

Supplementary Materials S2: All chemical structures of fungal meroterpenoids newly discovered in 2009-2019

The Chemistry and Biology of Fungal Meroterpenoids (2009-2019)

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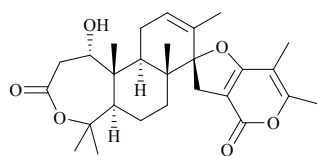
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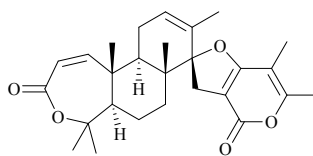
CONTENTS

1. Triketide-terpenoids (1-98)	2
2. Tetraketide-terpenoids (99-652).....	6
2.1 Derived from Orsellinic acid/orcinaldehyde (99-323)	6
2.2 Derived from 3,5-Dimethylorsellinic acid (DMOA) (324-492).....	14
2.3 Derived from 6-Methylsalicylic acid (6-MSA) (493-580)	20
2.4 Other tetraketide-terpenoids (581-652)	24
3. Other polyketide-terpenoids (653-730)	27
4. Indole-terpenoids (731-844).....	30
5. Shikimate-terpenoids (845-1544).....	35
6. Miscellaneous meroterpenoids (1545-1585).....	56

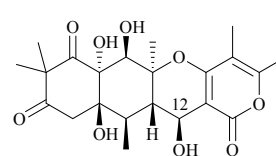
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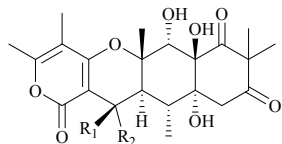
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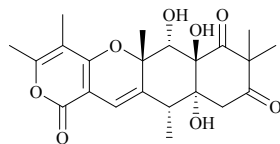
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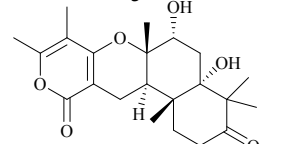
4 $R_1 = \text{OCH}_3$, $R_2 = \text{H}$

5 $R_1 = \text{OH}$, $R_2 = \text{H}$

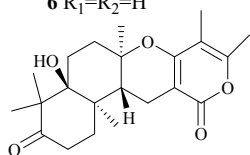
6 $R_1 = R_2 = \text{H}$



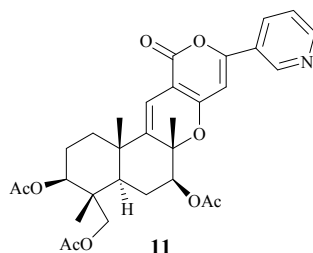
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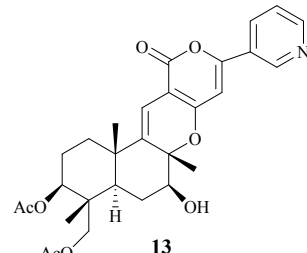
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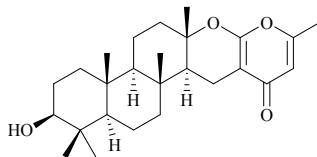
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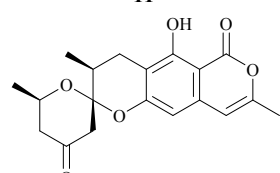
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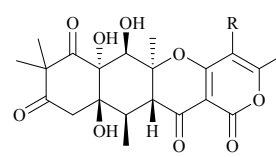
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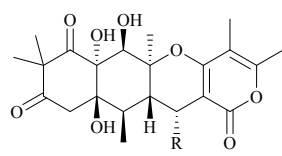


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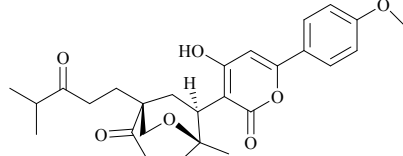
14 $R = \text{CH}_3$

16 $R = \text{H}$

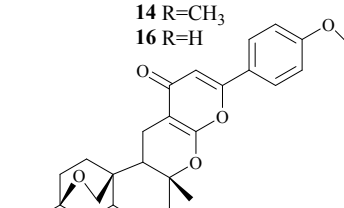


15 $R = \text{H}$

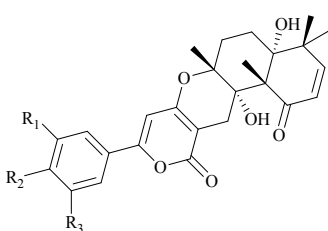
17 $R = \text{OH}$



18

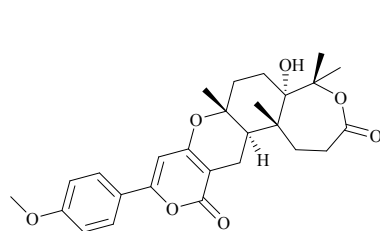


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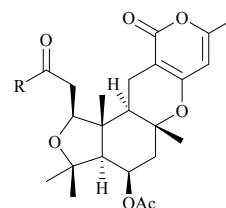


20 $R_1 = \text{OH}$, $R_2 = \text{OCH}_3$, $R_3 = \text{OCH}_3$

21 $R_1 = \text{OCH}_3$, $R_2 = \text{OH}$, $R_3 = \text{H}$

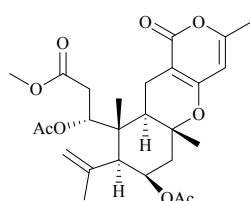


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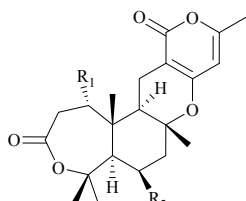


23 $R = \text{OCH}_3$

24 $R = \text{OH}$



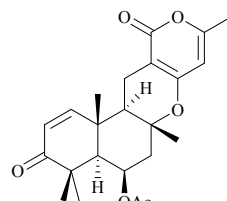
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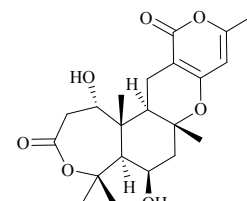
26 $R_1 = \text{OH}$, $R_2 = \text{H}$

27 $R_1 = \text{OH}$, $R_2 = \text{OAc}$

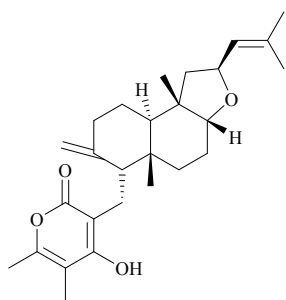
28 $R_1 = R_2 = \text{OAc}$



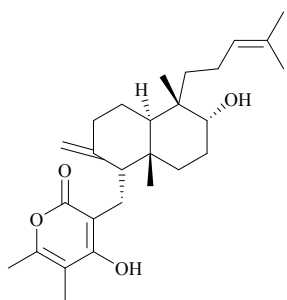
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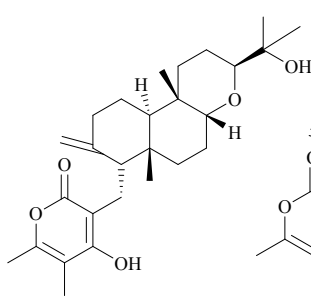
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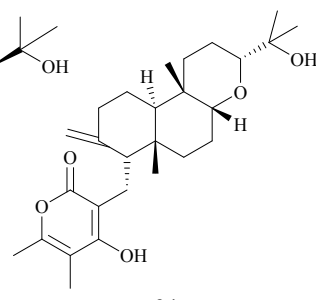
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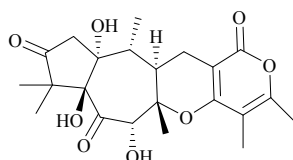
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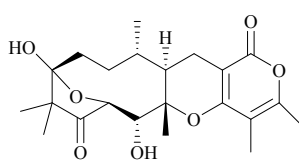
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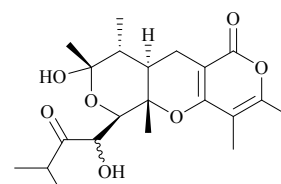
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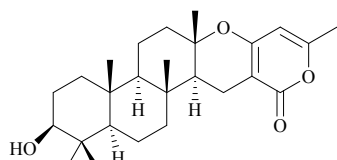
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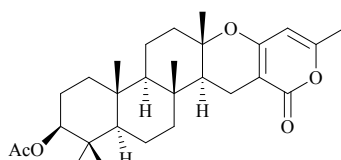
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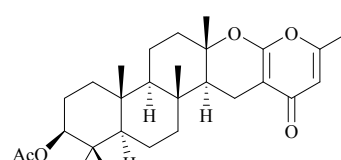
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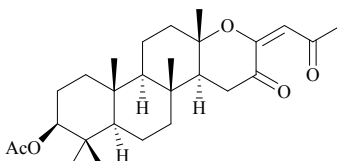
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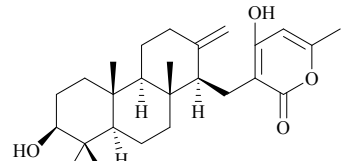
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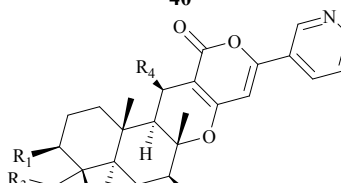
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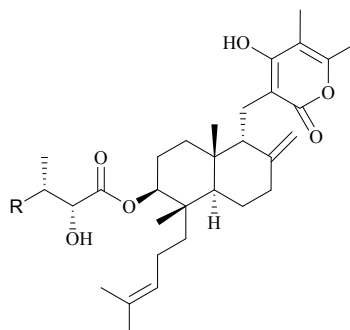
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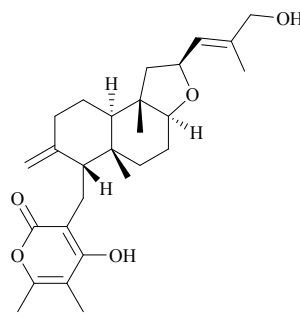
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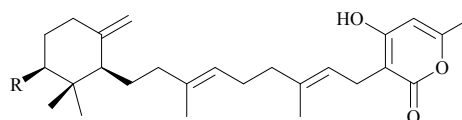
43 $R_1=OH$, $R_2=OAc$, $R_3=OH$, $R_4=H$
 44 $R_1=OH$, $R_2=OAc$, $R_3=OAc$, $R_4=OH$
 45 $R_1=OAc$, $R_2=H$, $R_3=OH$, $R_4=H$



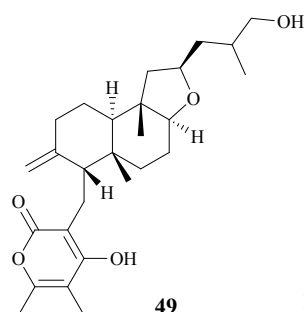
46 $R=CH_2CH_3$ 47 $R=CH_3$



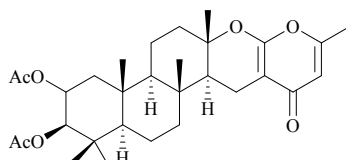
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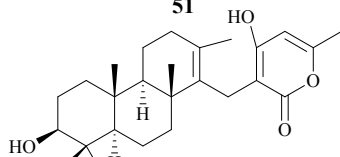
50 $R=OAc$ 53 $R=OH$



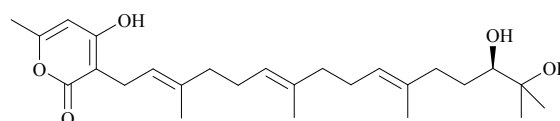
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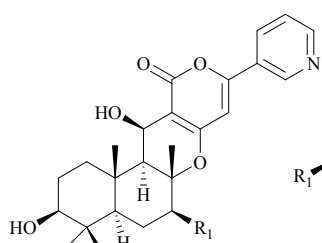
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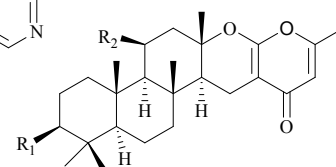
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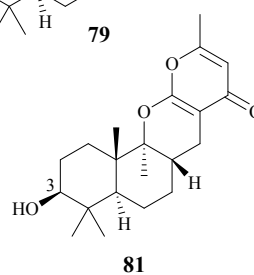
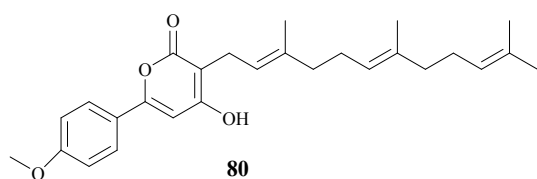
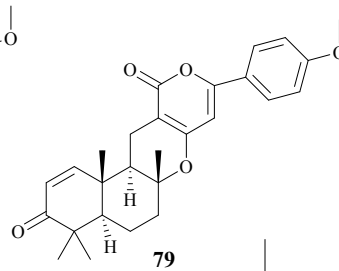
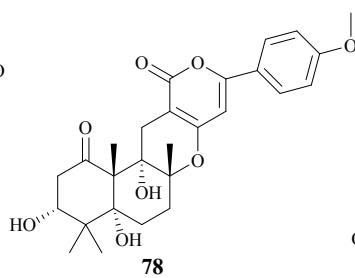
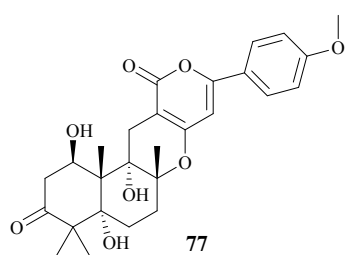
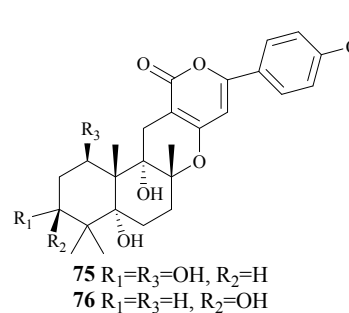
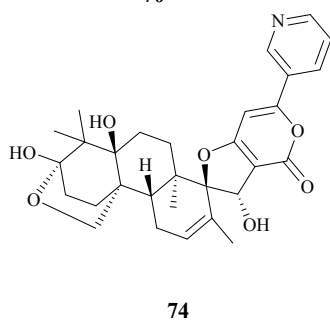
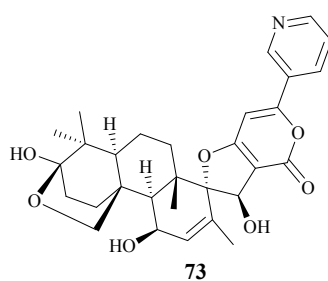
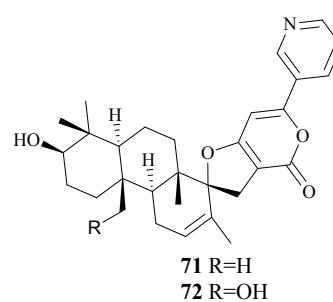
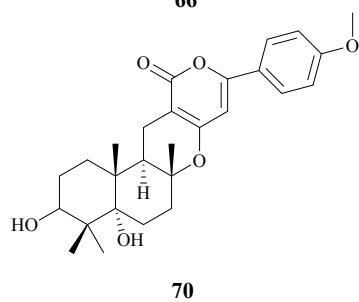
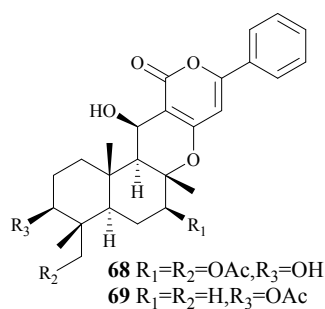
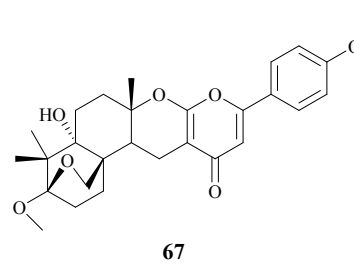
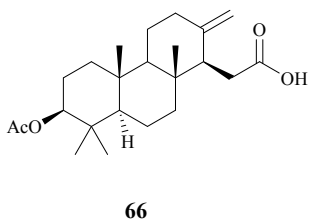
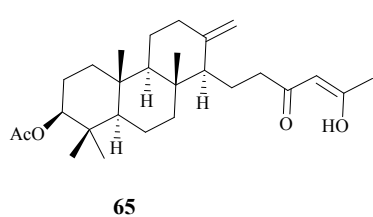
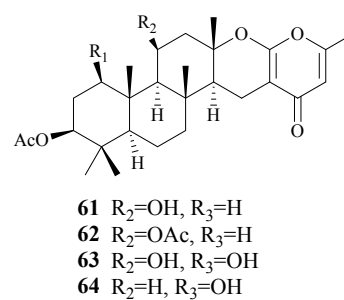
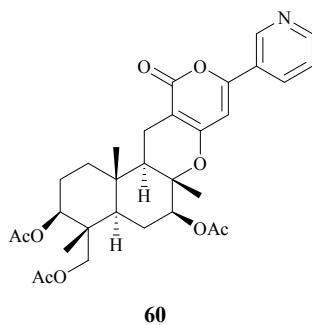
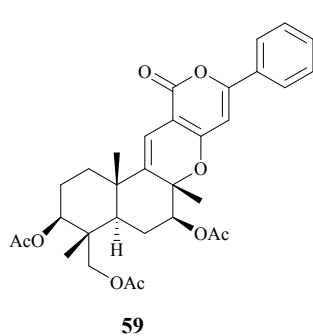
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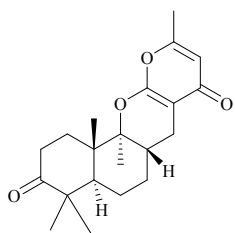


