

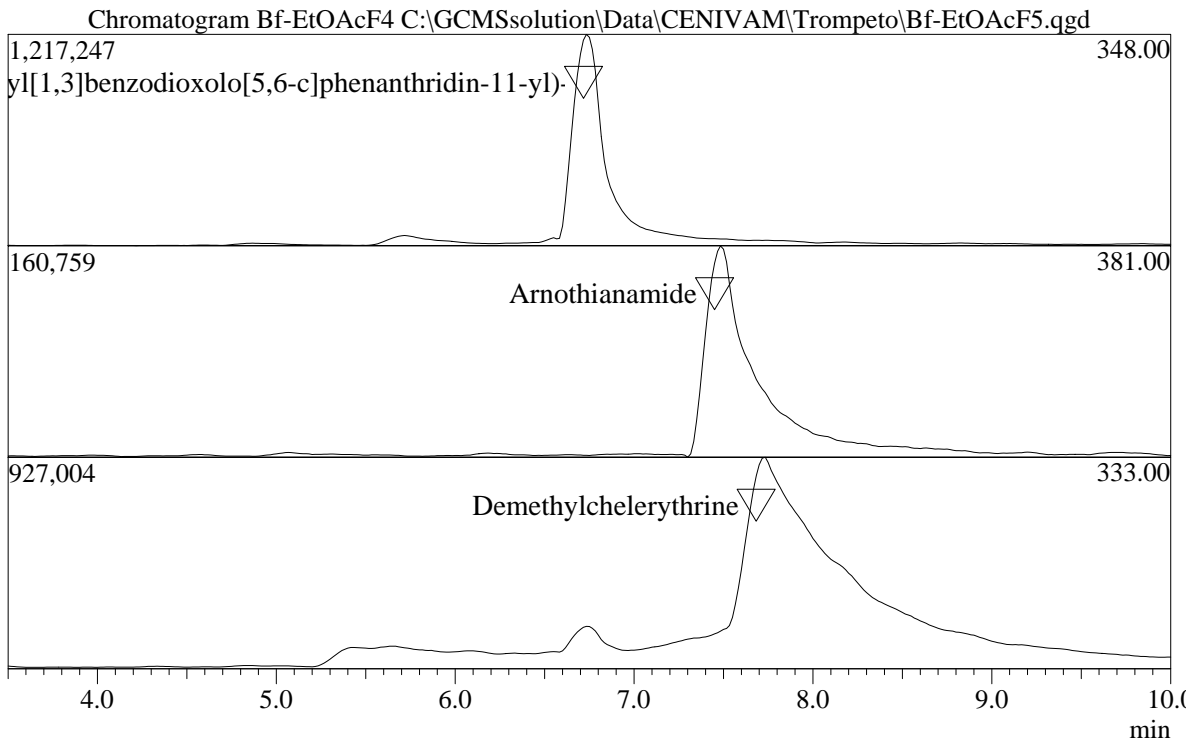


LABORATORIO DE CALIDAD DE PRODUCTOS NATURALES

CERTICADO DE CALIDAD/QUALITY CERTIFICATE

Reporte/Report : Bf-EtOAcF4
 Fecha/Date : 3/10/2007 3:36:32 PM

Identificación de la muestra/Sample ID : Bf-EtOAcF4
 Vol. Inyección/Injection Vol.: 2.000
 Archivo/Data File : C:\GCMSsolution\Data\CENIVAM\Trompeto\Bf-EtOAcF5.qgd
 Método/Method File : C:\GCMSsolution\Data\CENIVAM\linas\Justi\ESTEROIDESLINAS.qgm
 Tuning File : C:\GCMSsolution\System\Tune1\CPN-SEI10260CLOSE20070222.qgt
 Descripción muestra/Sample Descr. :



