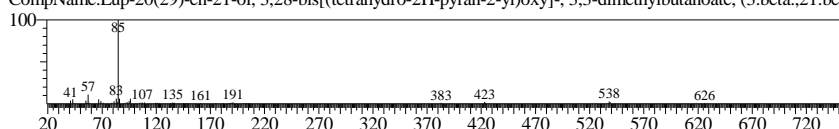
Hit#:2 Entry:186809 Library:NIST08.LIB
SI:48 Formula:C21H33BrINO2 CAS:40572-26-7 MolWeight:537 RetIndex:3079
CompName:Benzene, 1-bromo-2,4,6-tris(2,2-dimethylpropyl)-3-iodo-5-nitro- \$1-Bromo-3-iodo-2,4,6-trineopent



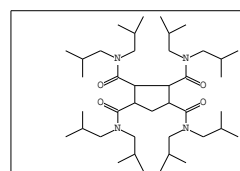
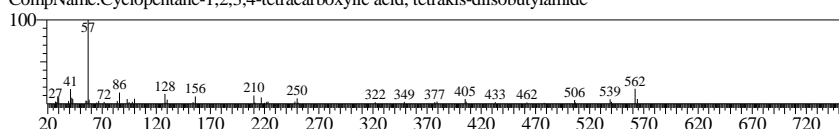
Hit#:3 Entry:184445 Library:NIST08.LIB
SI:44 Formula:C22H26O9S2 CAS:0-00-0 MolWeight:498 RetIndex:3723
CompName:.alpha.-D-Xylofuranose, 1,2-O-isopropylidene-3,5-di-O-(p-tosyl)-



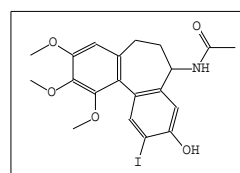
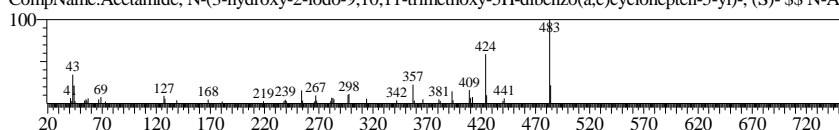
Hit#:4 Entry:190726 Library:NIST08.LIB
SI:41 Formula:C46H76O6 CAS:55401-92-8 MolWeight:724 RetIndex:4712
CompName:Lup-20(29)-en-21-ol, 3,28-bis[(tetrahydro-2H-pyran-2-yl)oxy]-, 3,3-dimethylbutanoate, (3.beta.,21.be



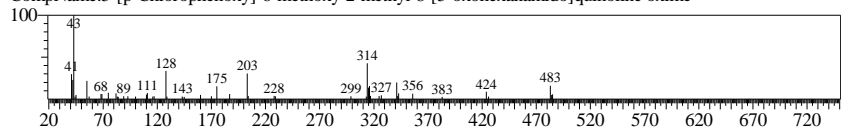
Hit#:5 Entry:190443 Library:NIST08.LIB
SI:40 Formula:C41H78N4O4 CAS:0-00-0 MolWeight:690 RetIndex:4316
CompName:Cyclopentane-1,2,3,4-tetracarboxylic acid, tetrakis-diisobutylamide



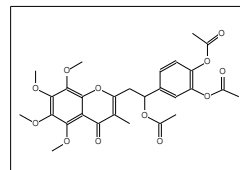
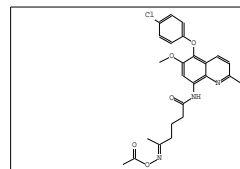
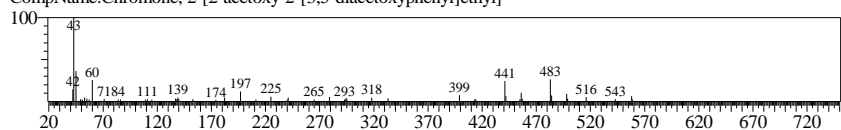
Hit#:6 Entry:183100 Library:NIST08.LIB
SI:33 Formula:C20H22INO5 CAS:38838-27-6 MolWeight:483 RetIndex:3486
CompName:Acetamide, N-(3-hydroxy-2-iodo-9,10,11-trimethoxy-5H-dibenzo(a,c)cyclohepten-5-yl)-, (S)- \$N-A



Hit#:7 Entry:183114 Library:NIST08.LIB
SI:33 Formula:C25H26ClN3O5 CAS:0-00-0 MolWeight:483 RetIndex:4040
CompName:5-[p-Chlorophenoxy]-6-methoxy-2-methyl-8-[5-oxohexanamido]quinoline oxime



Hit#:8 Entry:187733 Library:NIST08.LIB
SI:31 Formula:C28H30O12 CAS:0-00-0 MolWeight:558 RetIndex:4079
CompName:Chromone, 2-[2-acetoxy-2-[3,5-diacetoxyphenyl]ethyl]-



Method

[Comment]

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

[AOC-20i+s]

of Rinses with Presolvent :4
 # of Rinses with Solvent(post) :4
 # of Rinses with Sample :0
 Plunger Speed(Suction) :Middle
 Viscosity Comp. Time :0.2 sec
 Plunger Speed(Injection) :High
 Syringe Insertion Speed :High
 Injection Mode :Normal
 Pumping Times :0
 Inj. Port Dwell Time :0.0 sec
 Terminal Air Gap :No
 Plunger Washing Speed :High
 Washing Volume :6uL
 Syringe Suction Position :0.0 mm
 Syringe Injection Position :0.0 mm
 Solvent Selection :All A,B,C

[GC-2010]COLUMN: ZEBRON ZB5ms 30meter x 0.25mm I.D x 0.25µm film thickness

Column Oven Temp. :50.0 °C
 Injection Temp. :240.00 °C
 Injection Mode :Split
 Flow Control Mode :Linear Velocity
 Pressure :53.5 kPa
 Total Flow :14.0 mL/min
 Column Flow :1.00 mL/min
 Linear Velocity :36.3 cm/sec
 Purge Flow :3.0 mL/min
 Split Ratio :10.0
 High Pressure Injection :OFF
 Carrier Gas Saver :OFF
 Splitter Hold :OFF
 Oven Temp. Program

Rate	Temperature(°C)	Hold Time(min)
-	50.0	3.00
10.00	320.0	40.00

 Purge Flow Program

Rate	Flow(mL/min)	Hold Time(min)
-	3.0	3.00

< Ready Check Heat Unit >

Column Oven : Yes
 SPL1 : Yes
 MS : Yes

< Ready Check Detector(FTD) >

< Ready Check Baseline Drift >

< Ready Check Injection Flow >

SPL1 Carrier : Yes
 SPL1 Purge : Yes

< Ready Check APC Flow >

< Ready Check Detector APC Flow >

External Wait :No
 Equilibrium Time :0.5 min

[GC Program]

[GCMS-QP2010 Plus]

IonSourceTemp :240.00 °C
 Interface Temp. :320.00 °C
 Solvent Cut Time :2.50 min
 Detector Gain Mode :Absolute
 Detector Gain :1.00 kV
 Threshold :100

[MS Table]

--Group 1 - Event 1--

Start Time :3.00min
 End Time :70.00min
 ACQ Mode :Scan
 Event Time :0.50sec
 Scan Speed :1666
 Start m/z :35.00
 End m/z :750.00

Sample Inlet Unit :GC