55 $R_1=OH$
 56 $R_1=OAc$

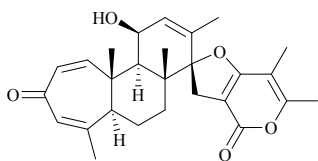


57 $R_1=O$, $R_2=H$
 58 $R_1=OH$, $R_2=OH$

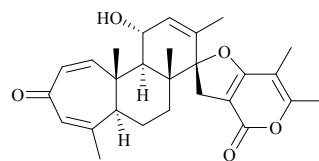




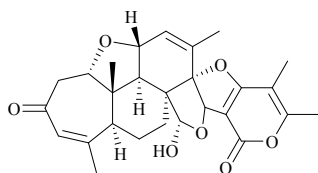
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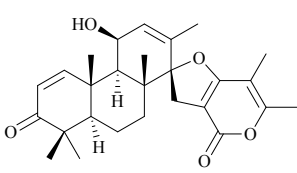
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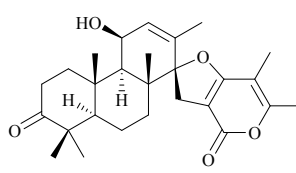
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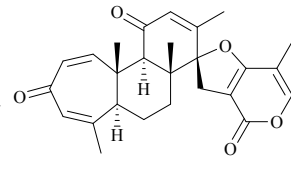
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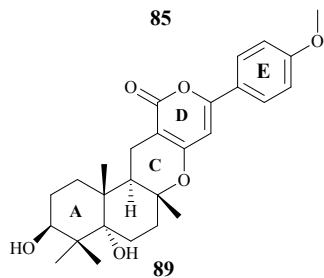
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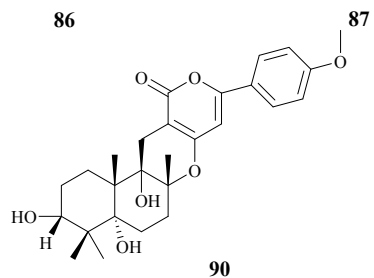
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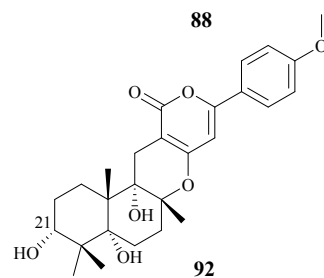
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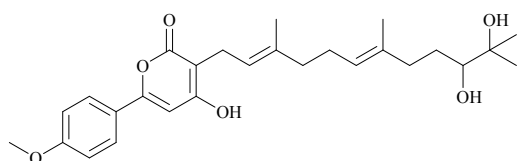
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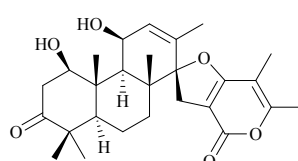
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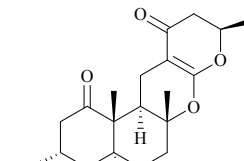
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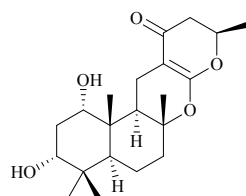
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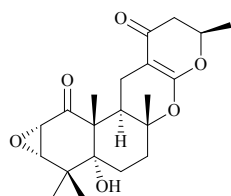
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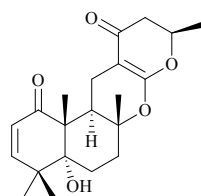
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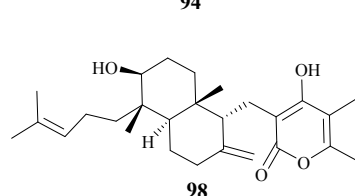
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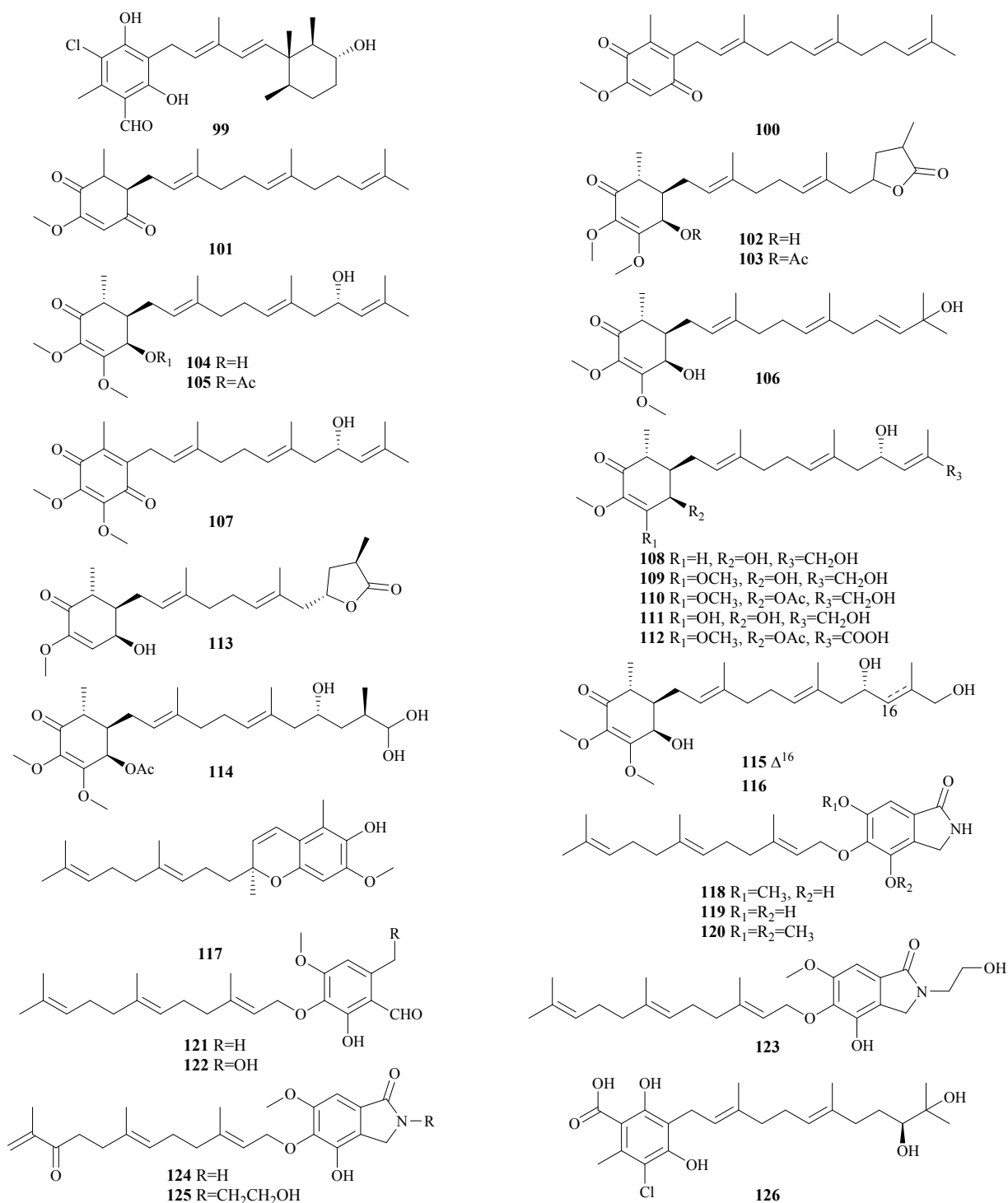
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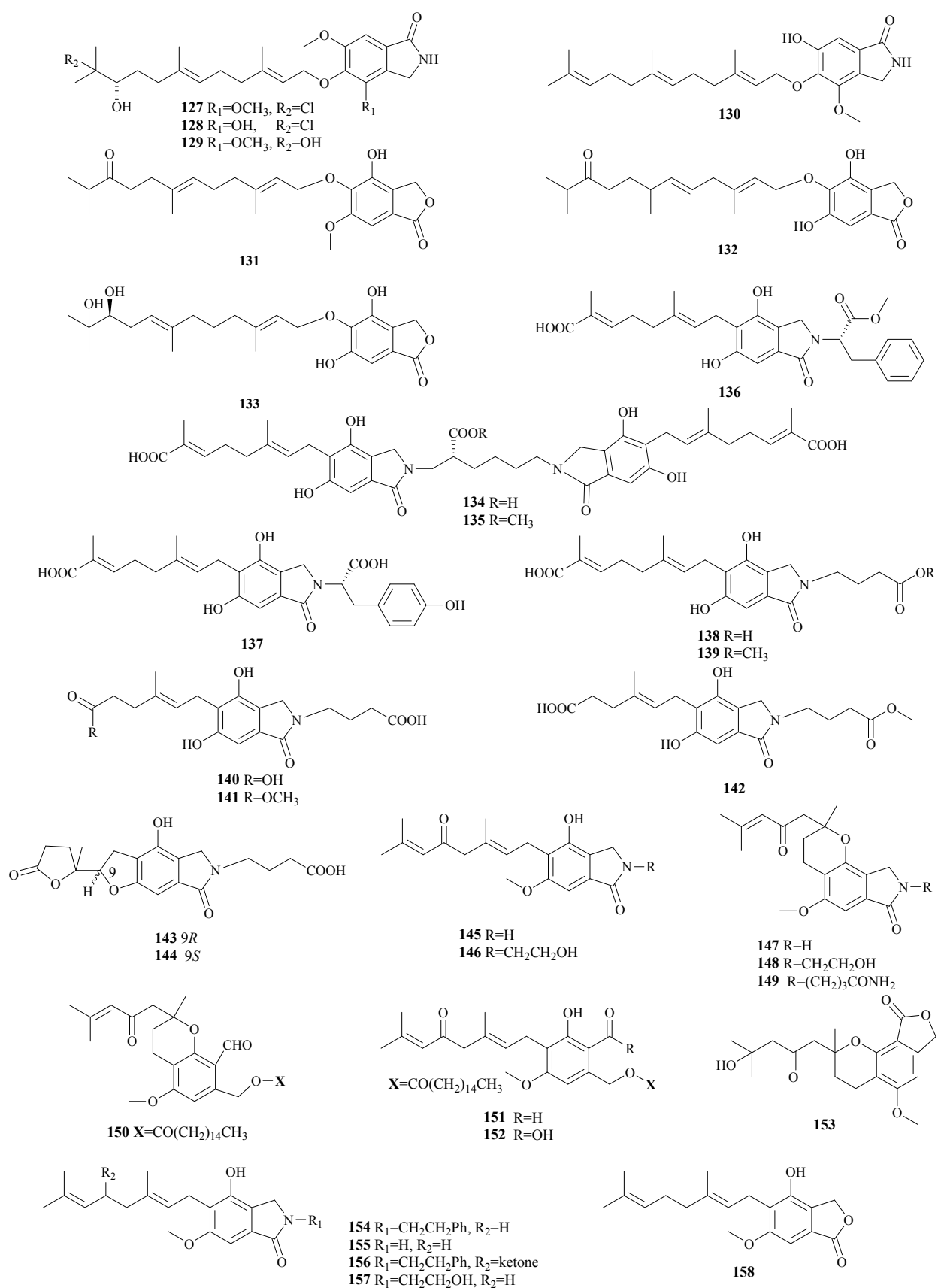


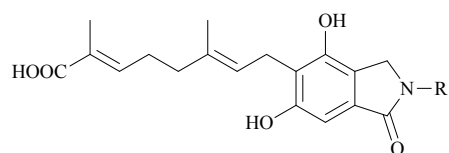
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2 Tetraketide-terpenoids (99-652)

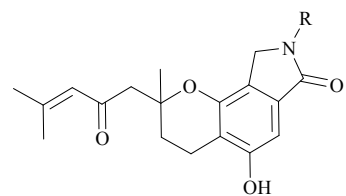
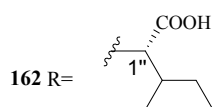
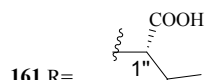
2.1 Derived from Orsellinic acid/orcinaldehyde (99-323)



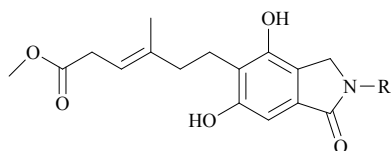




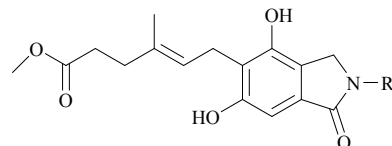
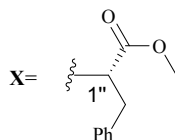
159 R=H
160 R=CH₂CH₂OH
171 R=CH₂CH₂Ph



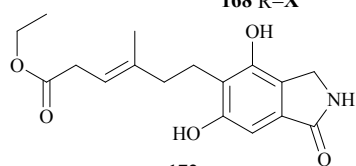
169 R=CH₃
170 R=(CH₂)₃COOH



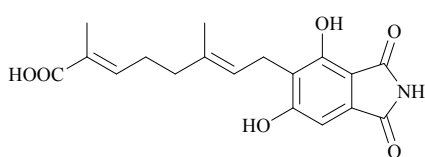
164 R=H
166 R=CH₂CH₂OH
168 R=X



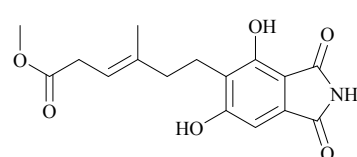
163 R=H
165 R=CH₂CH₂OH
167 R=X



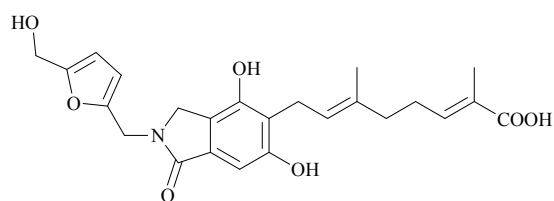
172



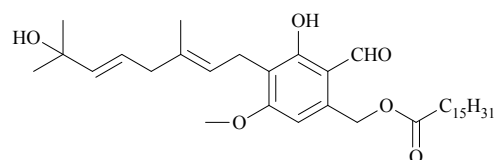
173



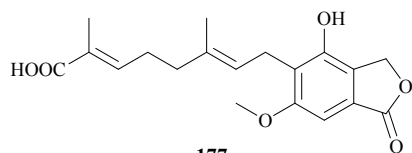
174



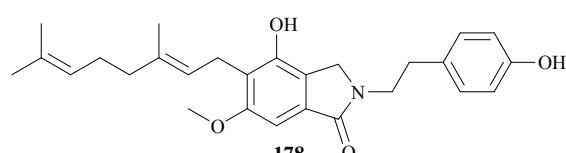
175



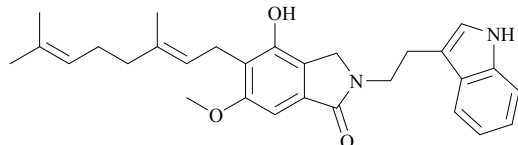
176



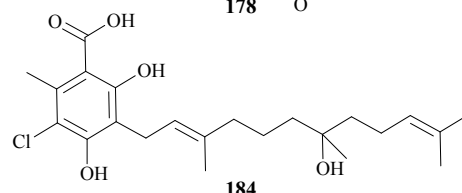
177



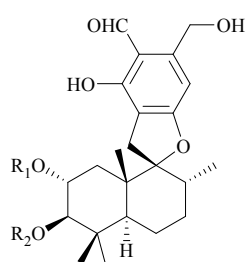
178



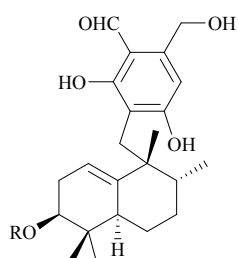
179



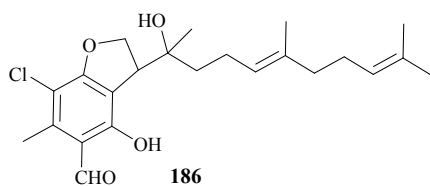
184



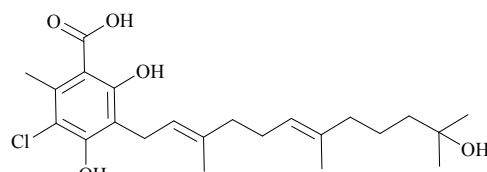
180 R₁=Ac, R₂=H
181 R₁=H, R₂=Ac



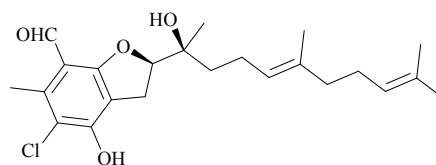
182 R=Ac
183 R=H



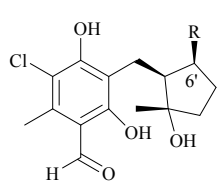
186



185



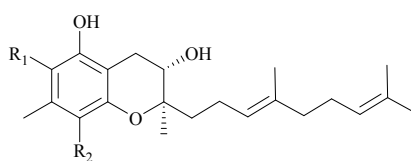
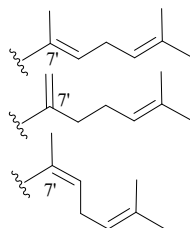
187



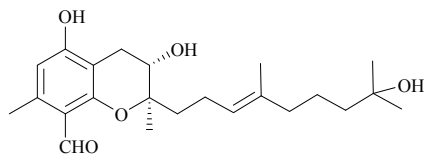
188 R=

189 R=

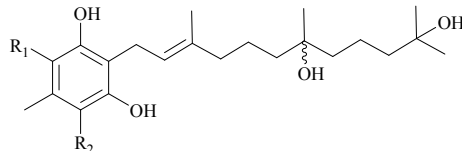
190 R=



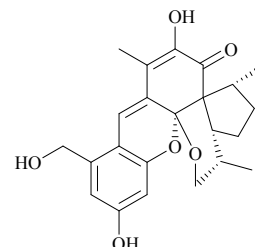
191 $R_1=CHO$, $R_2=Cl$
192 $R_1=CHO$, $R_2=H$
193 $R_1=Cl$, $R_2=CHO$



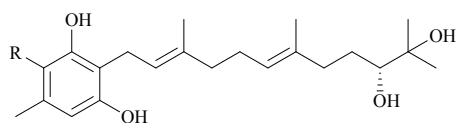
194



195/196 $R_1=Cl$, $R_2=COOH$
197 $R_1=H$, $R_2=COOH$
198 $R_1=Cl$, $R_2=H$

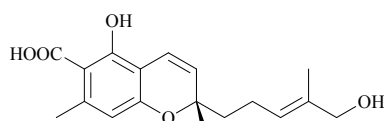


199

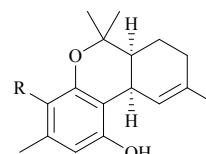


200 R=COOH

203 R=H

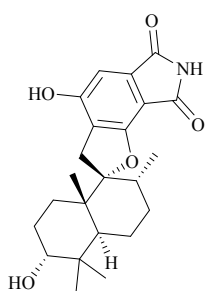


202

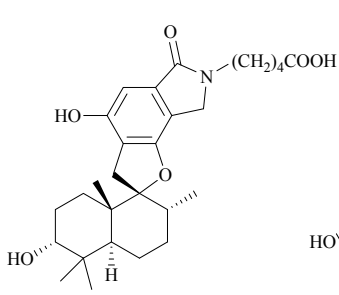


201 R=COOH

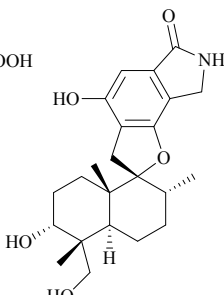
204 R=H



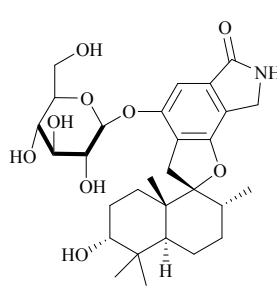
205



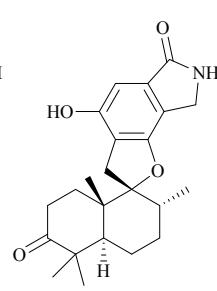
206



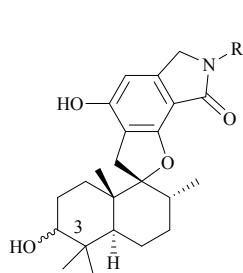
207



208



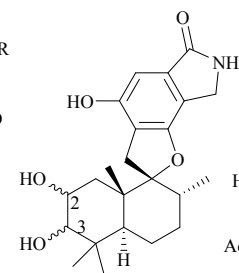
209



210 3R, R=H

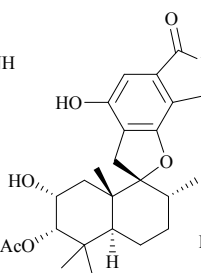
211 3S, R=H

212 3R, R=(CH₂)₂OH

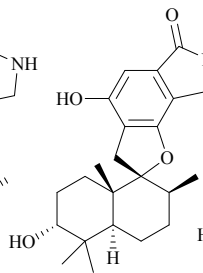


213 2R, 3R

214 2S, 3S

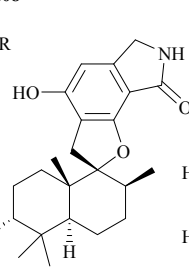


215

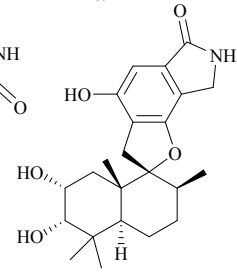


216 R=H

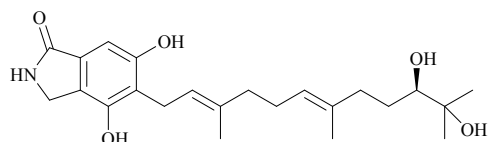
217 R=(CH₂)₂OH



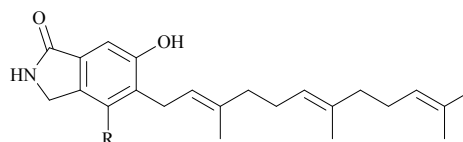
218



219

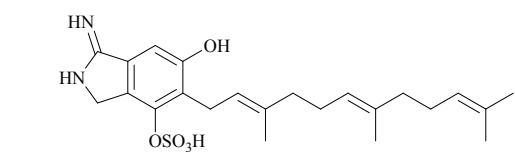


220

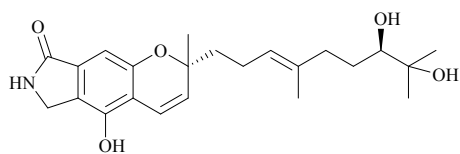


221 R=OH

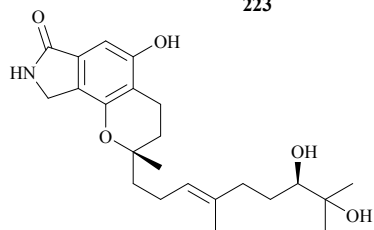
222 R=OSO₃H



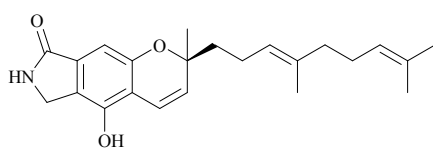
223



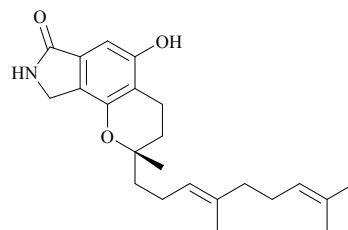
224



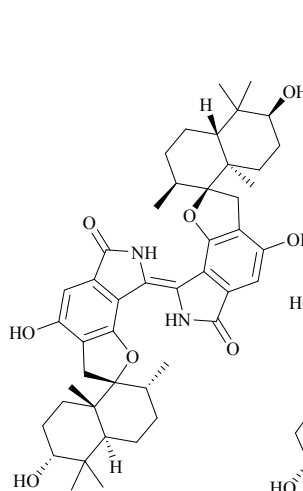
225



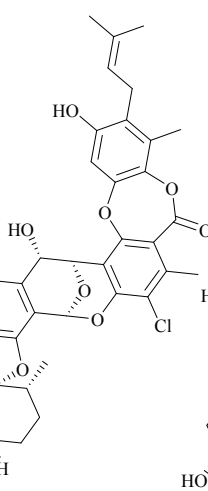
226



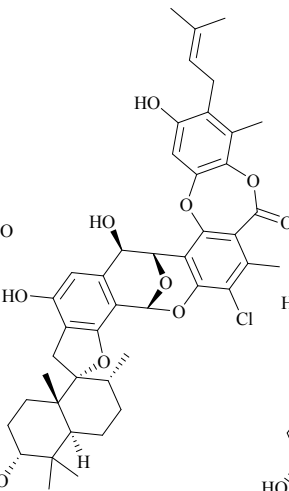
227



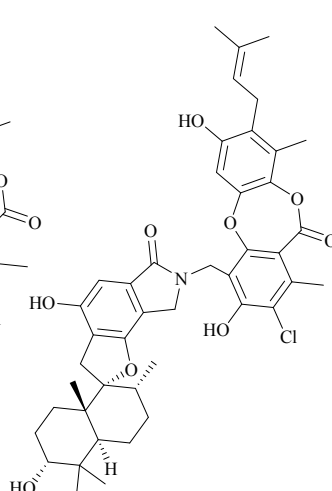
228



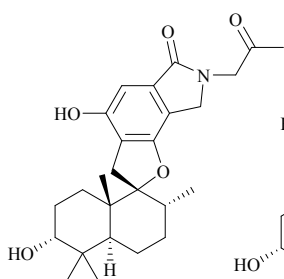
229



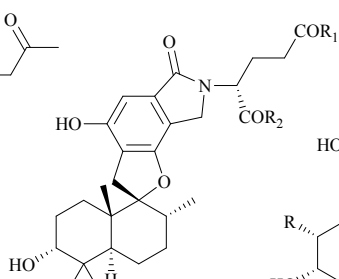
230



231

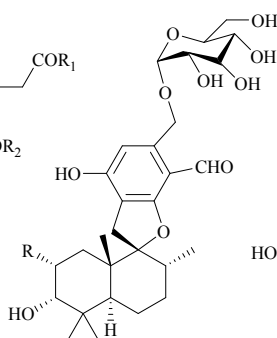


232



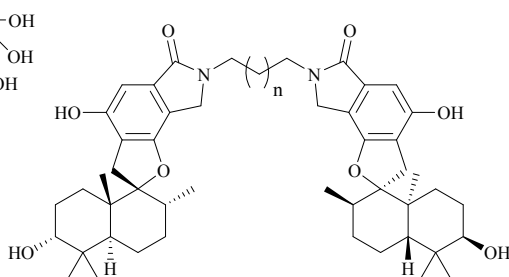
233 $R_1 = \text{OCH}_3$, $R_2 = \text{OH}$