Quantitative Result Table

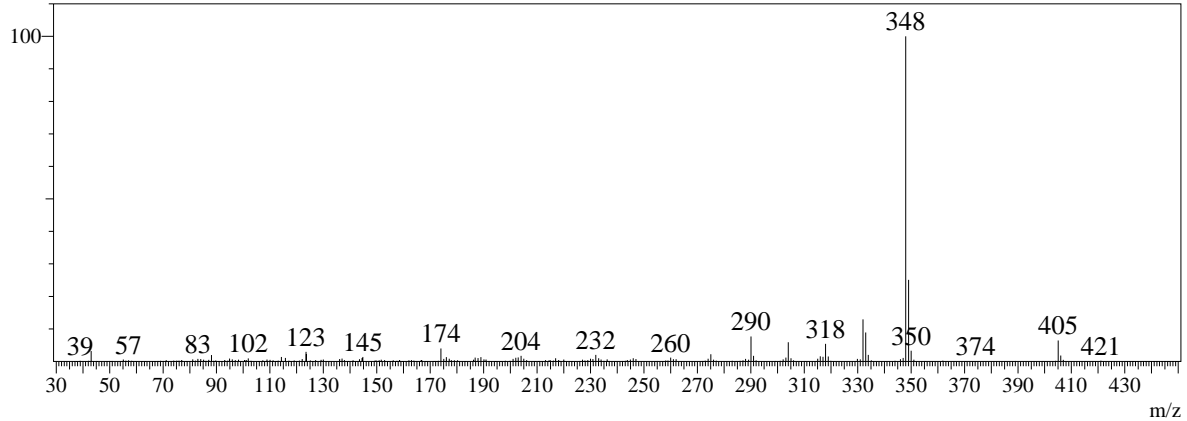
ID#	R.Time	Name	m/z	Area	Conc.	ilarity
1	6.736	2-Propanone, 1-(12,13-dihydro-1,2-dimethoxy-12-methyl[1,3]ben	348	14580402	41.5	96
2	7.485	Arnothianamide	381	2630067	7.5	80
3	7.727	Demethylchelerythrine	333	17927372	51.0	89