234 $R_1 = \text{OH}$, $R_2 = \text{OCH}_3$



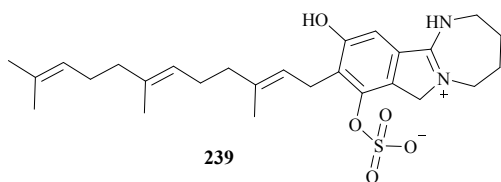
235 $R = \text{H}$

236 $R = \text{OH}$

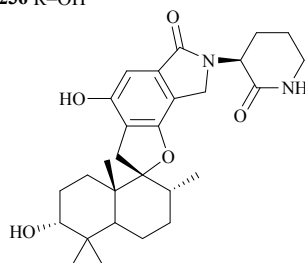


237 $n=2$

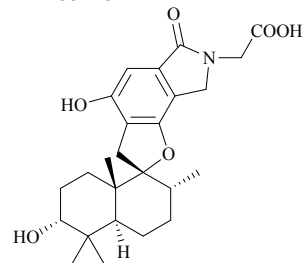
238 $n=3$



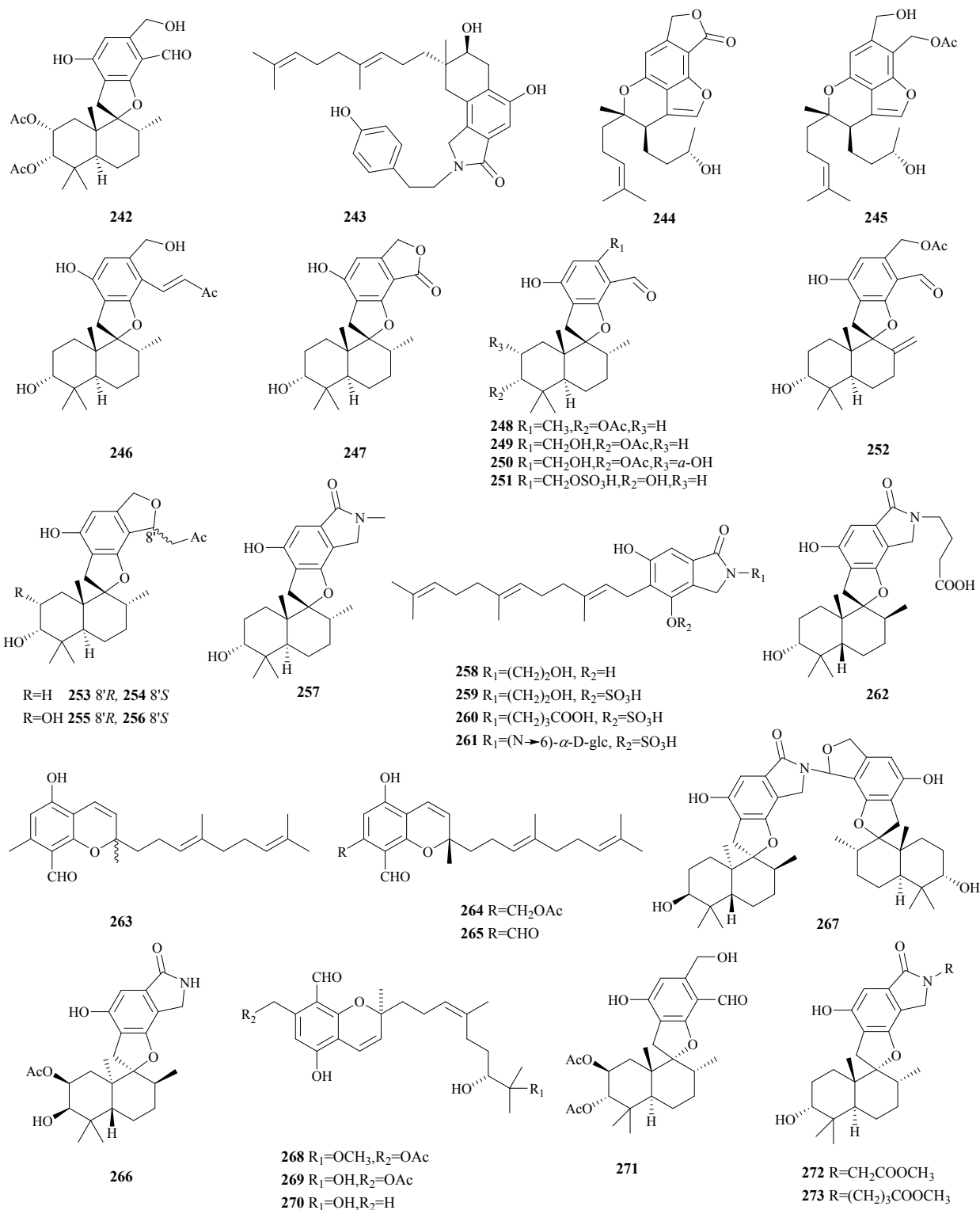
239

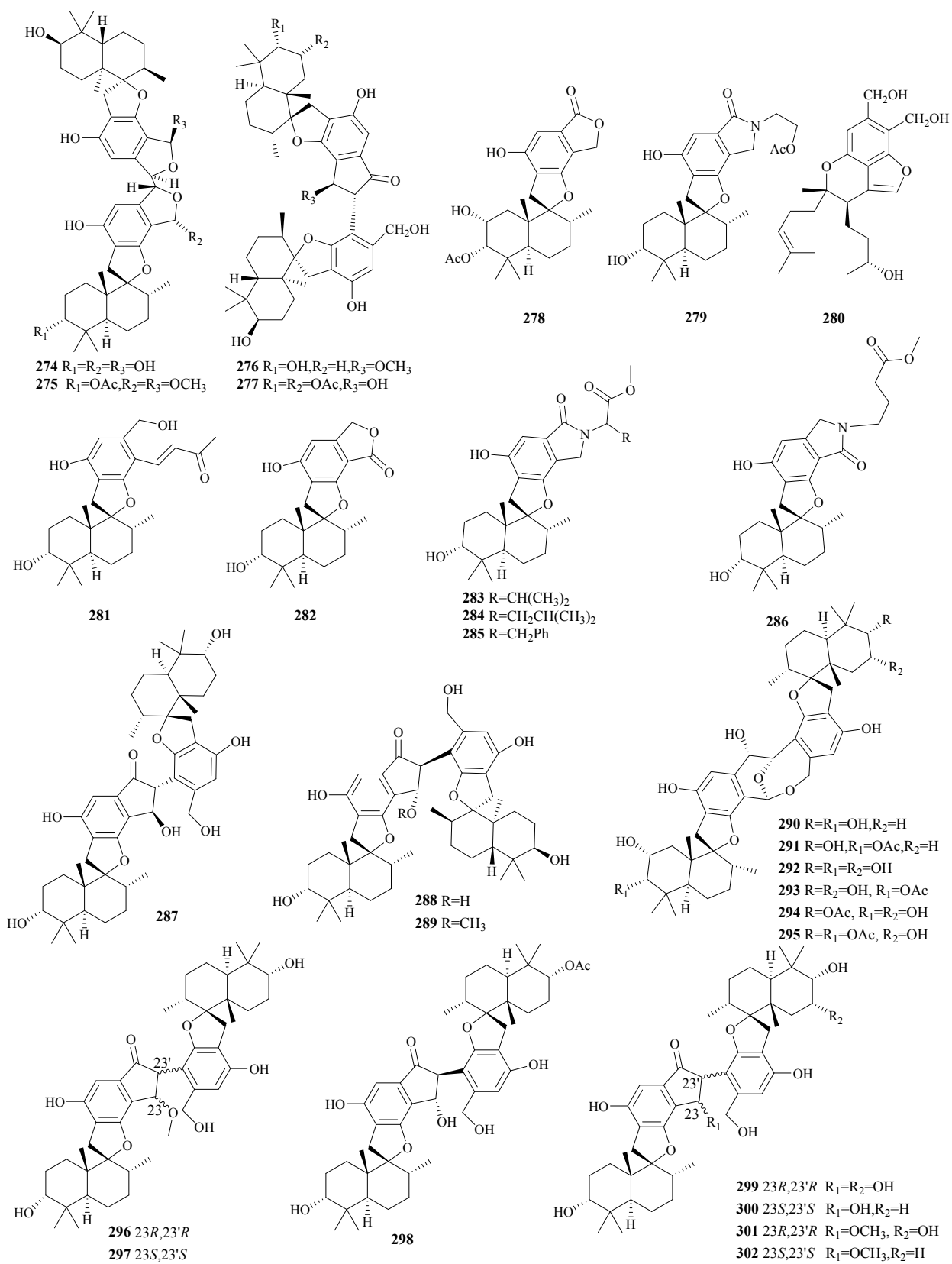


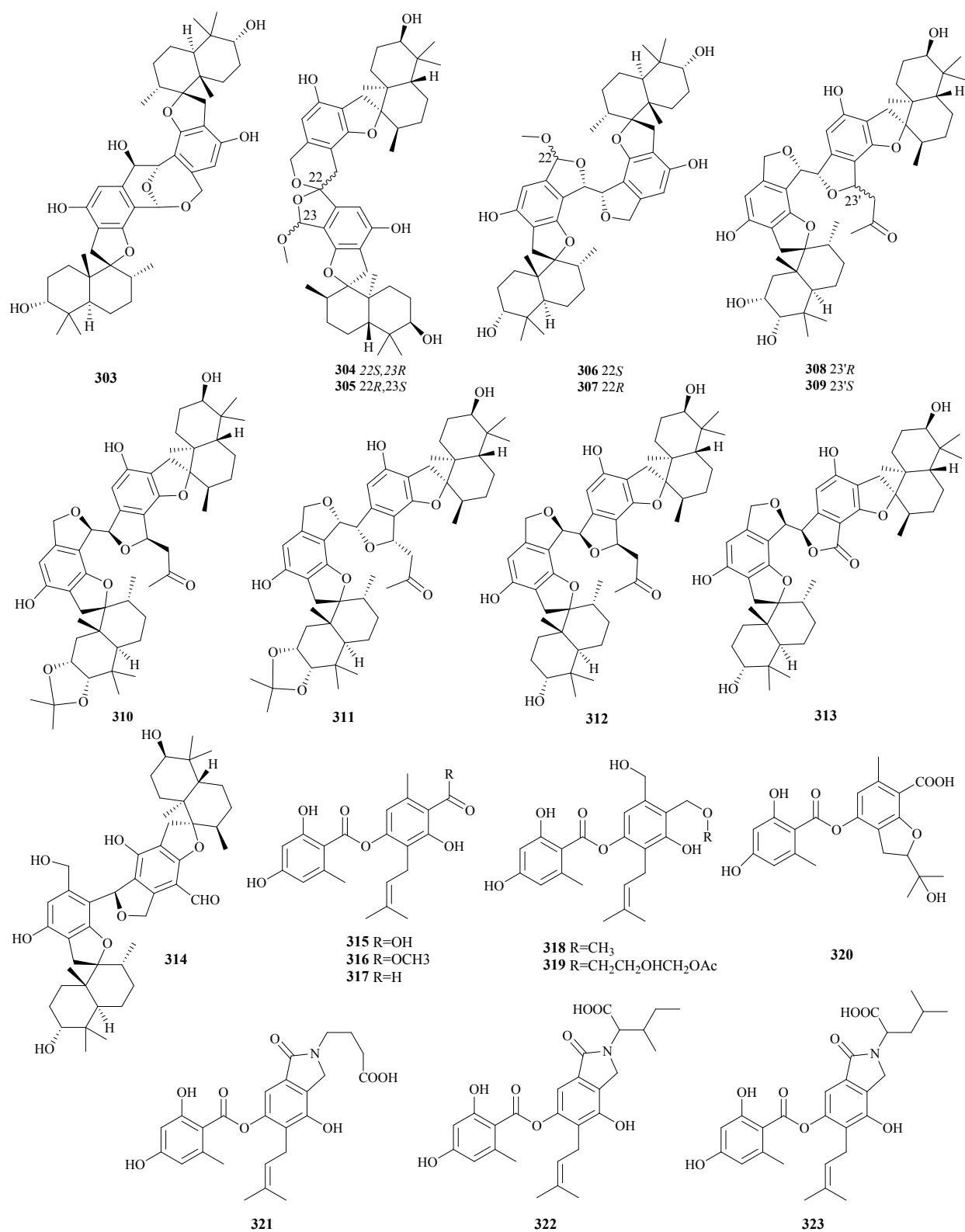
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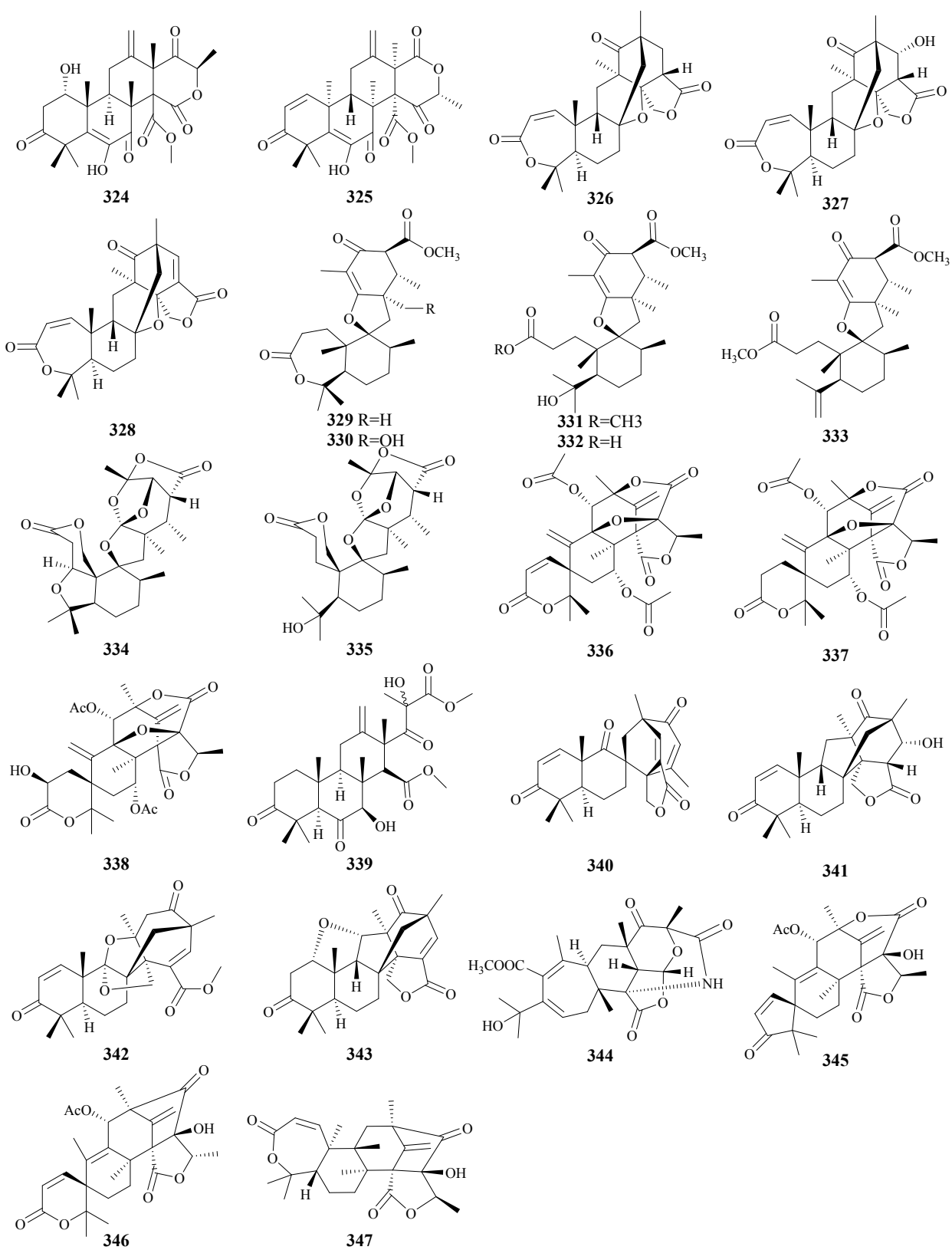
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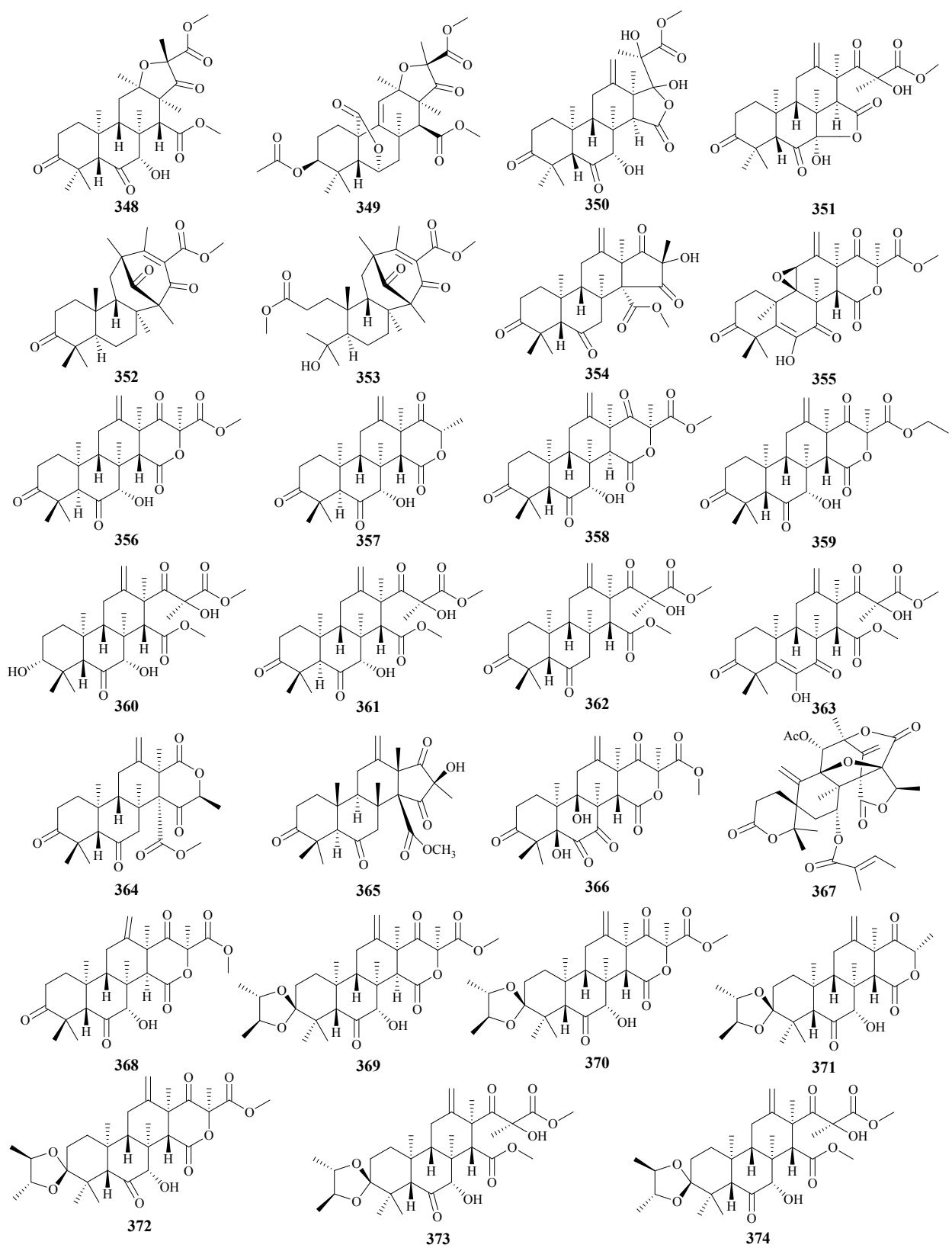


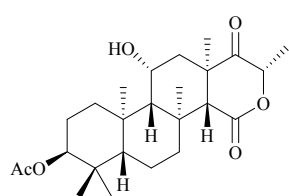




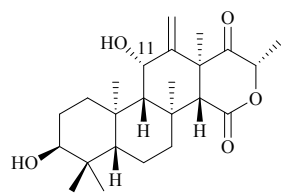
2.2 Derived from 3,5-Dimethylorsellinic acid (DMOA) (324-492)



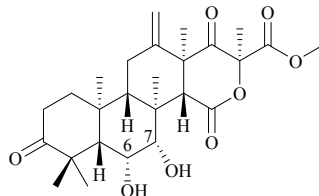




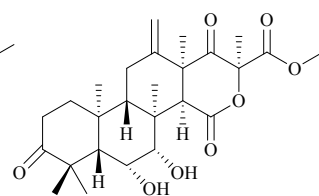
375



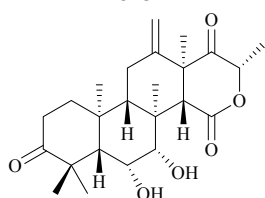
376



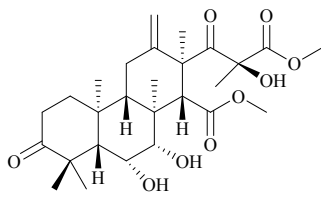
377



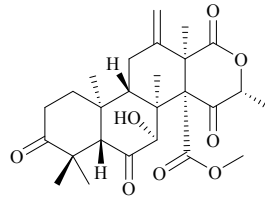
378



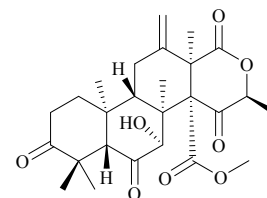
379



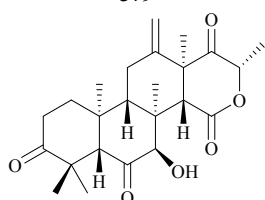
380



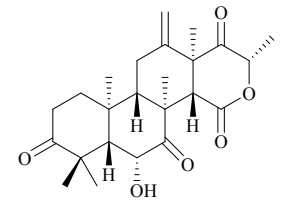
381



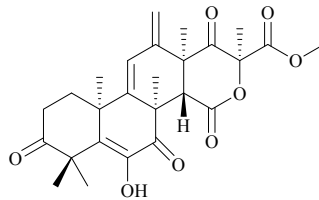
382



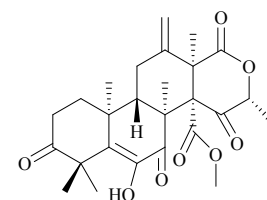
383



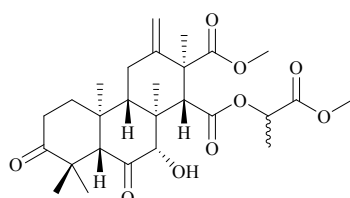
384



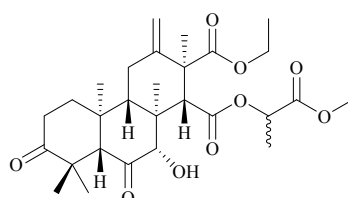
385



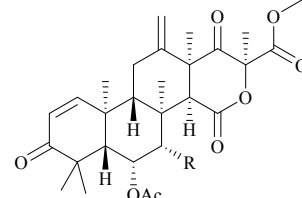
388



386

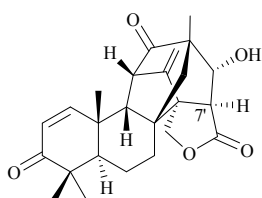


387

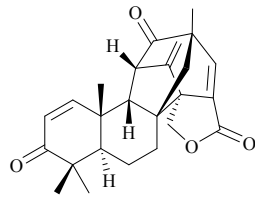


389 R=OH

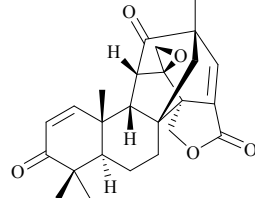
390 R=H



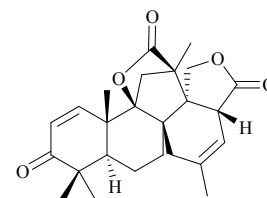
391



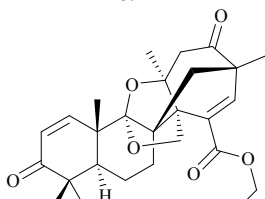
392



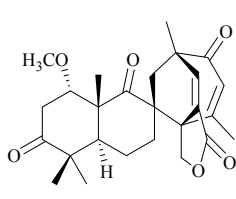
393



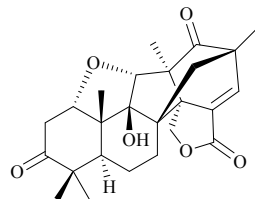
394



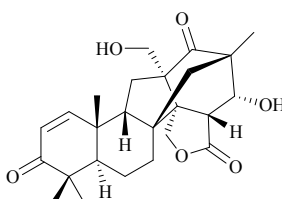
395



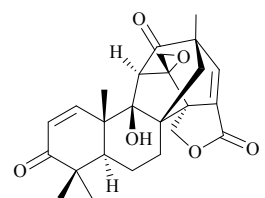
396



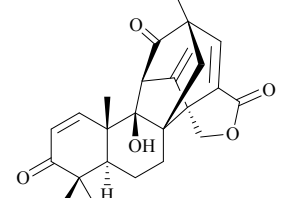
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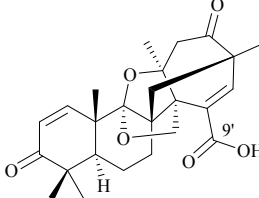
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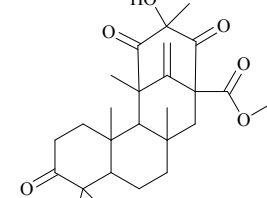
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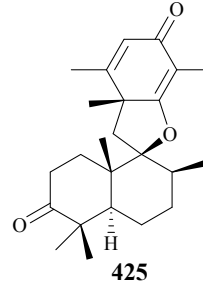
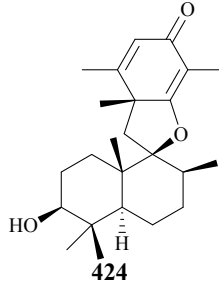
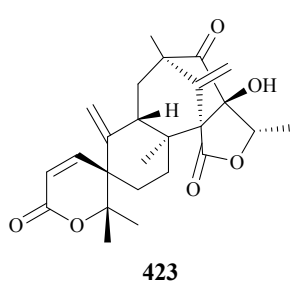
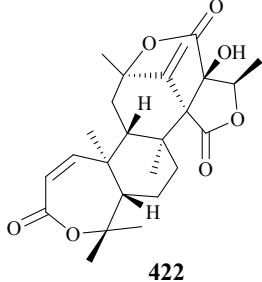
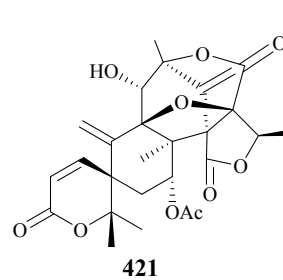
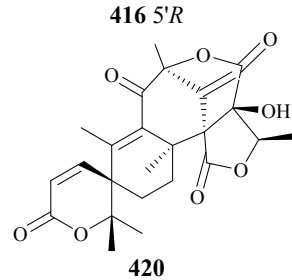
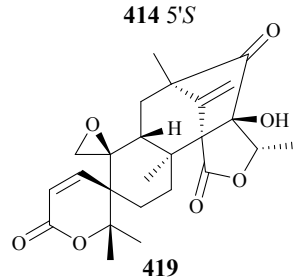
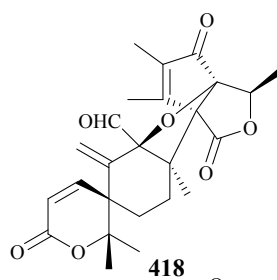
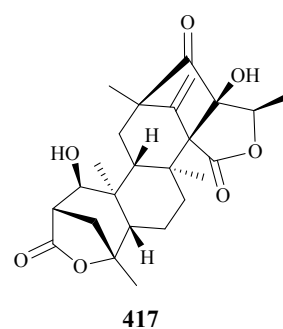
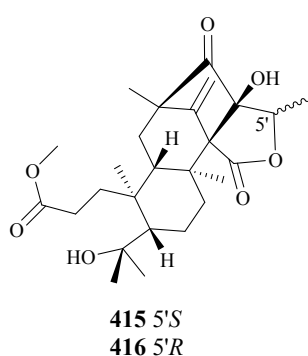
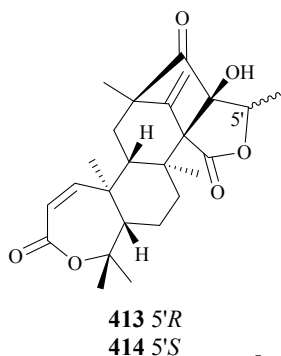
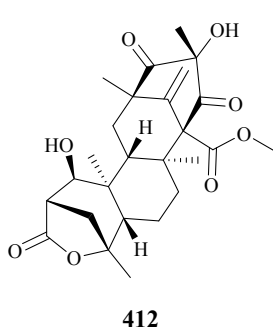
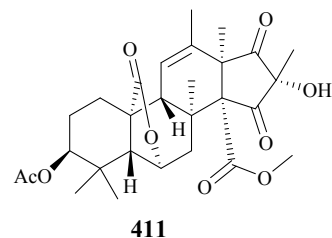
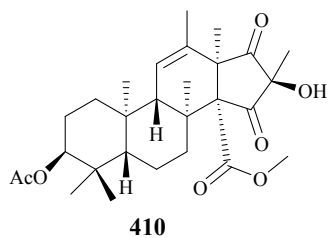
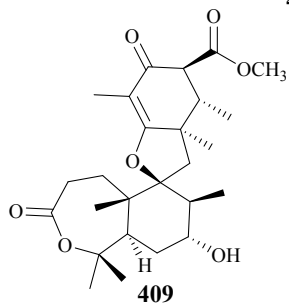
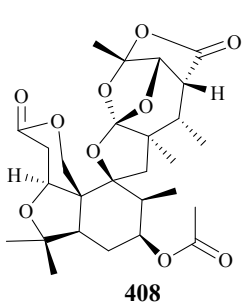
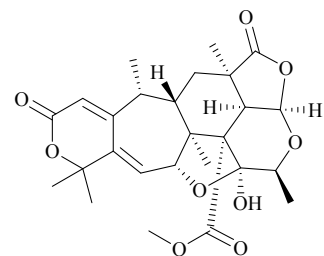
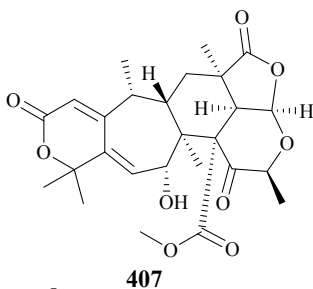
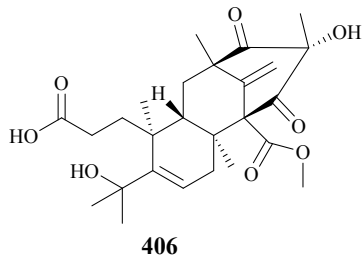
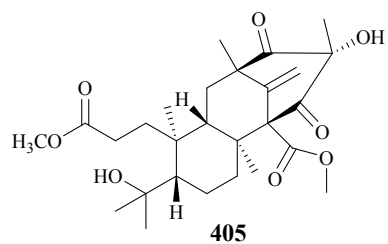
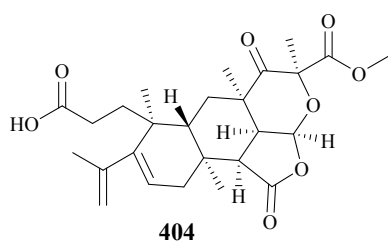
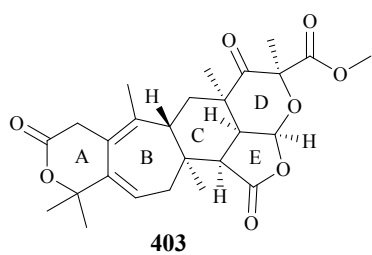
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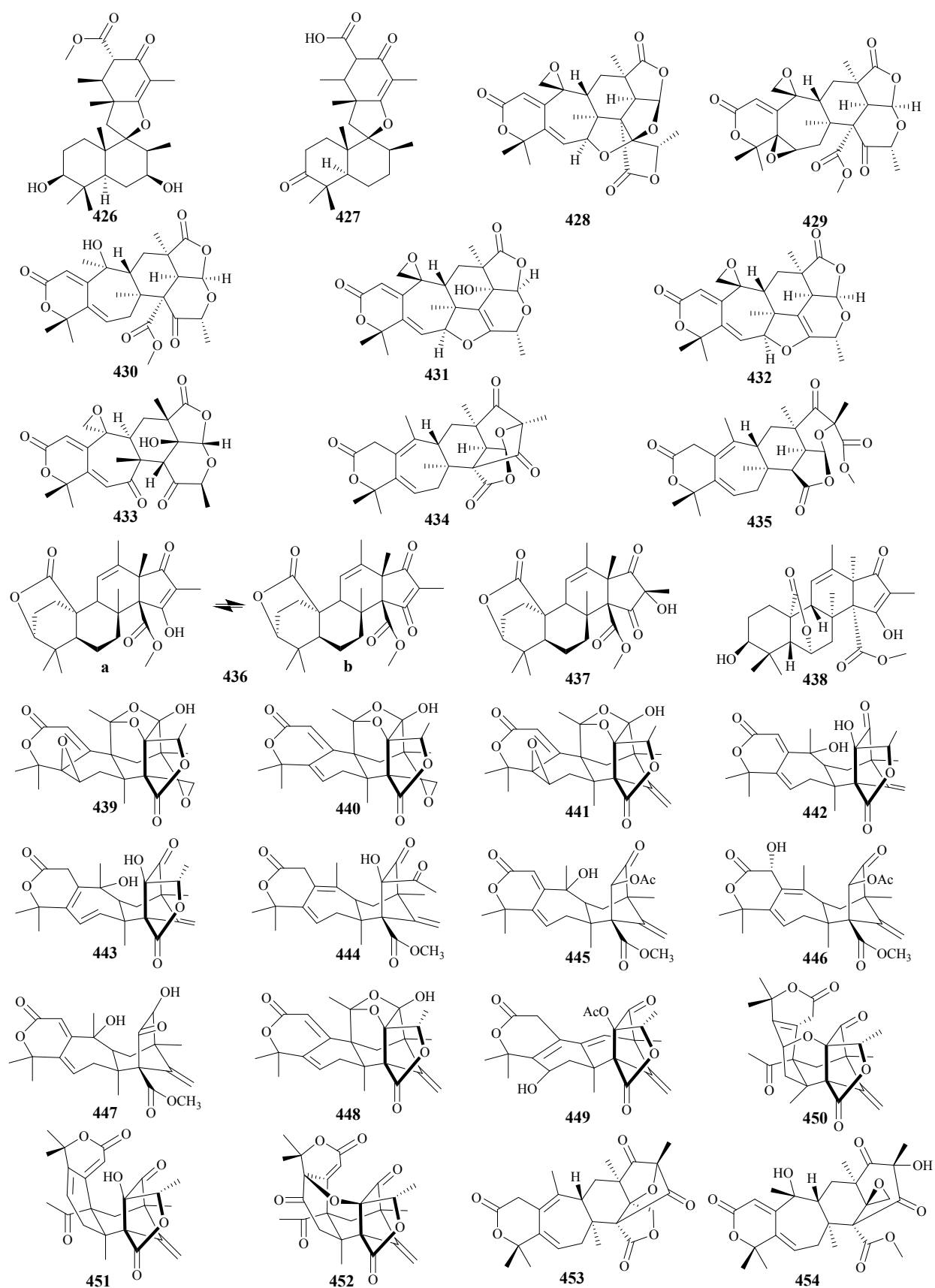


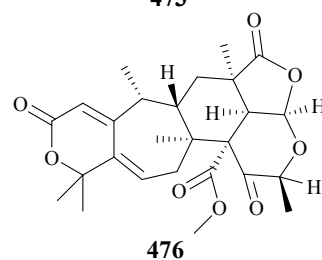
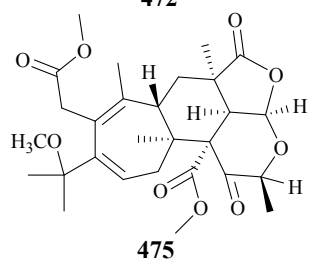
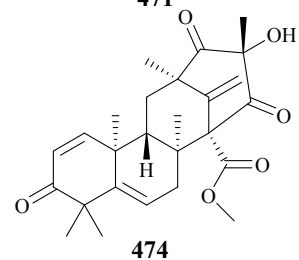
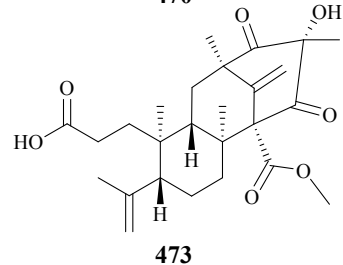
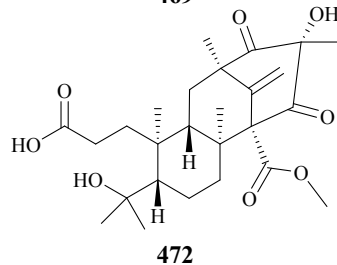
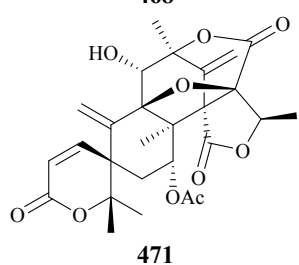
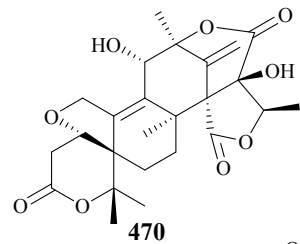
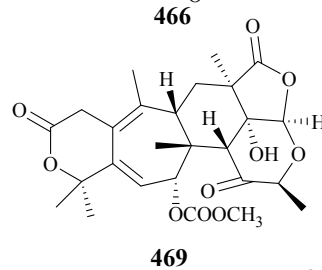
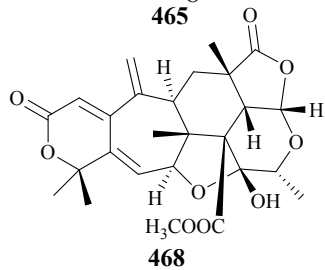
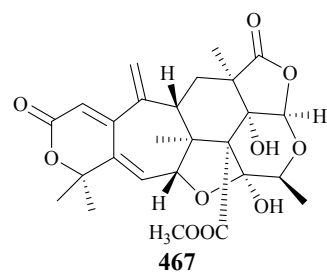
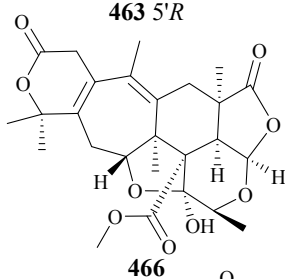
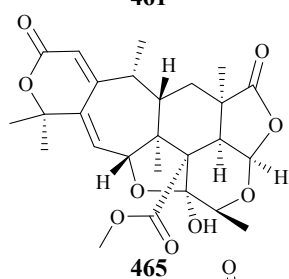
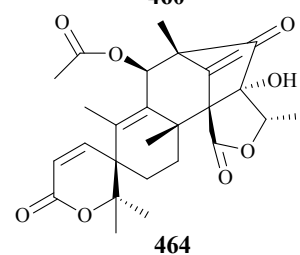
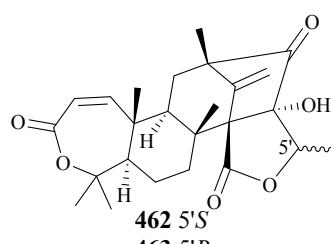
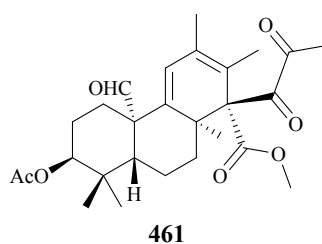
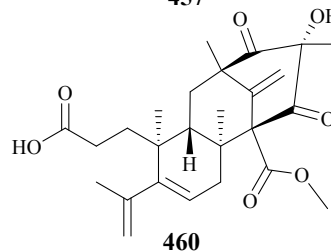
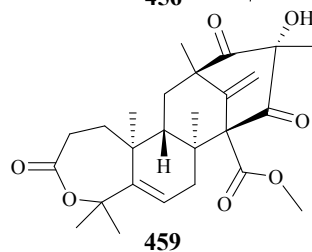
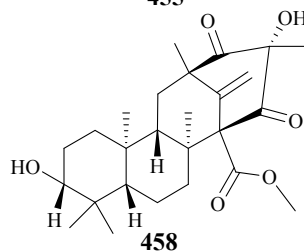
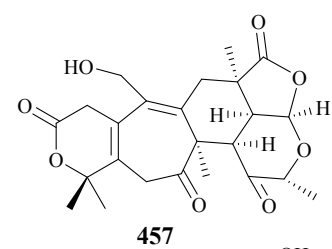
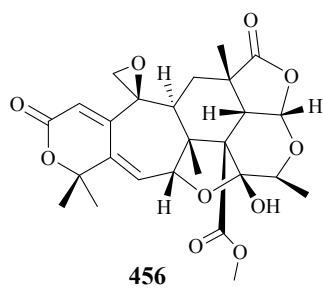
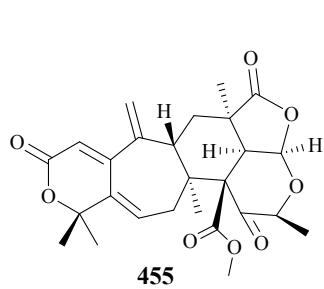
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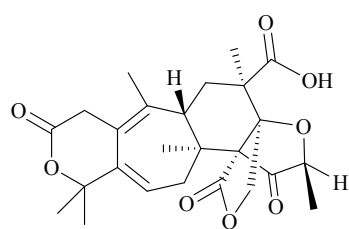


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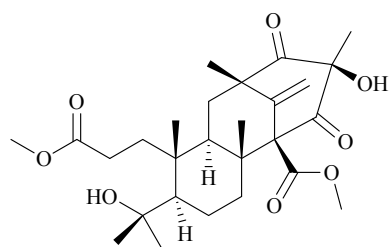