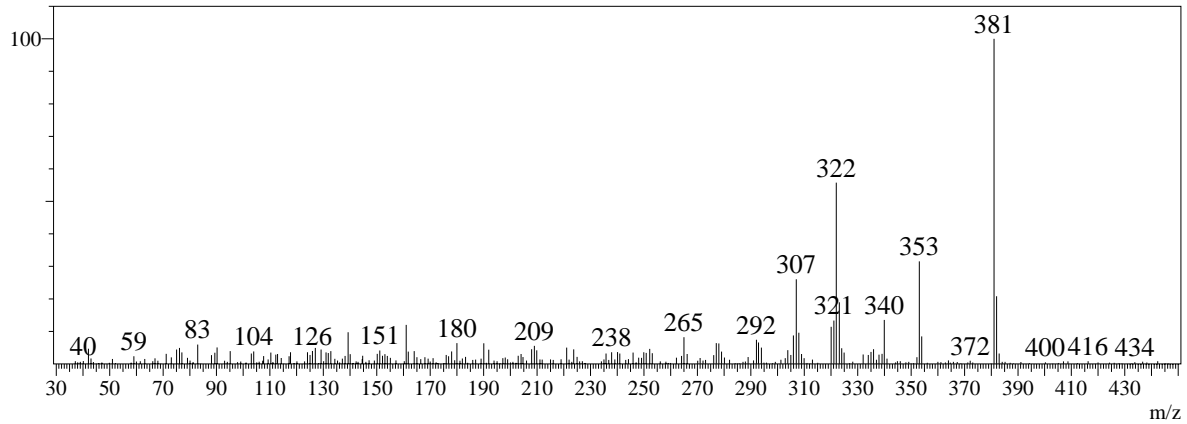
LABORATORIO DE CALIDAD DE PRODUCTOS NATURALES

Spectrum

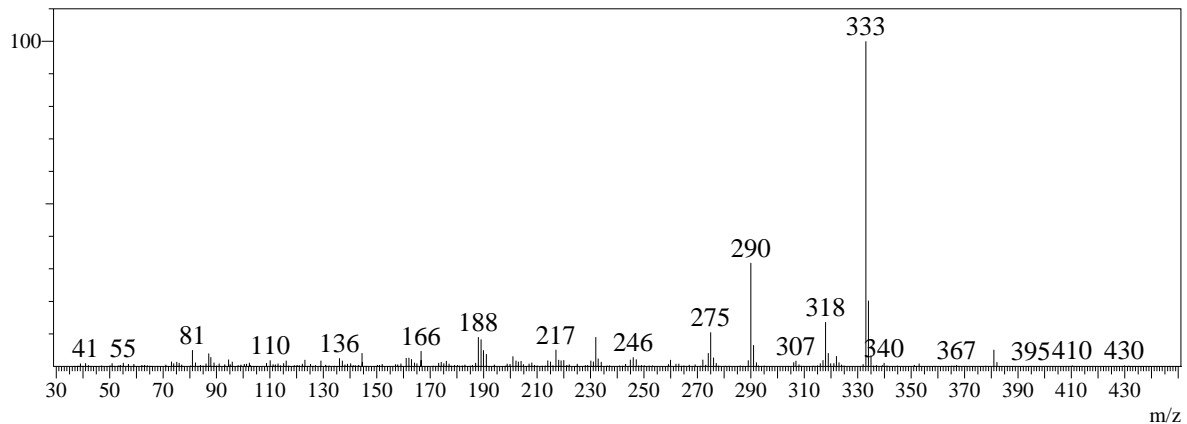
ID#:1 R.Time:6.7(Scan#:195)
MassPeaks:408
RawMode:Averaged 6.7-6.8(194-196)
BG Mode:Calc. from Peak



ID#:2 R.Time:7.5(Scan#:240)
MassPeaks:399
RawMode:Averaged 7.5-7.5(239-241)
BG Mode:Calc. from Peak



ID#:3 R.Time:7.7(Scan#:255)
MassPeaks:429
RawMode:Averaged 7.7-7.8(254-256)
BG Mode:Calc. from Peak



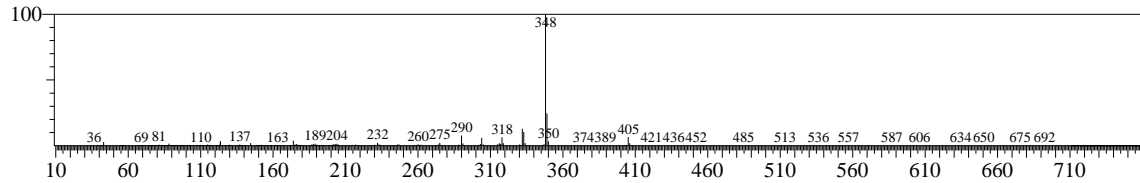


LABORATORIO DE CALIDAD DE PRODUCTOS NATURALES

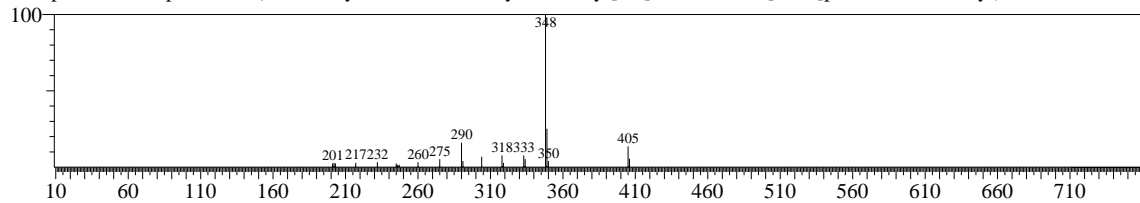
Library

<< Target >>

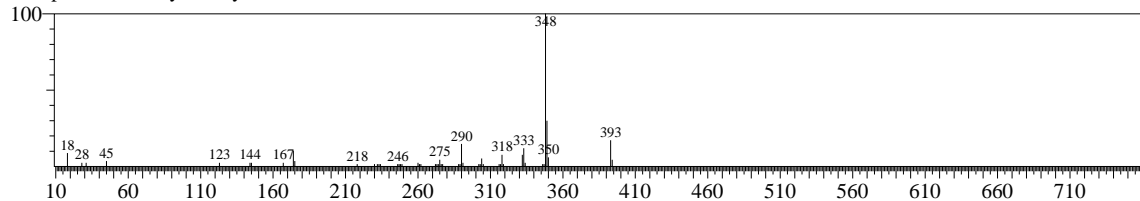
Line#:1 R.Time:6.717(Scan#:194) MassPeaks:408
RawMode:Averaged 6.700-6.733(193-195) BasePeak:348.00(2111105)
BG Mode:Calc. from Peak