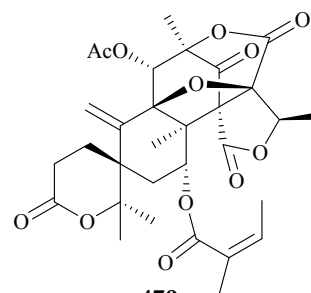




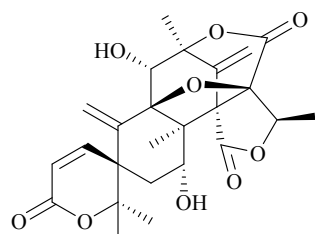
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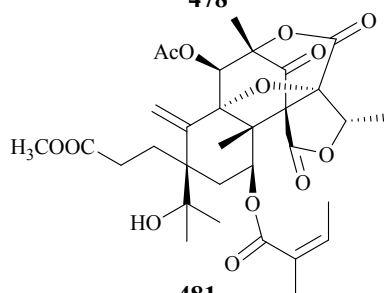
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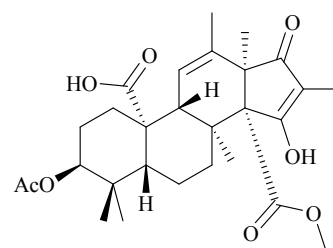
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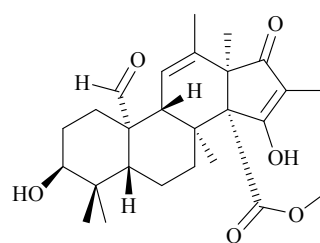
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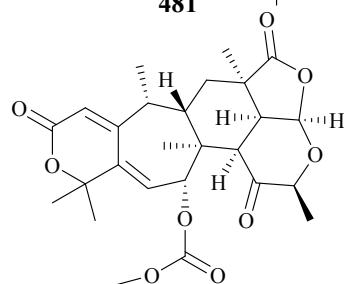
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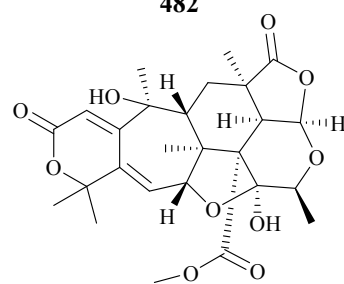
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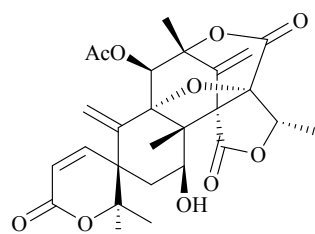
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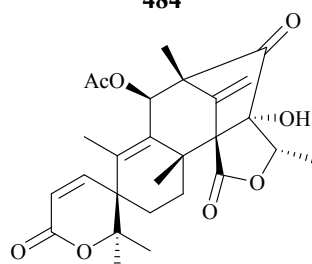
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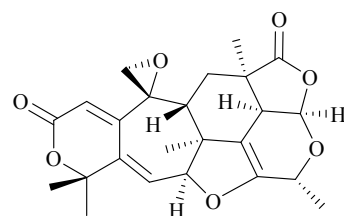
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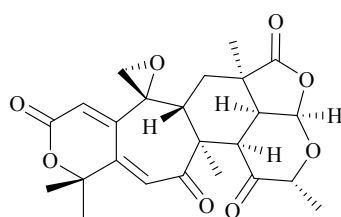
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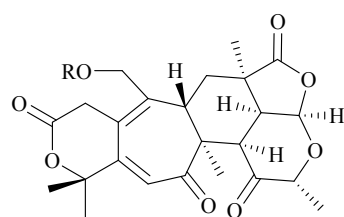
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488

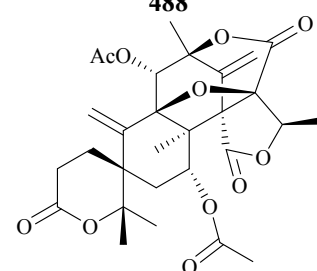


489



490 R=Ac

491 R=COCH₂CH(CH₃)₂



492

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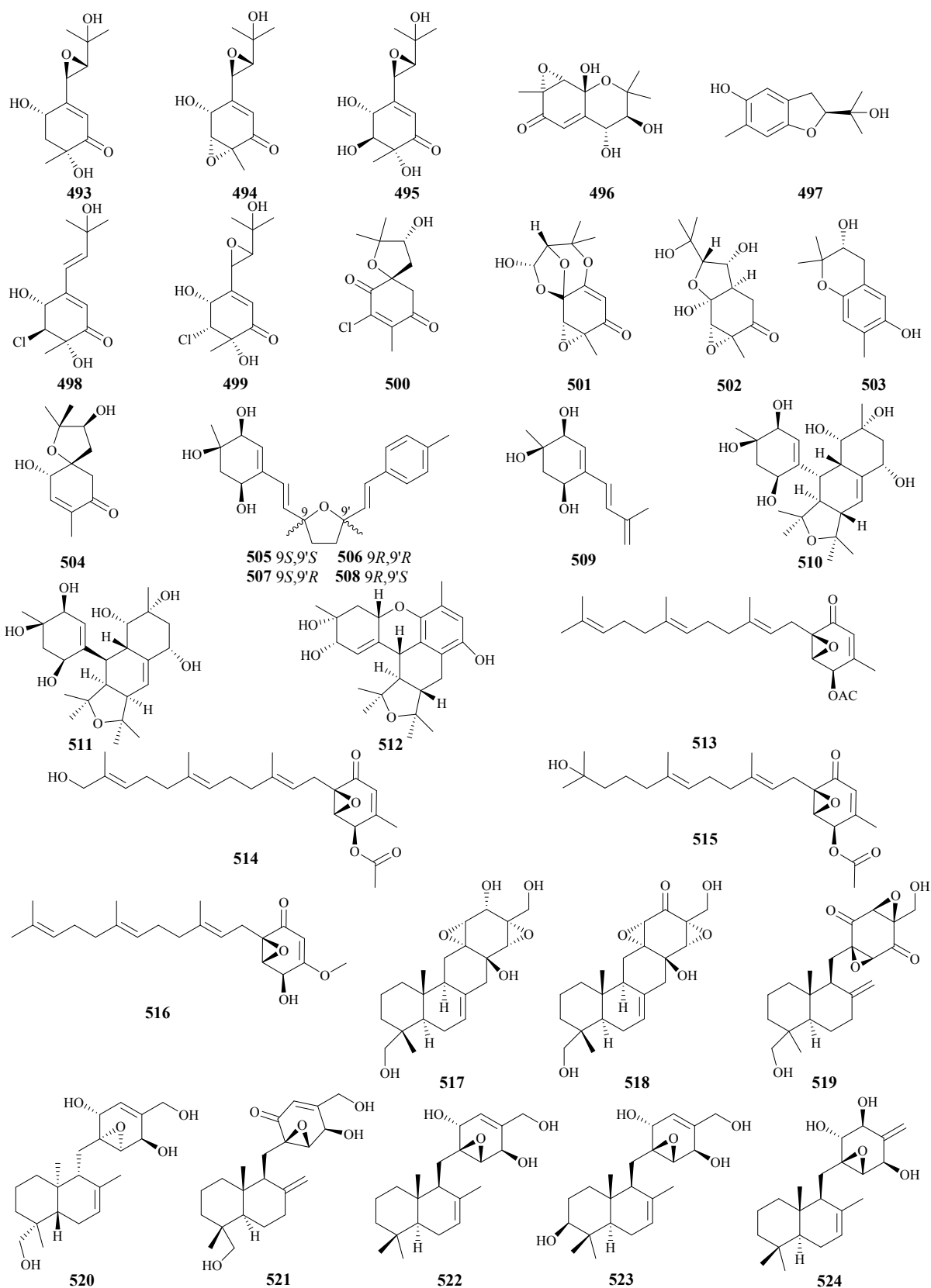
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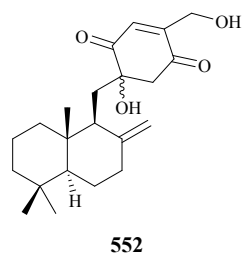
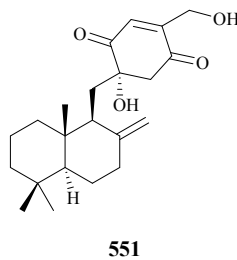
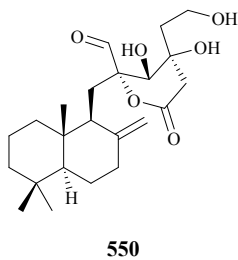
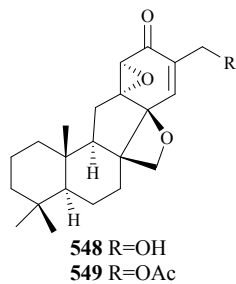
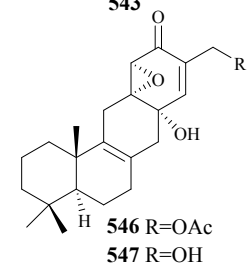
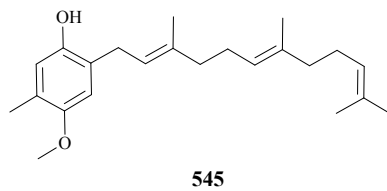
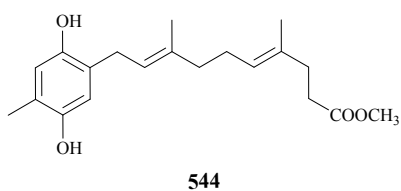
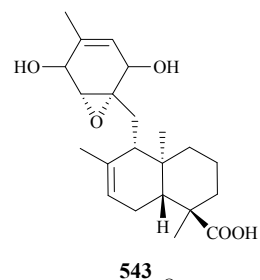
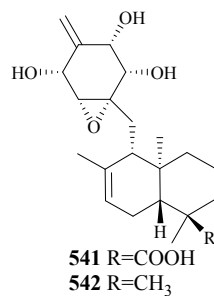
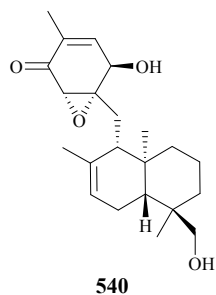
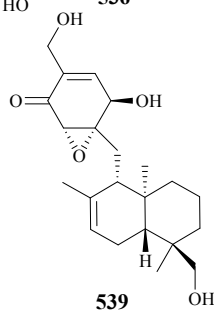
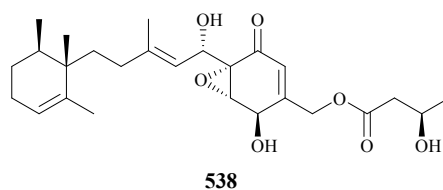
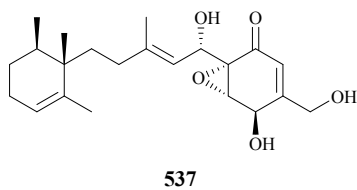
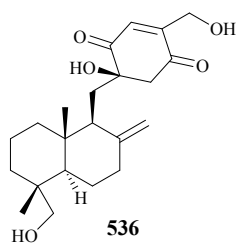
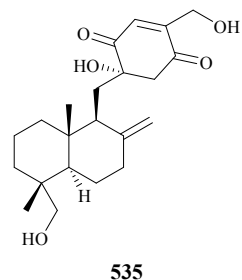
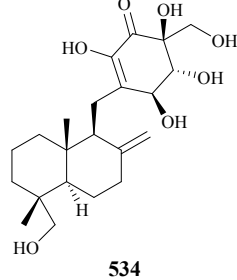
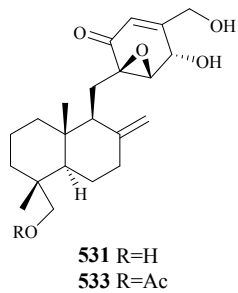
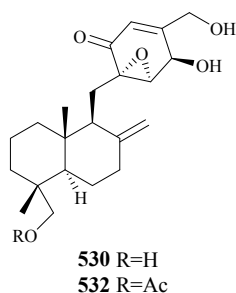
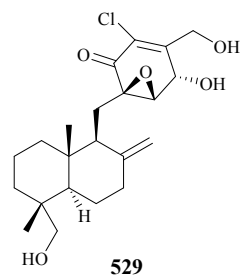
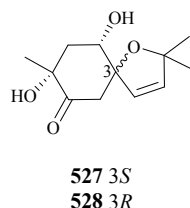
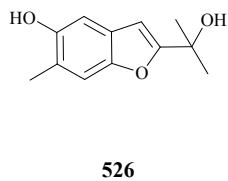
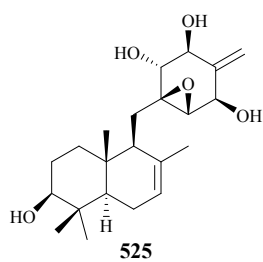
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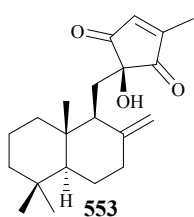
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(493-

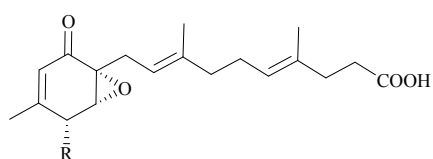
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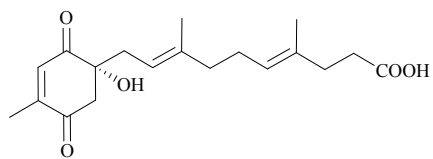




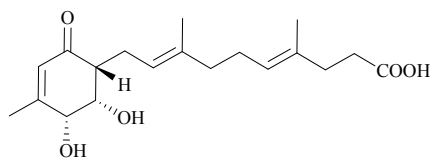
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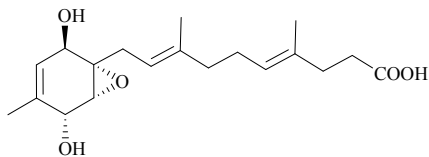
554 R=OH
558 R=OAc



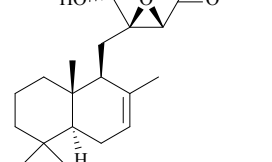
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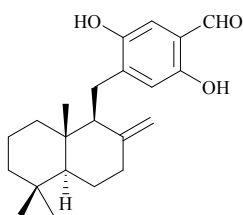
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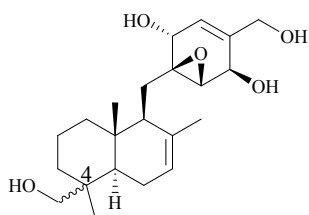
557



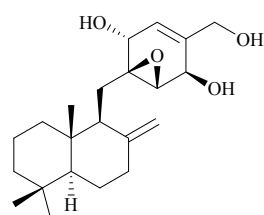
559



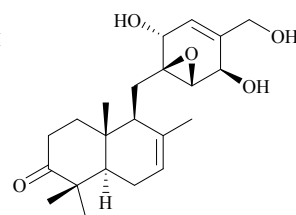
560



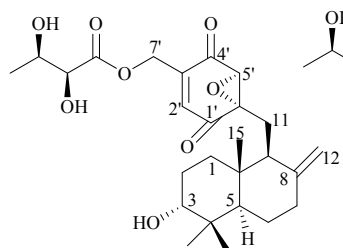
561 4S
562 4R



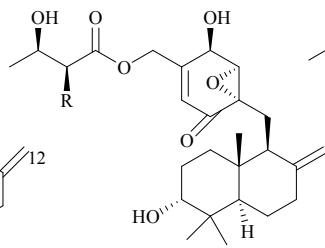
563



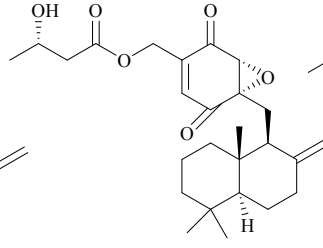
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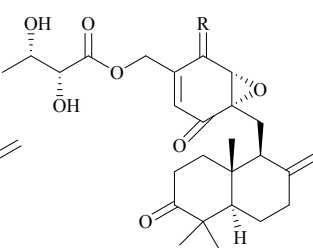
565



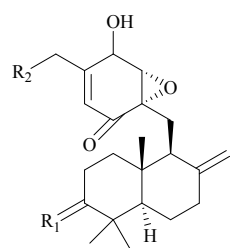
566 R=OH
567 R=H



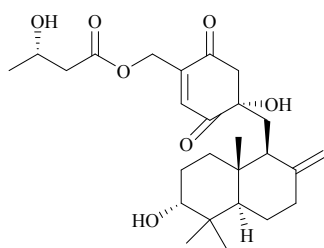
568



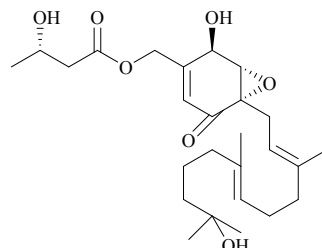
569 R=O
570 R=α-H/β-OH



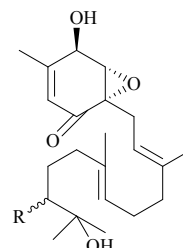
571 R₁=α-OH/β-H, R₂=H
572 R₁=O, R₂=OH



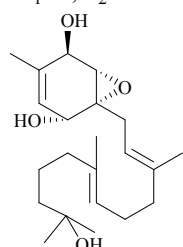
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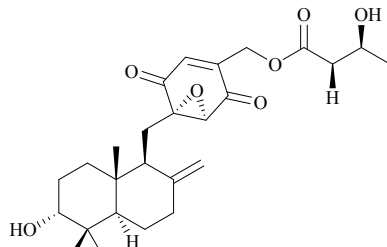
574



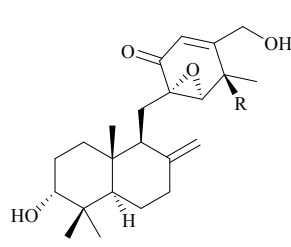
575 R=H
576 R=OH



577

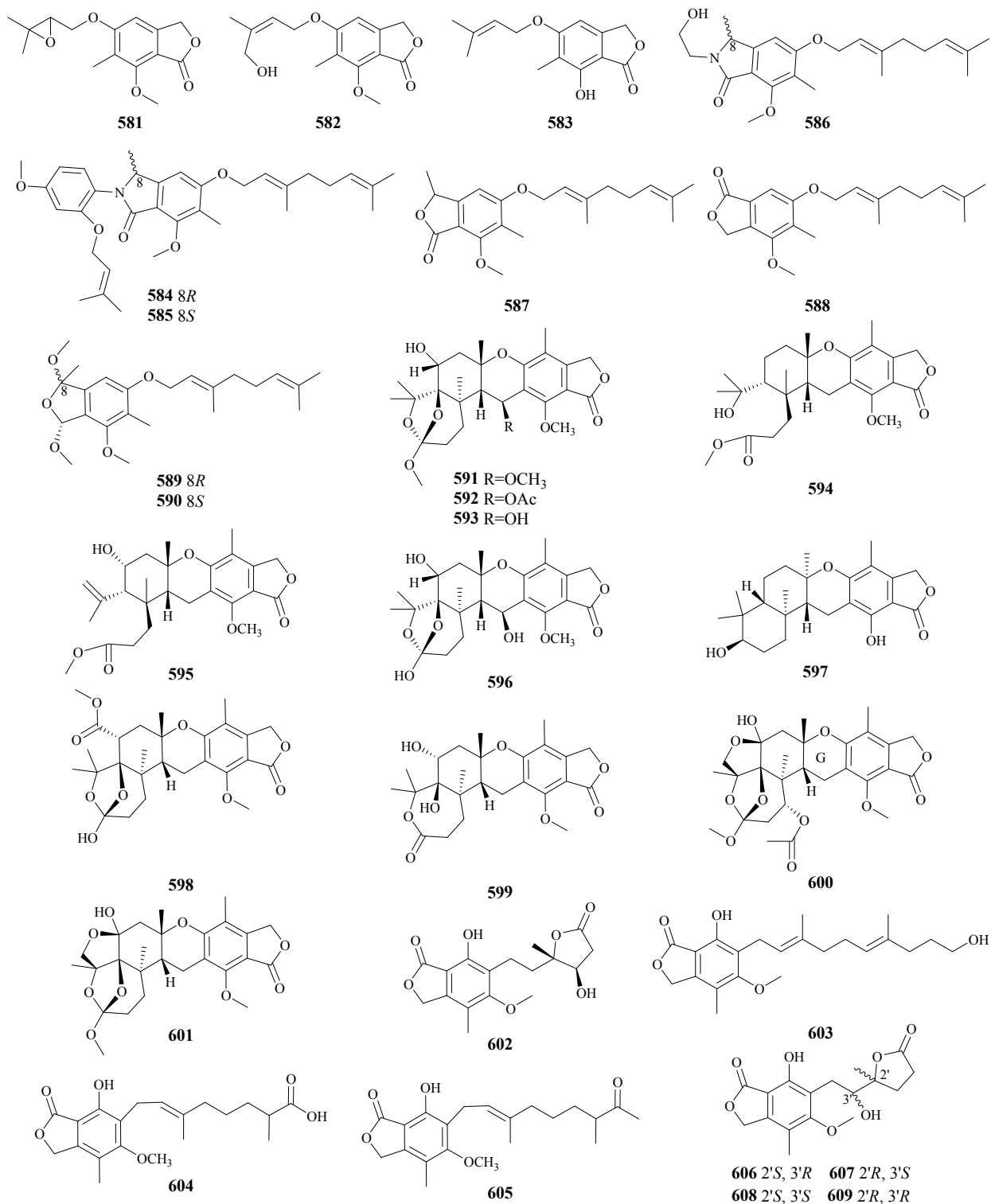


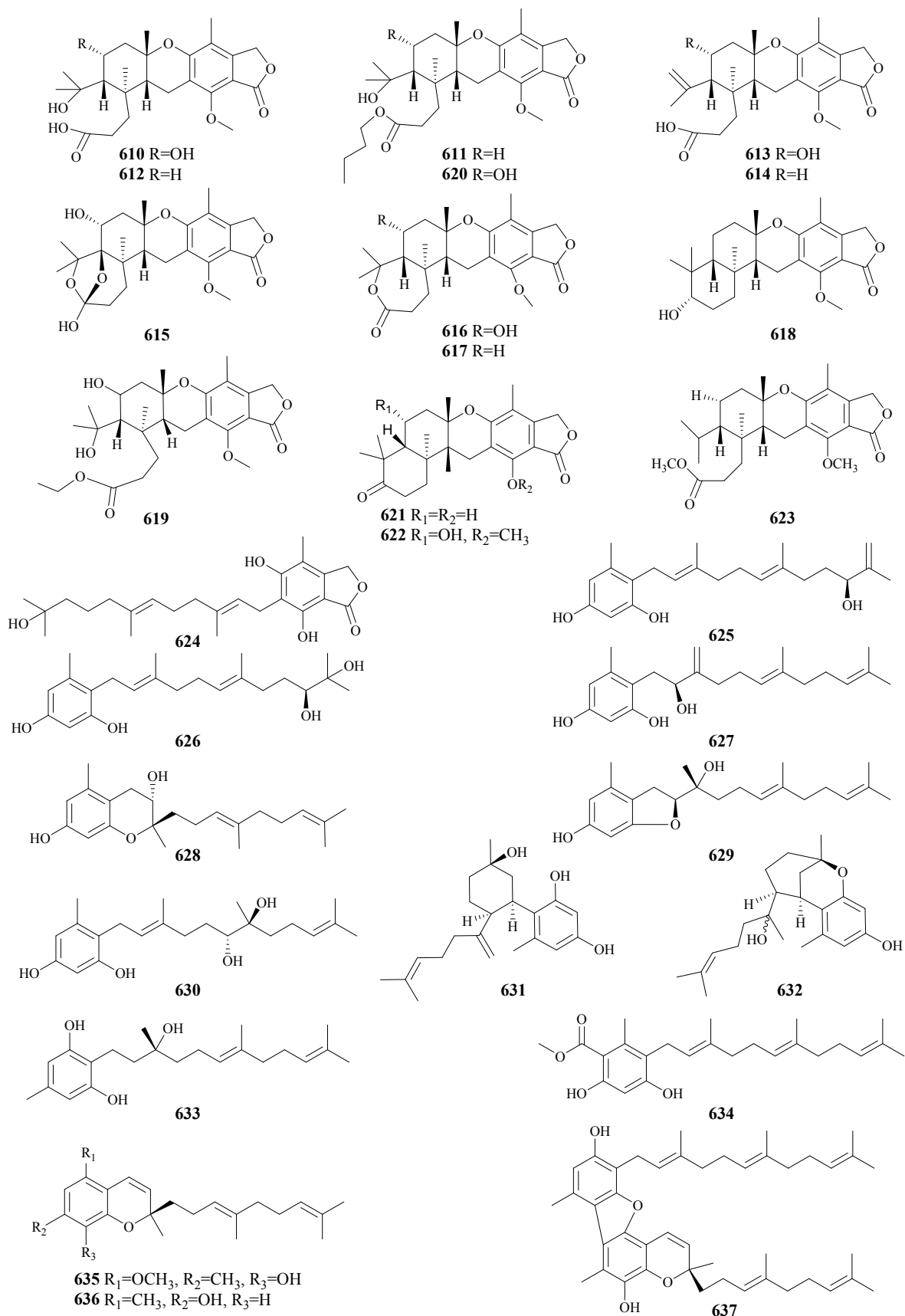
578

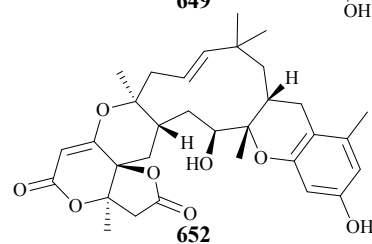
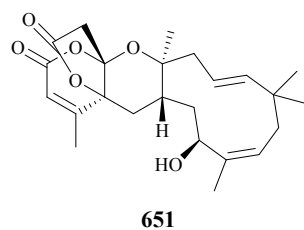
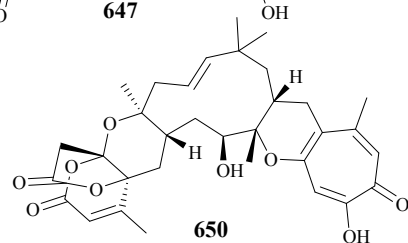
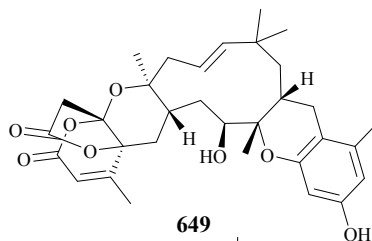
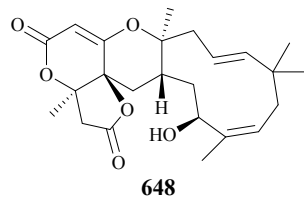
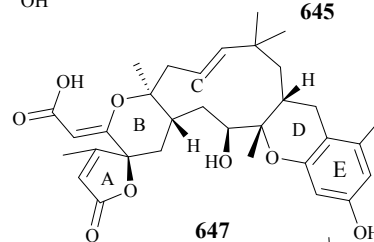
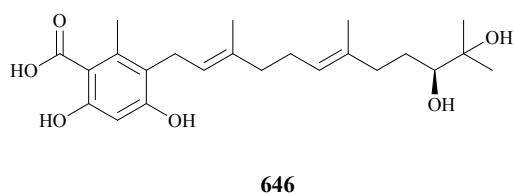
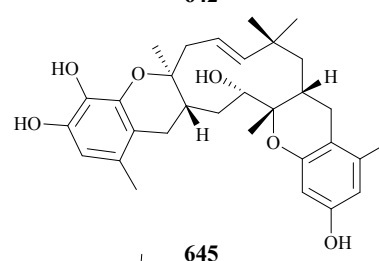
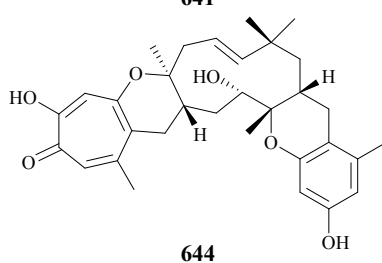
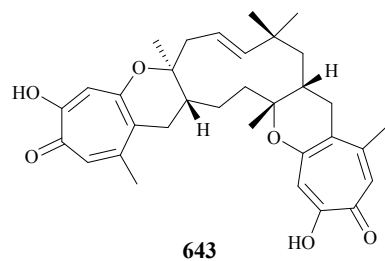
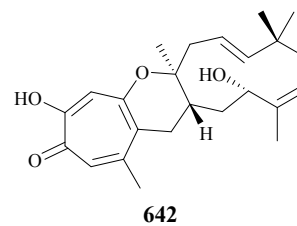
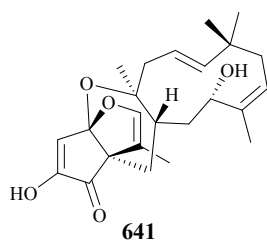
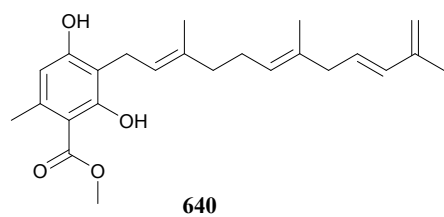
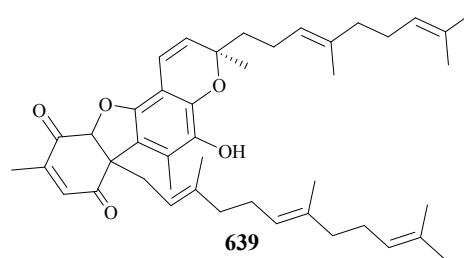
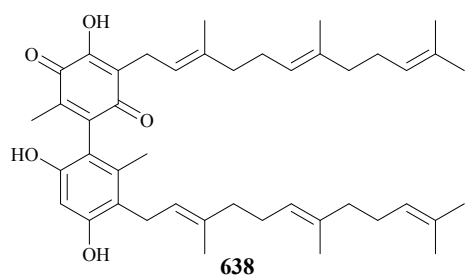


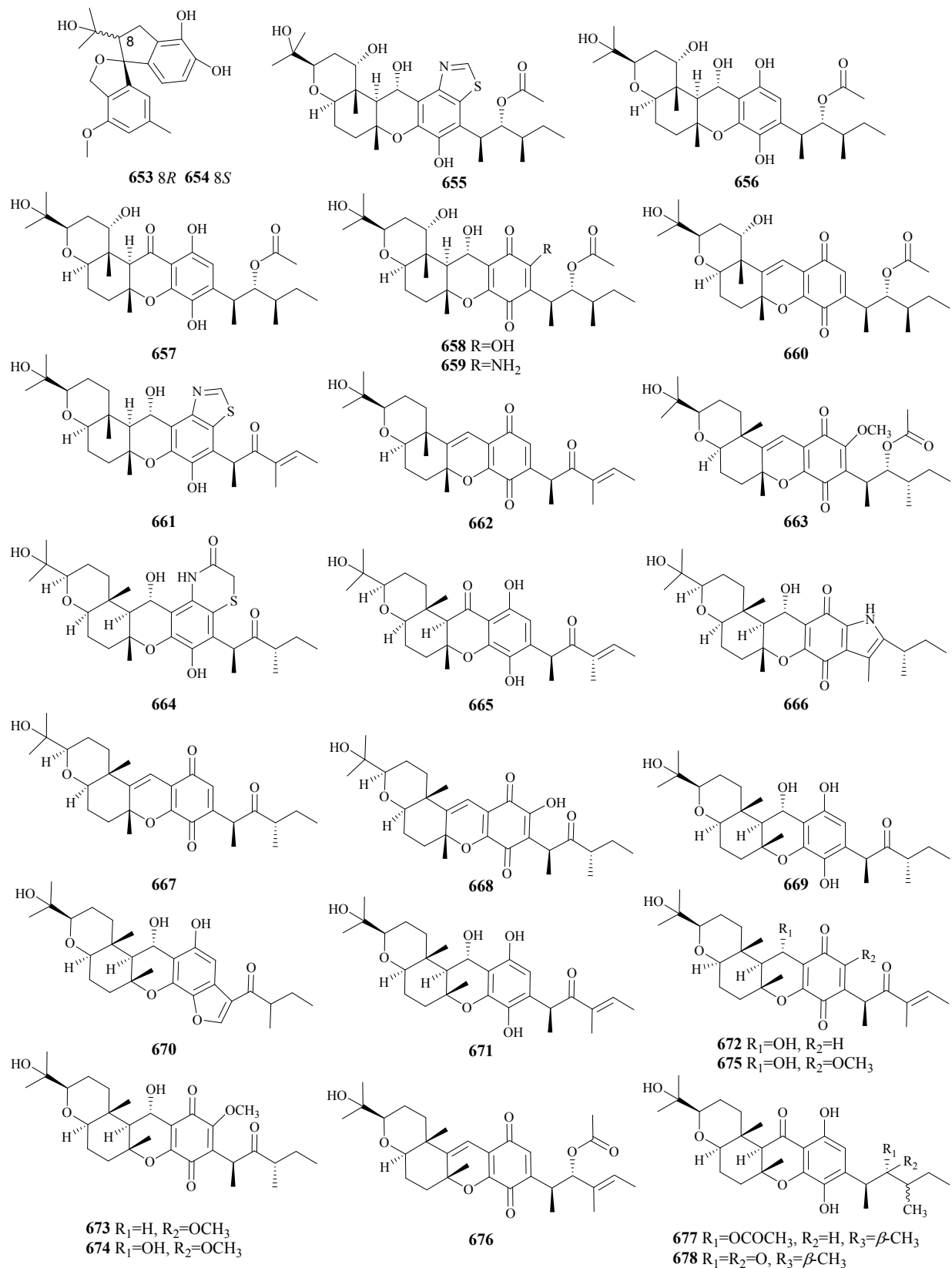
579 R=OH
580 R=keto

2.4. Other tetraketide-terpenoids (581-652)

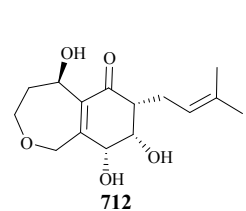
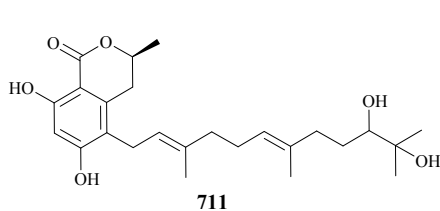
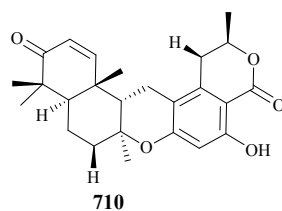
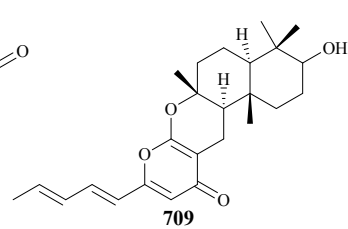
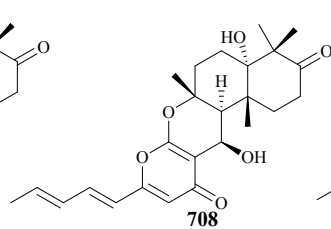
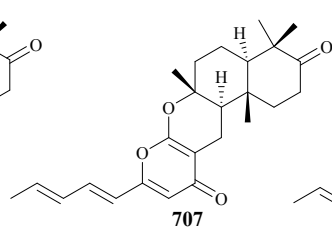
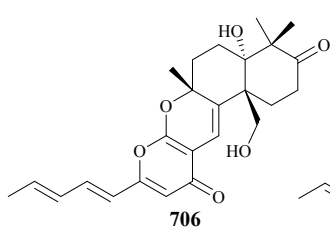
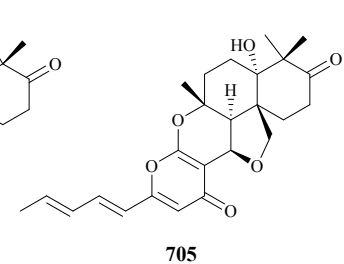
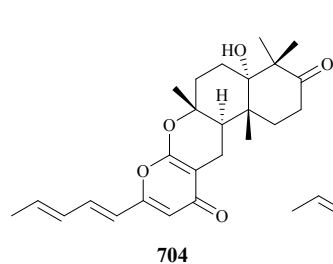
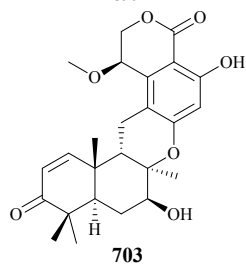
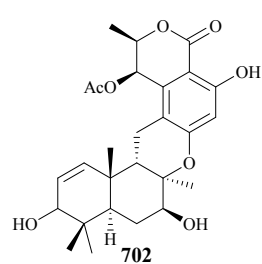
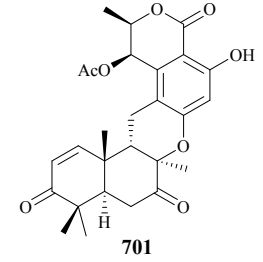
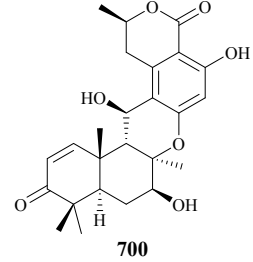
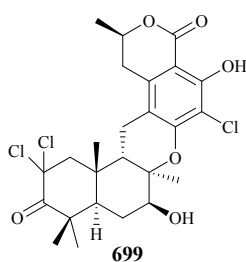
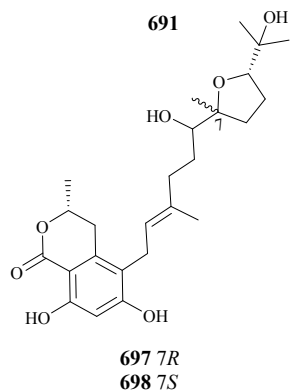
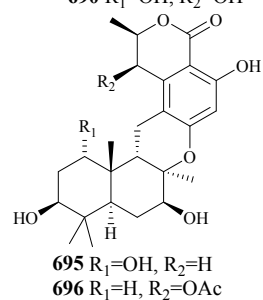
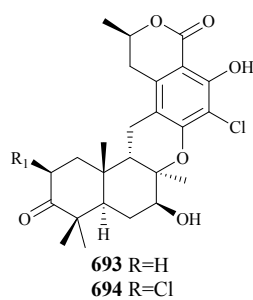
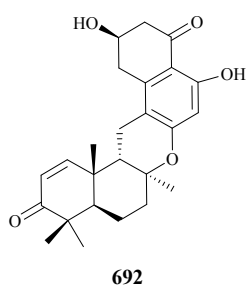
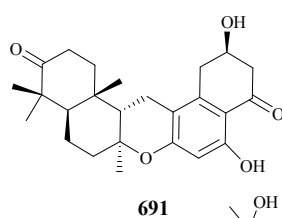
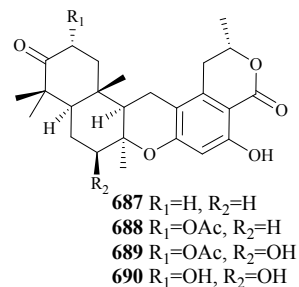
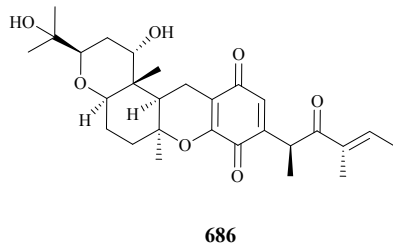
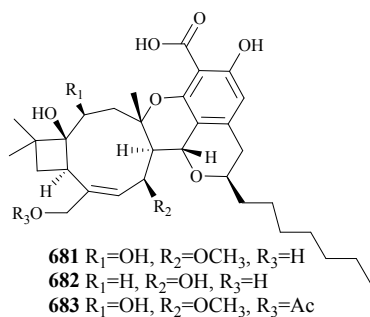
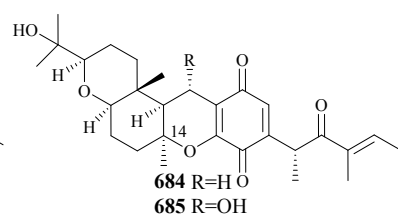
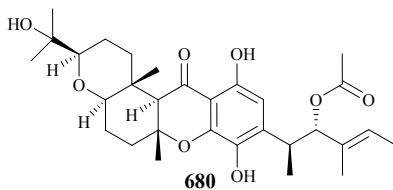
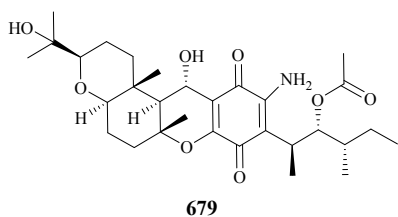