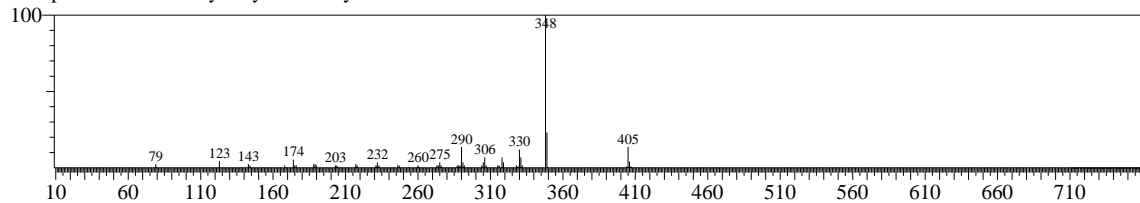
Hit#:1 Entry:383031 Library:WILEY7N2.LIB
SI:83 Formula:C₂₄H₂₃NO₅ CAS:56051-44-6 MolWeight:405 RetIndex:0
CompName:2-Propanone, 1-(12,13-dihydro-1,2-dimethoxy-12-methyl[1,3]benzodioxolo[5,6-c]phenanthridin-11-yl)-



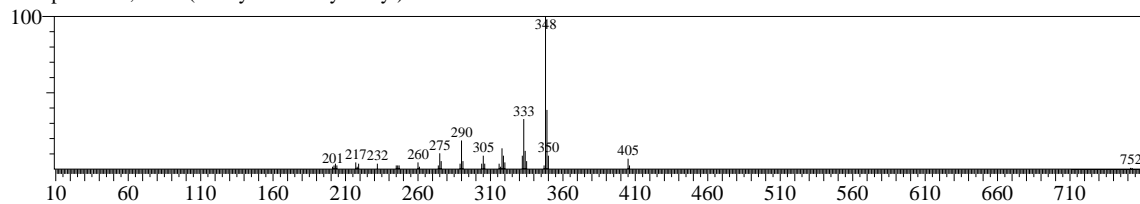
Hit#:2 Entry:375419 Library:WILEY7N2.LIB
SI:80 Formula:C₂₃H₂₃NO₅ CAS:0-0-0 MolWeight:393 RetIndex:0
CompName:Ethoxychelerythrine



Hit#:3 Entry:383030 Library:WILEY7N2.LIB
SI:79 Formula:C₂₄H₂₃NO₅ CAS:56051-44-6 MolWeight:405 RetIndex:0
CompName:11-Acetyl-dihydrochelerythrine

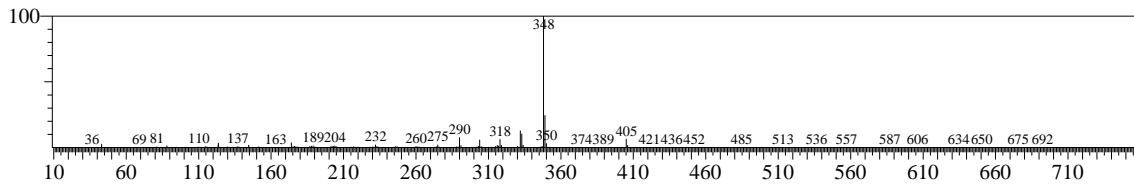


Hit#:4 Entry:439162 Library:WILEY7N2.LIB
SI:78 Formula:C₄₅H₄₀N₂O₉ CAS:0-0-0 MolWeight:752 RetIndex:0
CompName:1,3-Bis-(11-Hydrochelerythrinyl)-Acetone