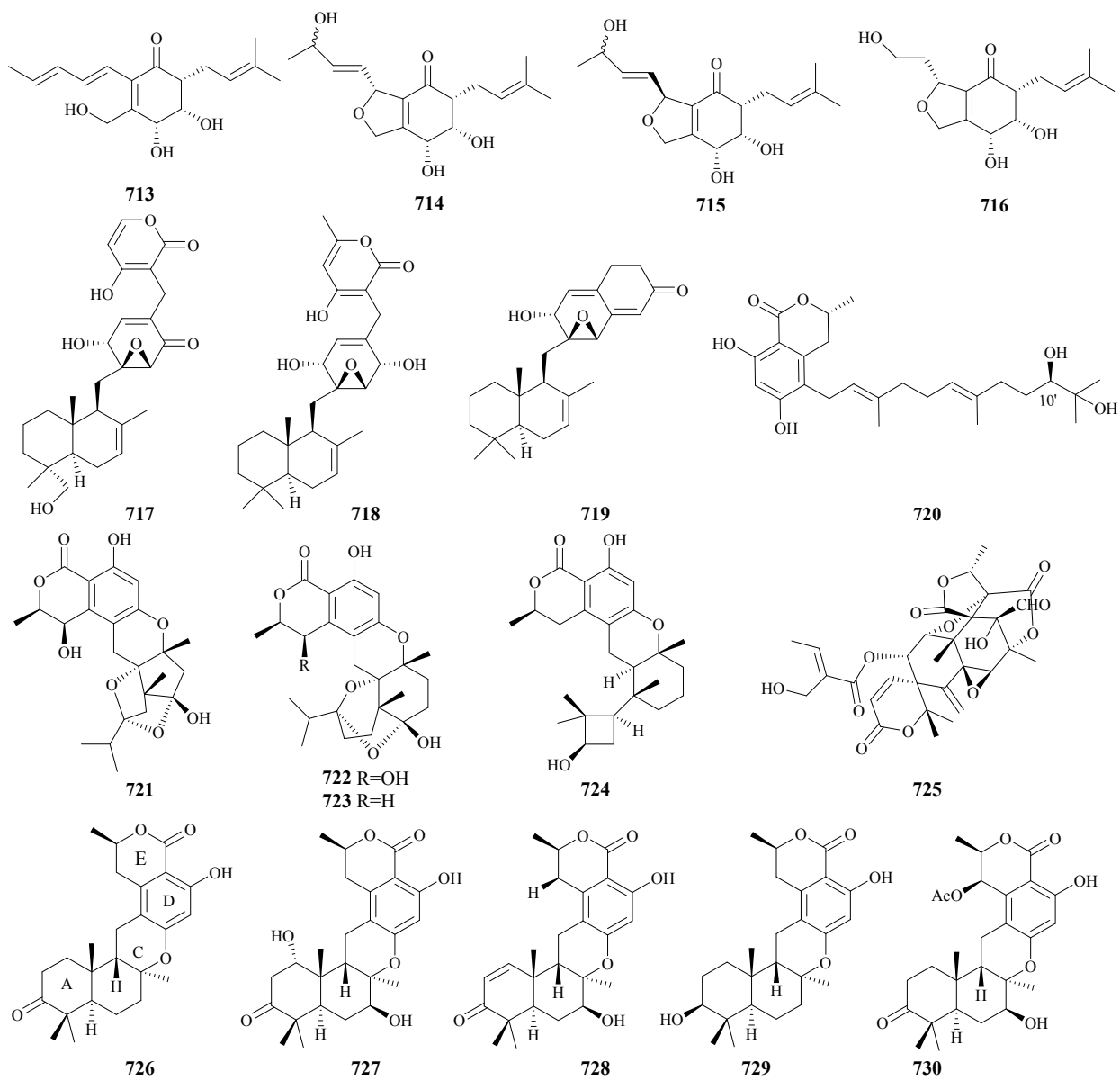


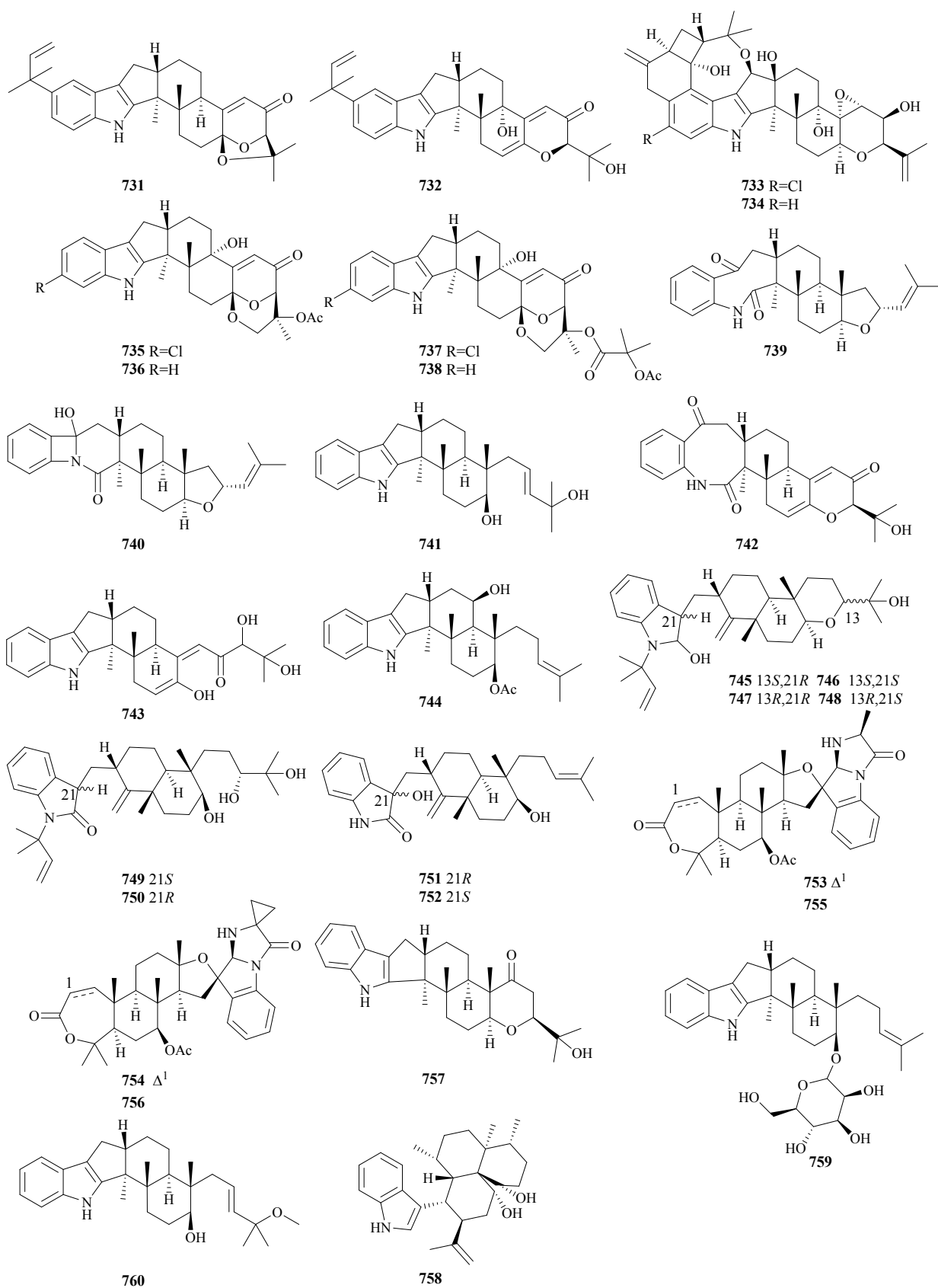




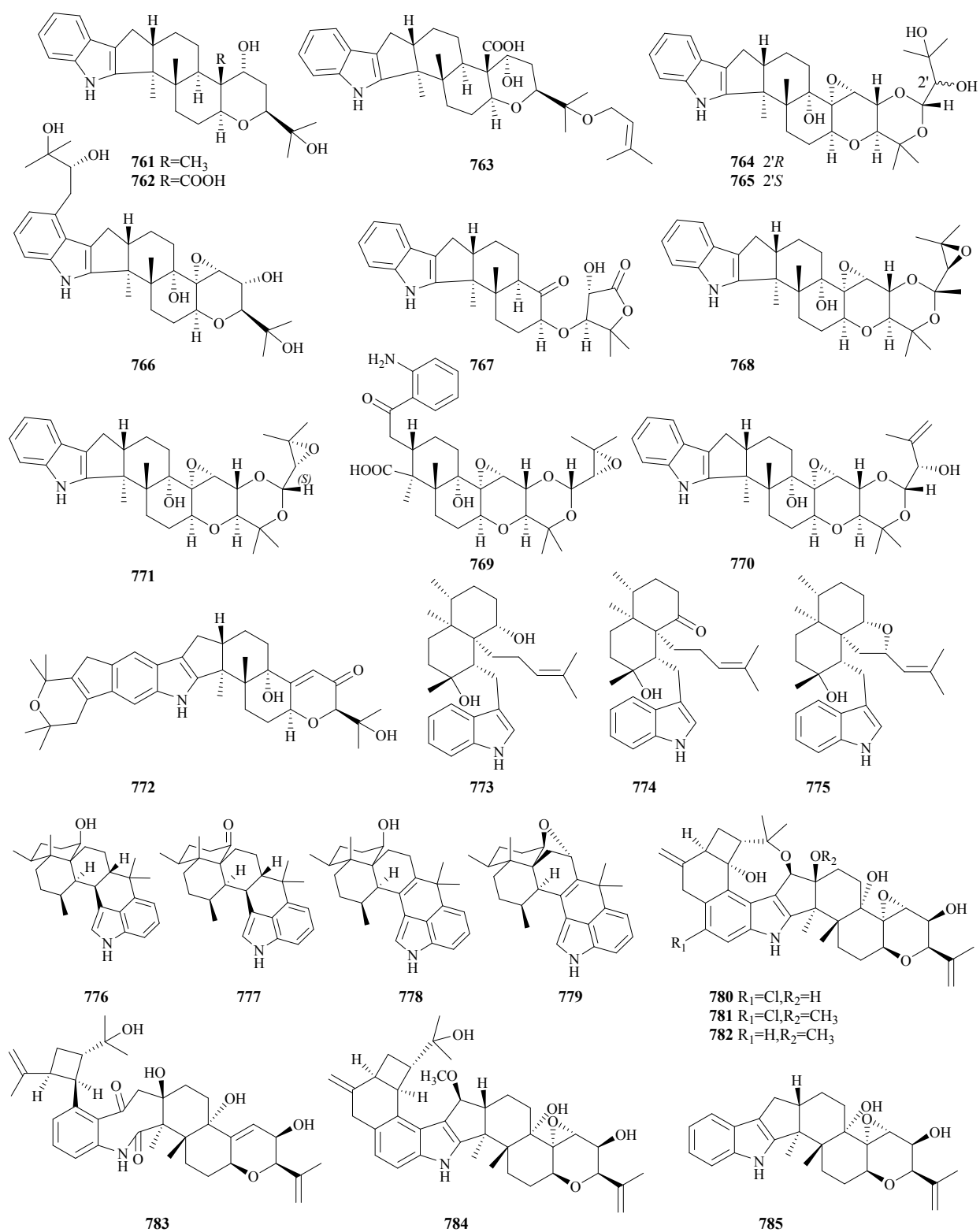
3. Other polyketide-terpenoids (653-730)

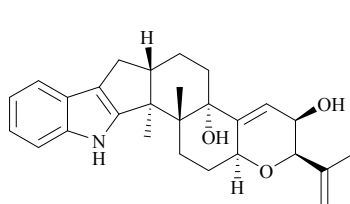




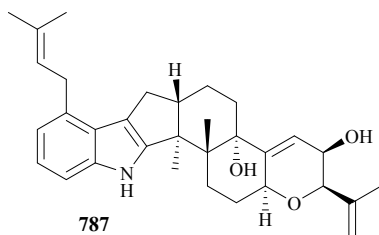


4. Indole-terpenoids (731-844)

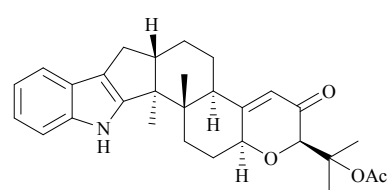




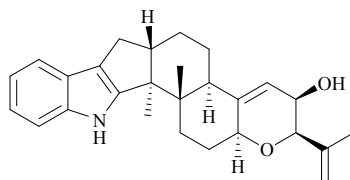
786



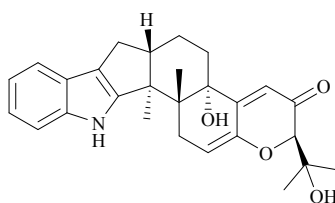
787



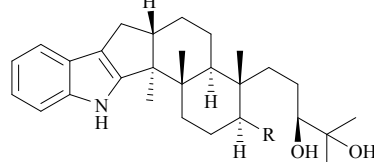
788



789

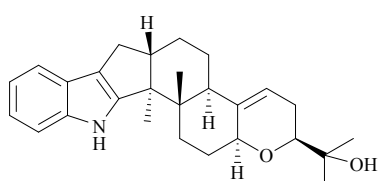


790

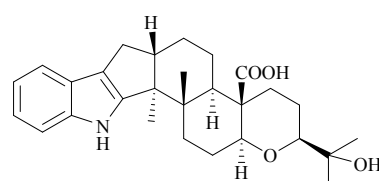


791 R=OH

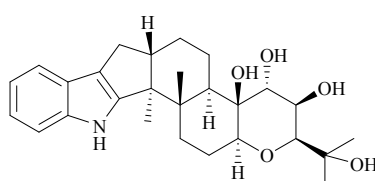
792 R=OAc



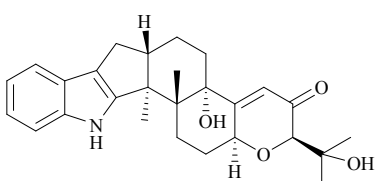
793



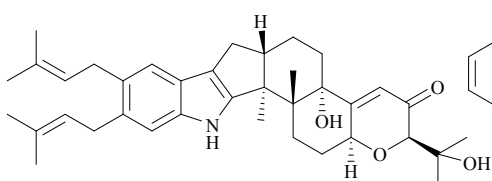
794



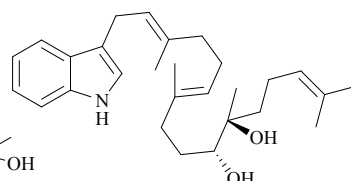
795



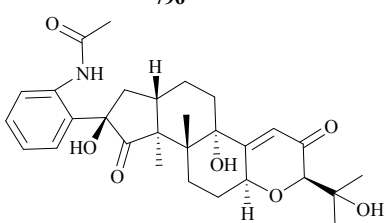
796



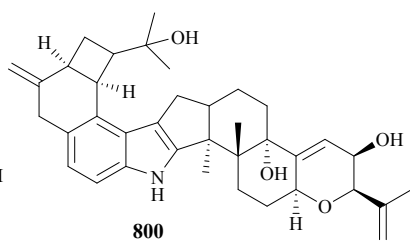
797



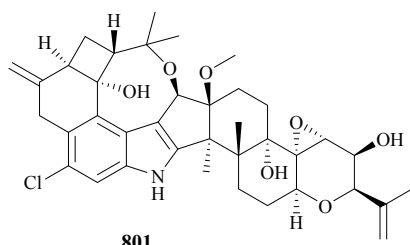
798



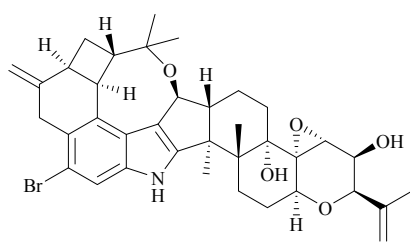
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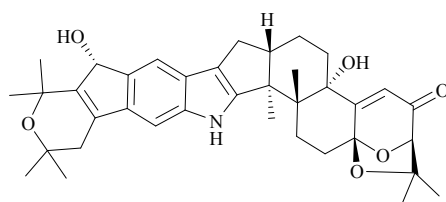
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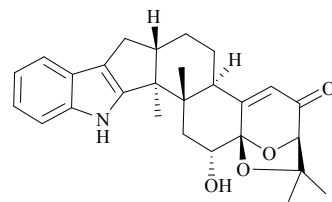
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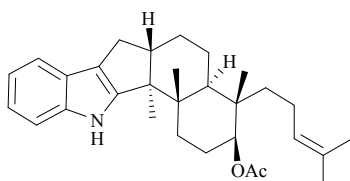
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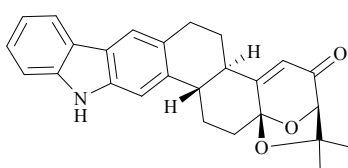
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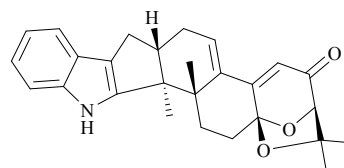
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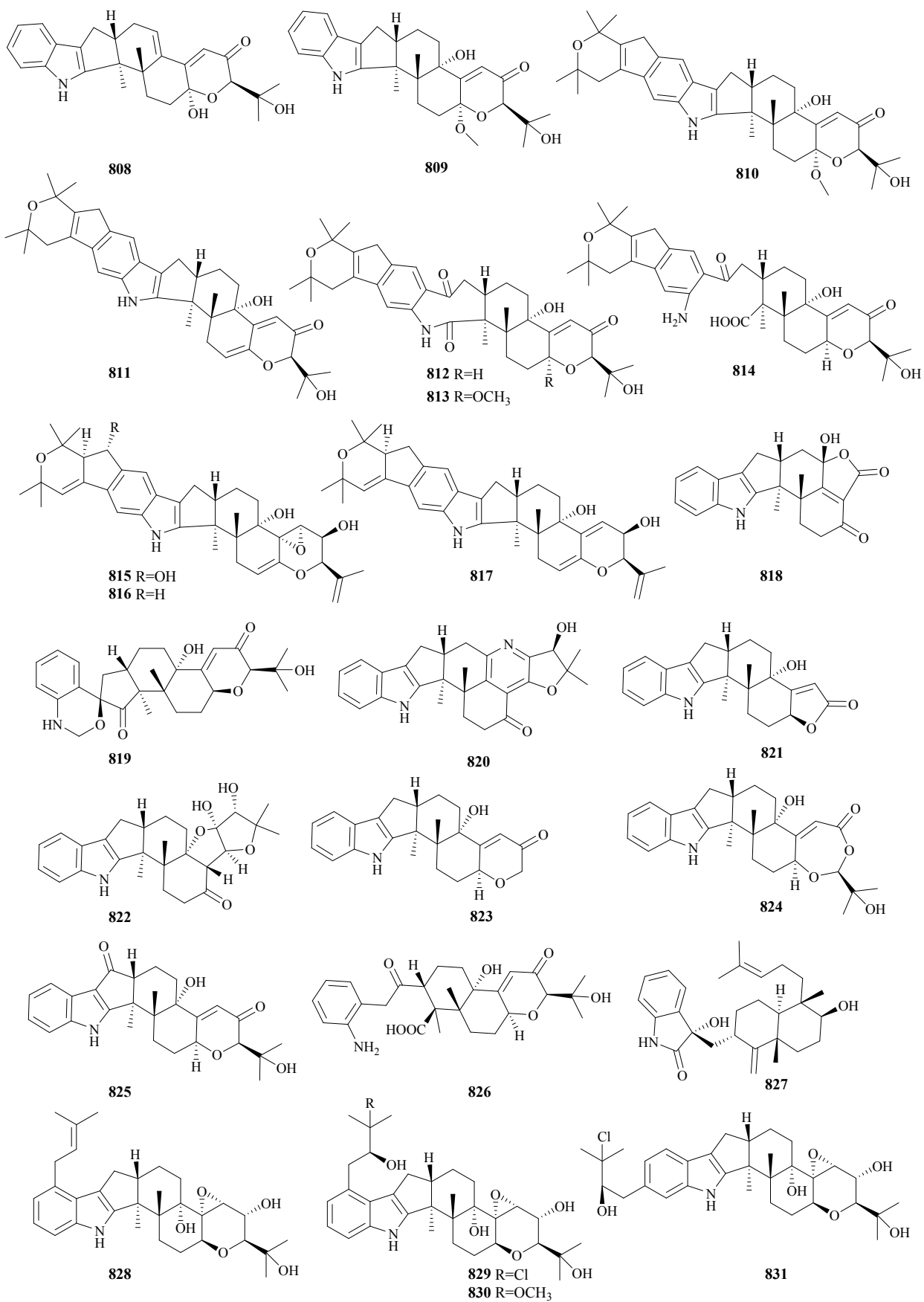
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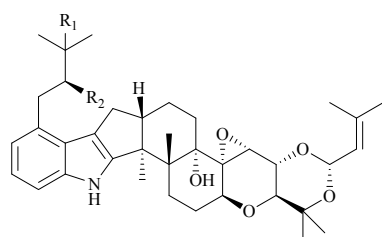


806

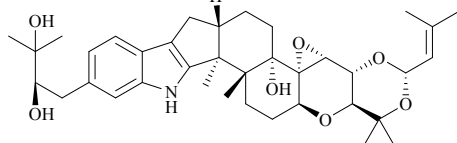


807

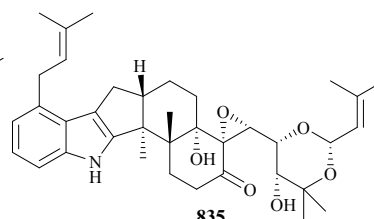




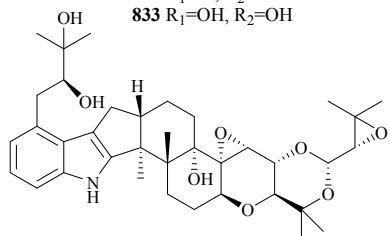
832 $R_1=Cl$, $R_2=OH$
833 $R_1=OH$, $R_2=OH$



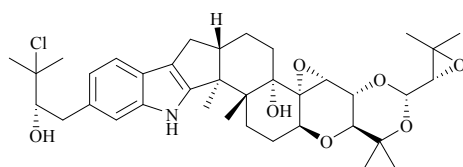
834



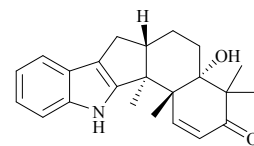
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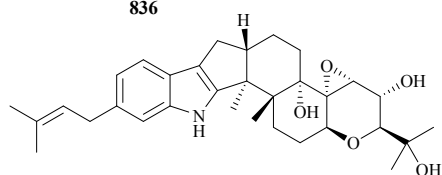
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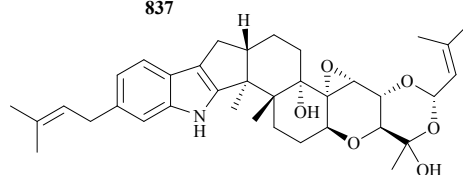
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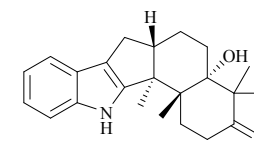
840



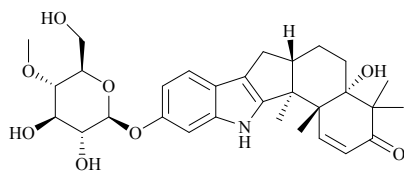
838



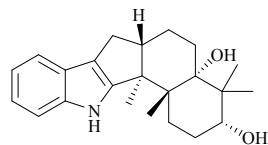
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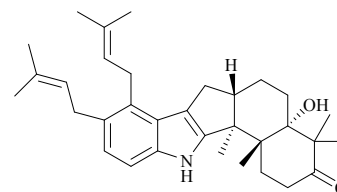
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842