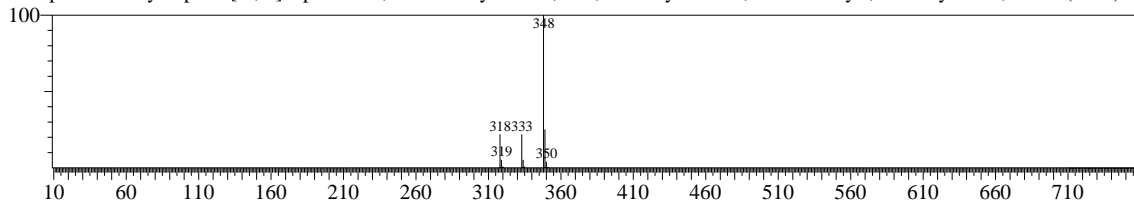


<< Target >>

Line#:1 R.Time:6.717(Scan#:194) MassPeaks:408
RawMode:Averaged 6.700-6.733(193-195) BasePeak:348.00(2111105)
BG Mode:Calc. from Peak

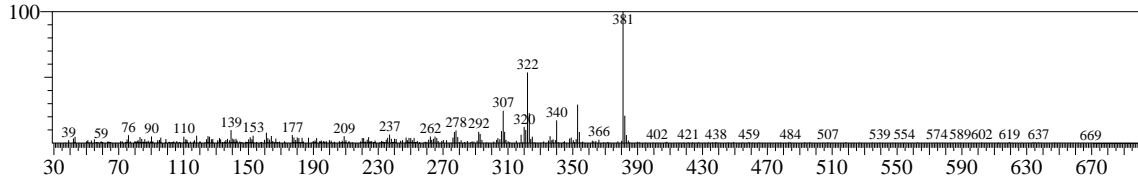


Hit#:5 Entry:338479 Library:WILEY7N2.LIB
SI:75 Formula:C₂₂H₂₀O₄ CAS:102521-7-3 MolWeight:348 RetIndex:0
CompName:Dicyclopenta[ef,kl]heptalene-1,2-dicarboxylic acid, 10b,10c-dihydro-10b,10c-dimethyl-, dimethyl ester, trans- (CAS)

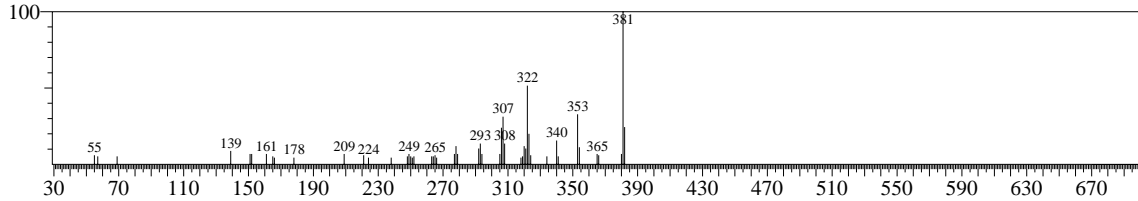


<< Target >>

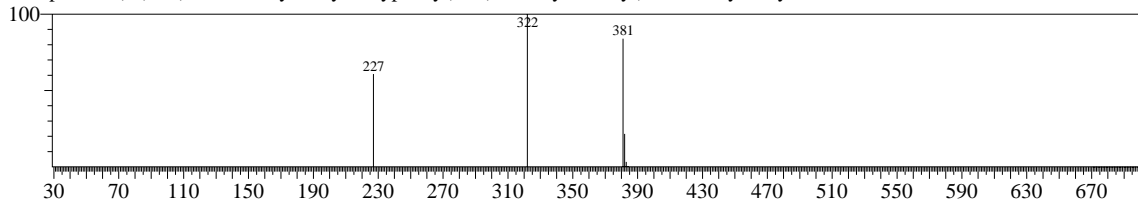
Line#:2 R.Time:7.450(Scan#:238) MassPeaks:396
RawMode:Averaged 7.433-7.467(237-239) BasePeak:381.00(200000)
BG Mode:Calc. from Peak



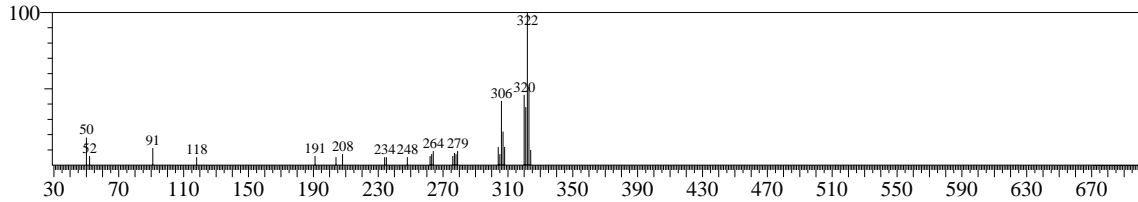
Hit#:1 Entry:366749 Library:WILEY7N2.LIB
SI:74 Formula:C21H19NO6 CAS:0-0-0 MolWeight:381 RetIndex:0
CompName:Arnothianamide



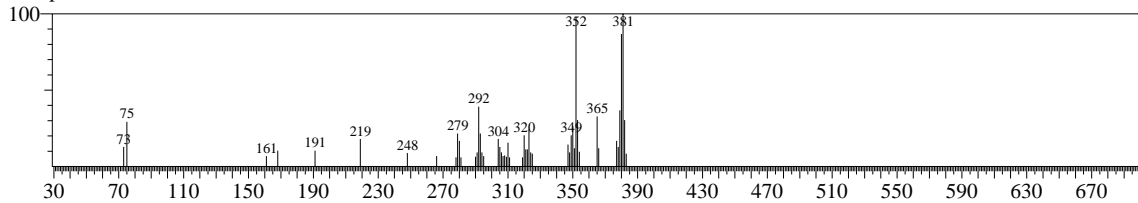
Hit#:2 Entry:366798 Library:WILEY7N2.LIB
SI:45 Formula:C22H23NO5 CAS:0-0-0 MolWeight:381 RetIndex:0
CompName:(+)-1-(3,5-Dimethyl-4-hydroxyphenyl)-9-(methoxycarbonyl)-1,2,3,4-hydroxycarbazole



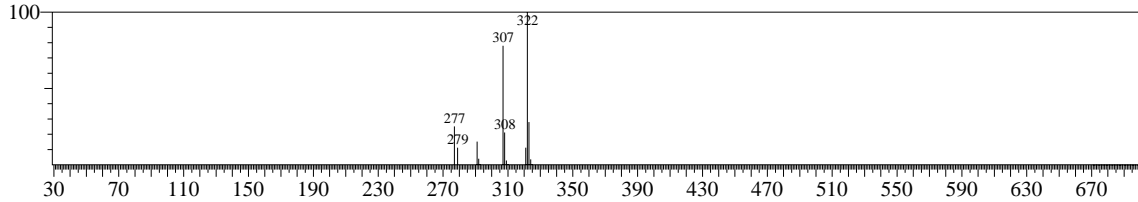
Hit#:3 Entry:346615 Library:WILEY7N2.LIB
SI:44 Formula:C20H20ClNO3 CAS:19717-36-3 MolWeight:357 RetIndex:0
CompName:Berbium, 7,8,13,13a-tetrahydro-2,3,11-trimethoxy-, chloride (CAS)



Hit#:4 Entry:366844 Library:WILEY7N2.LIB
SI:44 Formula:C23H31NO2Si CAS:73885-59-3 MolWeight:381 RetIndex:0
CompName:TMS DERIVATIVE OF N-N-PROPYLNORAPOCODEINE

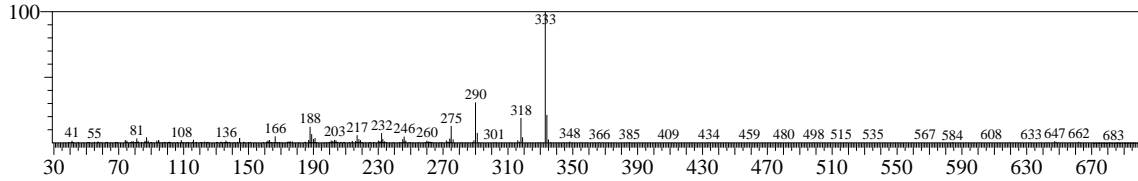


Hit#:5 Entry:310397 Library:WILEY7N2.LIB
SI:43 Formula:C25H22 CAS:76804-52-9 MolWeight:322 RetIndex:0
CompName:9-(2,6-Bis(1-methylethenyl)phenyl)fluorene \$ 9H-Fluorene, 9-[2,6-bis(1-methylethenyl)phenyl]- (CAS)

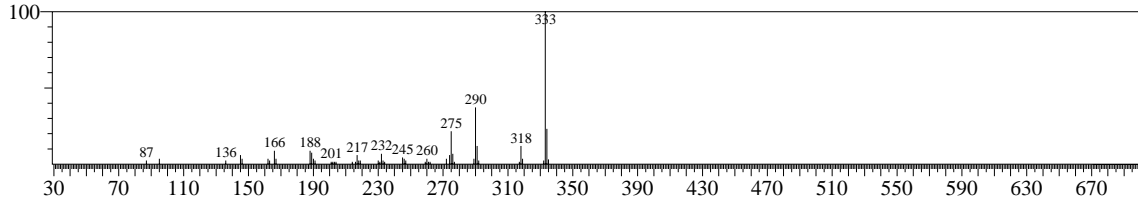


<< Target >>

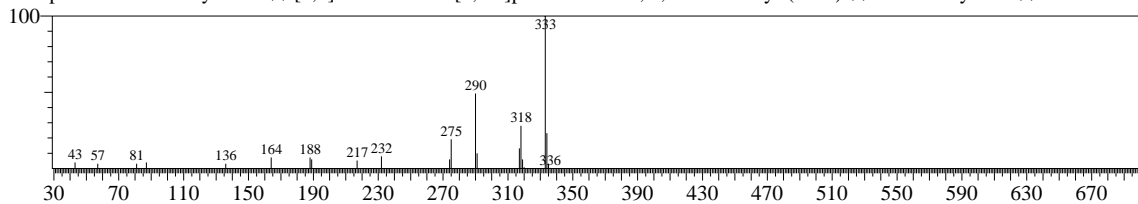
Line#3 R.Time:7.683(Scan#:252) MassPeaks:371
RawMode:Averaged 7.667-7.700(251-253) BasePeak:333.00(781015)
BG Mode:Calc. from Peak



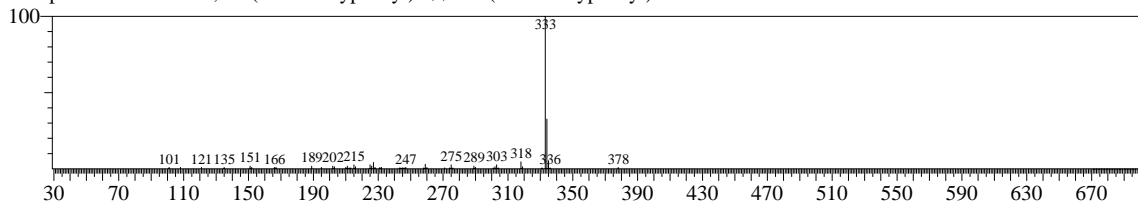
Hit#1 Entry:322689 Library:WILEY7N2.LIB
SI:87 Formula:C20H15NO4 CAS:0-0-0 MolWeight:333 RetIndex:0
CompName:Demethylchelerythrine



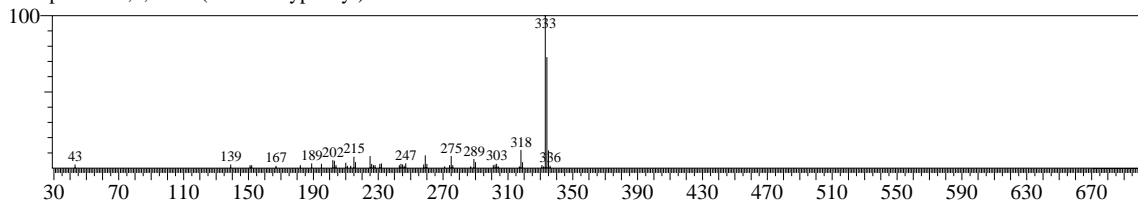
Hit#2 Entry:322688 Library:WILEY7N2.LIB
SI:79 Formula:C20H15NO4 CAS:6900-99-8 MolWeight:333 RetIndex:0
CompName:Norchelerythrine \$\$ [1,3]Benzodioxolo[5,6-c]phenanthridine, 1,2-dimethoxy- (CAS) \$ Norchelerythrin \$ N-Norchele



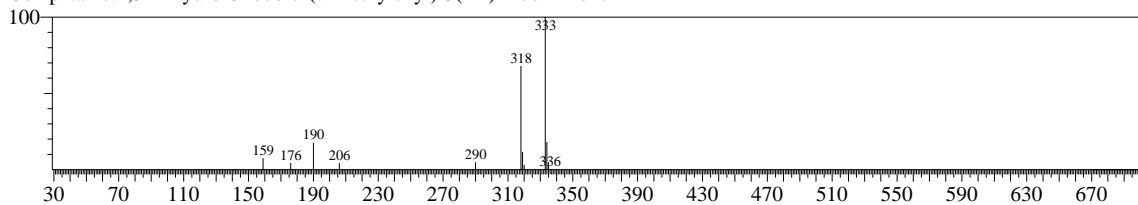
Hit#3 Entry:364909 Library:WILEY7N2.LIB
SI:65 Formula:C23H22O5 CAS:95431-48-4 MolWeight:378 RetIndex:0
CompName:Acetic acid, tris(4-methoxyphenyl)- \$ Tris(4-methoxyphenyl)acetic acid #



Hit#4 Entry:363542 Library:WILEY7N2.LIB
SI:63 Formula:C24H24O4 CAS:0-0-0 MolWeight:376 RetIndex:0
CompName:1,1,1-Tris(4-methoxyphenyl)acetone



Hit#5 Entry:322210 Library:WILEY7N2.LIB
SI:62 Formula:C11H16INOSi CAS:88761-39-1 MolWeight:333 RetIndex:0
CompName:2,3-Dihydro-6-iodo-7-(trimethylsilyl)-5(1H)-indolizinone



Method

[Comment]

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

[AOC-20i+s]

of Rinses with Presolvent :3
 # of Rinses with Solvent(post) :5
 # of Rinses with Sample :3
 Plunger Speed(Suction) :Middle
 Viscosity Comp. Time :0.0 sec
 Plunger Speed(Injection) :High
 Syringe Insertion Speed :High
 Injection Mode :Normal
 Pumping Times :5
 Inj. Port Dwell Time :0.0 sec
 Terminal Air Gap :No
 Plunger Washing Speed :High
 Washing Volume :8uL
 Syringe Suction Position :0.0 mm
 Syringe Injection Position :0.0 mm
 Solvent Selection :All A,B,C

[GC-2010]

Column Oven Temp. :275.0 °C
 Injection Temp. :310.00 °C
 Injection Mode :Split
 Flow Control Mode :Linear Velocity
 Pressure :139.5 kPa
 Total Flow :6.0 mL/min
 Column Flow :1.00 mL/min
 Linear Velocity :39.6 cm/sec
 Purge Flow :1.0 mL/min
 Split Ratio :4.0
 High Pressure Injection :OFF
 Carrier Gas Saver :OFF
 Splitter Hold :OFF
 Oven Temp. Program

Rate	Temperature(°C)	Hold Time(min)
-	275.0	2.00
8.0	320.0	10.00

< 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 :Yes
 Equilibrium Time :0.1 min

[GC Program]

[GCMS-QP2010]

IonSourceTemp :260.00 °C
 Interface Temp. :280.00 °C
 Solvent Cut Time :2.50 min
 Ionization Mode :SEI
 Detector Gain Mode :Relative
 Detector Gain :1.00 kV
 Threshold :5

[MS Table]

Group : 1
 Start Time :3.50min
 End Time :17.00min
 ACQ Mode :Scan
 Interval :1.00sec
 Scan Speed :714
 Start m/z :35.00
 End m/z :700.00

Sample Inlet Unit :GC

[MS Program]

Use MS Program :OFF