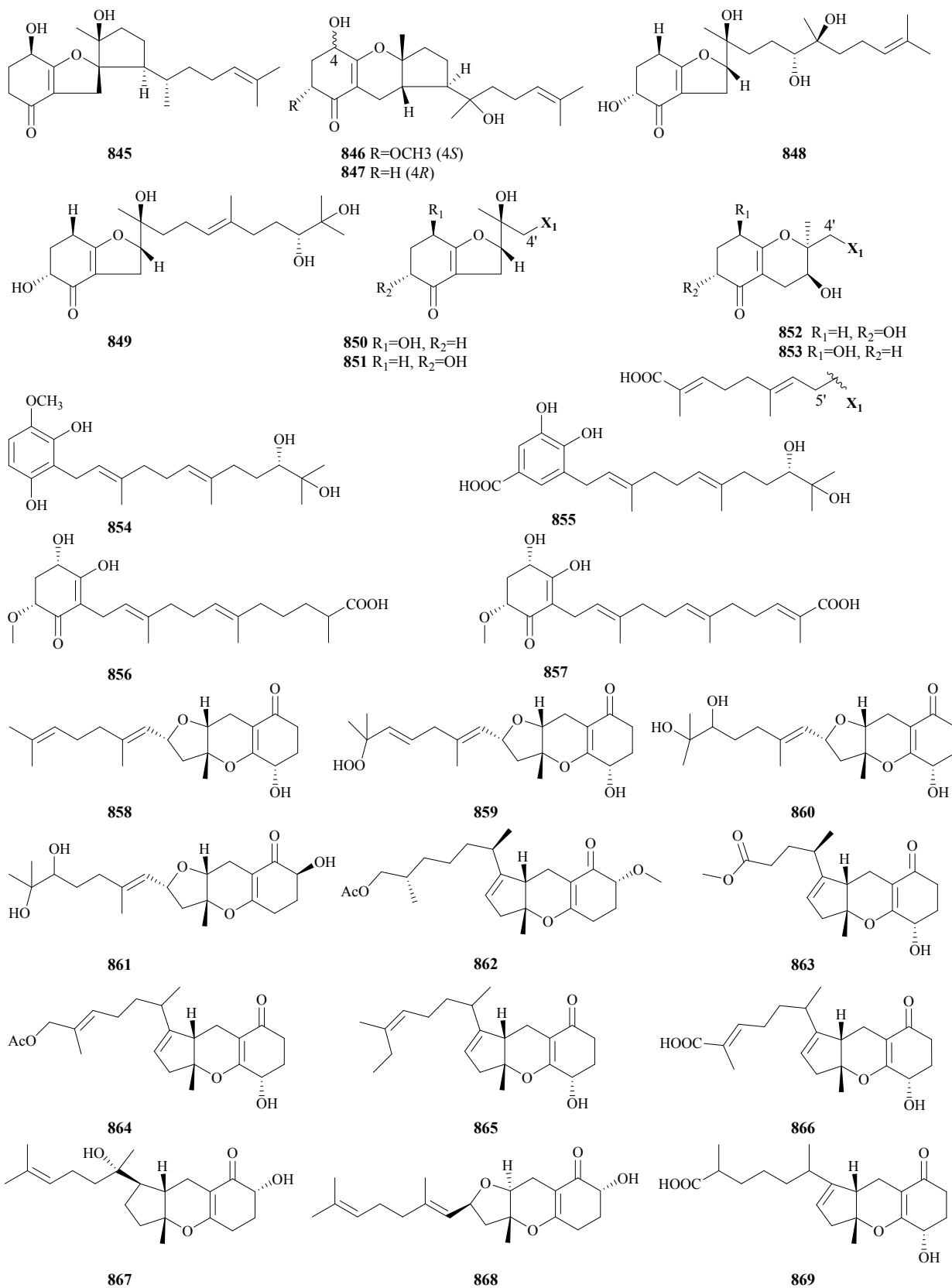


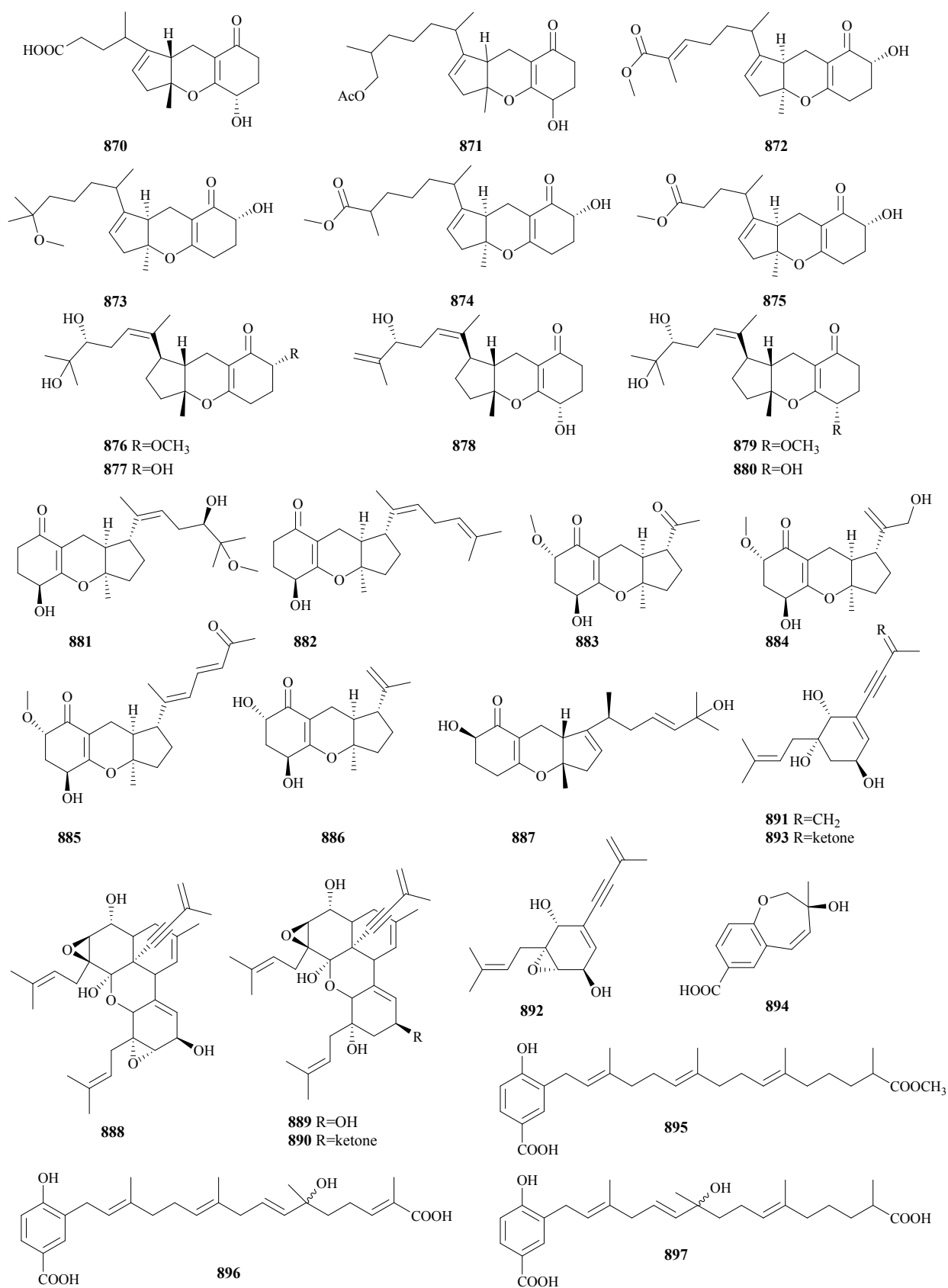
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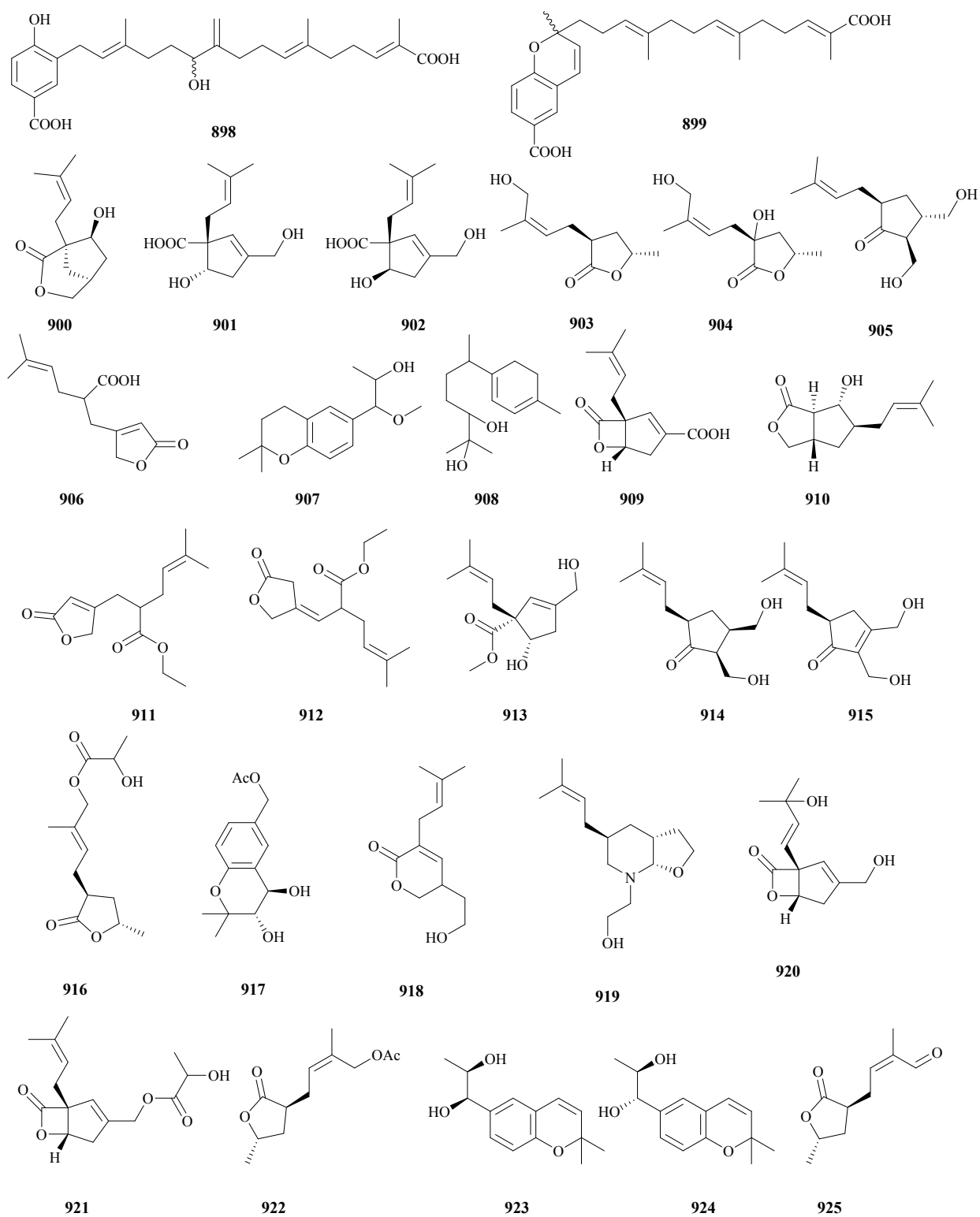


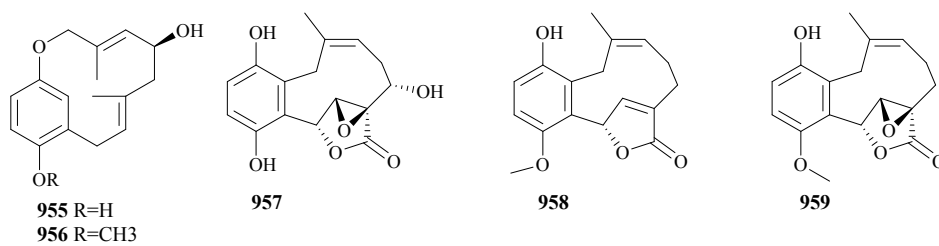
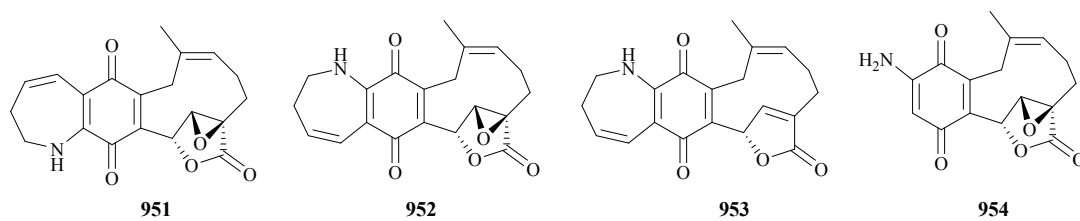
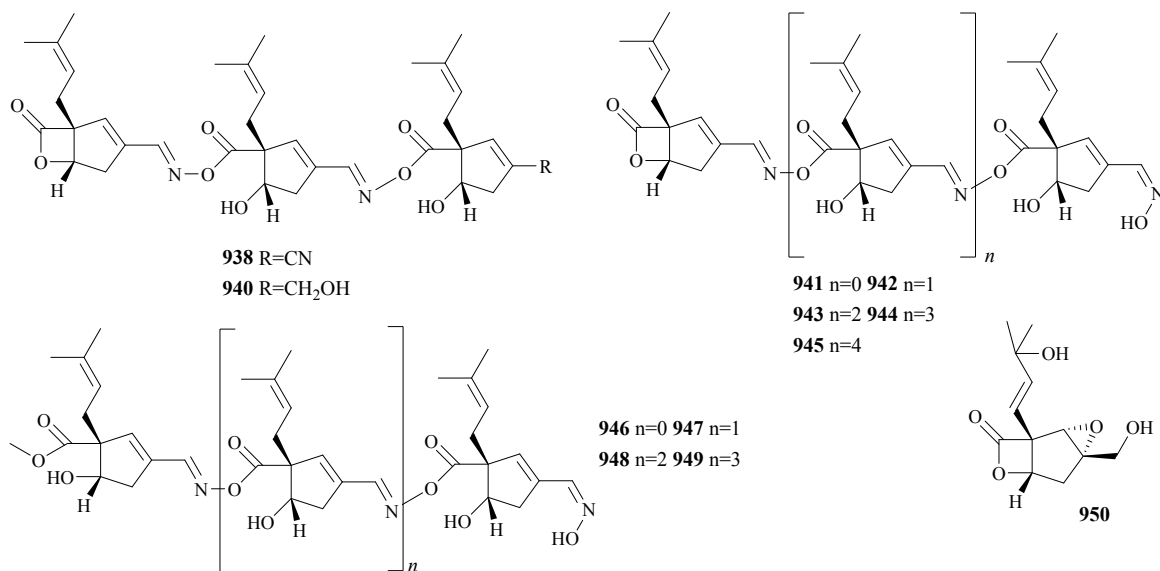
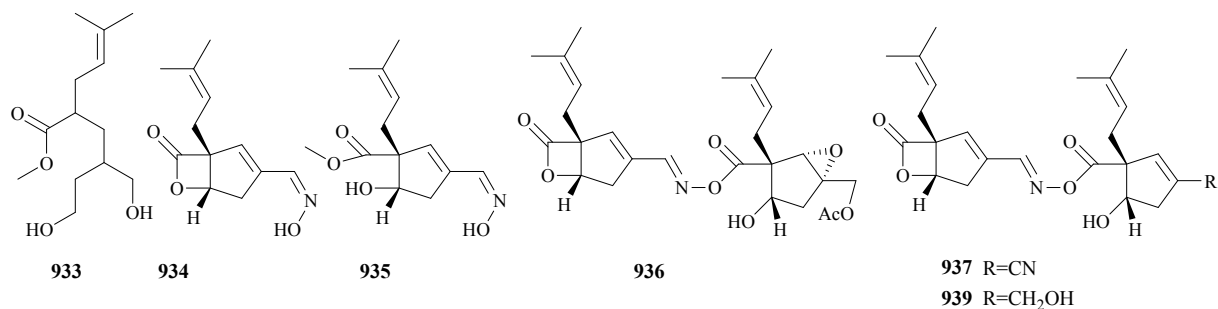
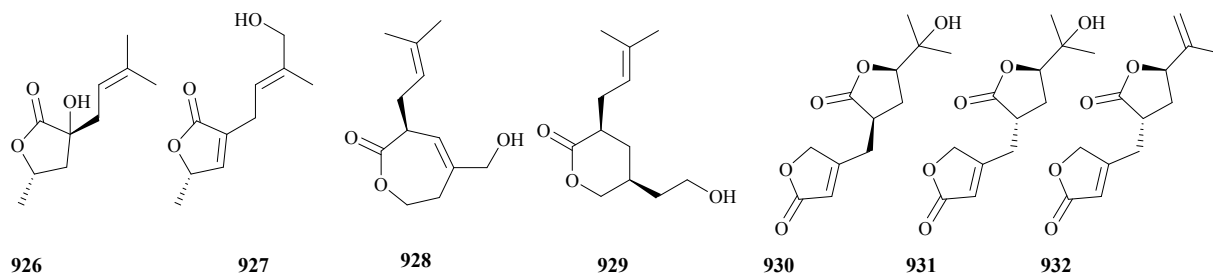
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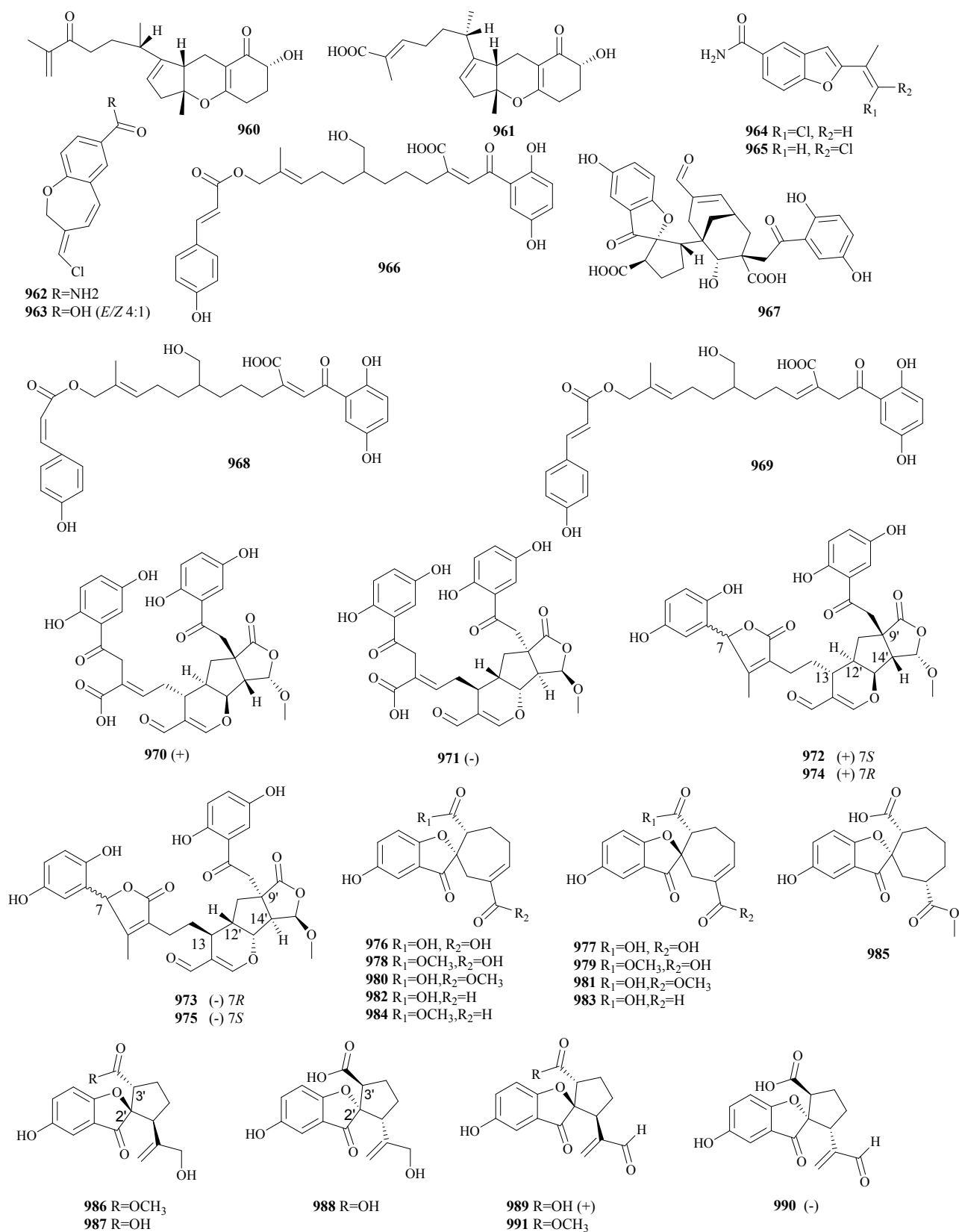
5. Shikimate-terpenoids (845-1544)

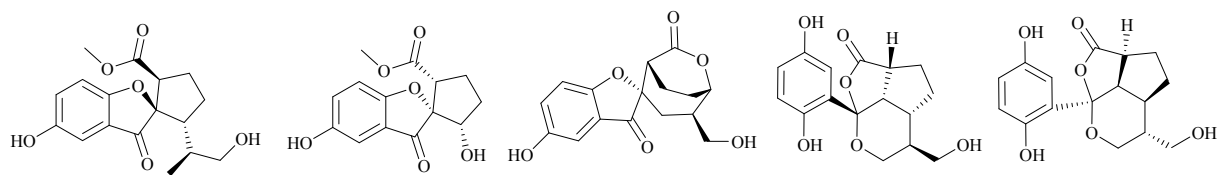












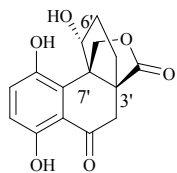
992

993

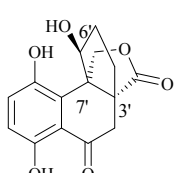
994

995 (+)

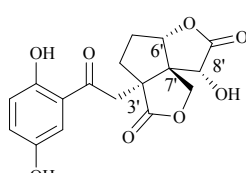
996 (-)



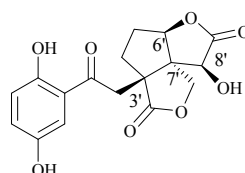
997 (+)



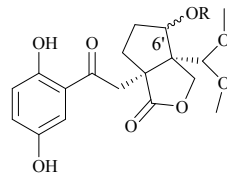
998 (-)



999 (+)



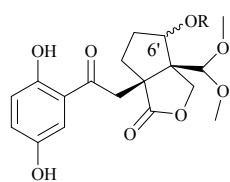
1000 (-)



1001 (+) R=H 3 6'R

1003 (+) R=H 6'S

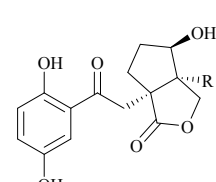
1005 (+) R=Et 6'R



1002 (-) R=H 6'S

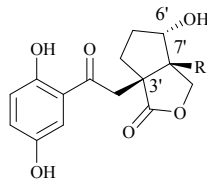
1004 (-) R=H 6'R

1006 (-) R=Et 6'R



1007 (+) R=CH₂OH

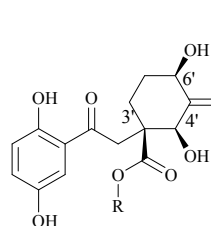
1009 (+) R=COOH



1008 (-) R=CH₂OH

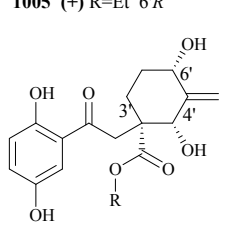
1010 (-) R=COOH

1011 (±) R=Cl



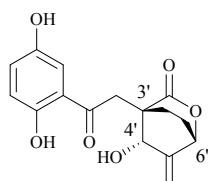
1012 (+) R=H

1014 (+) R=CH₃

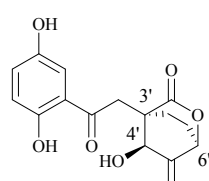


1013 (-) R=H

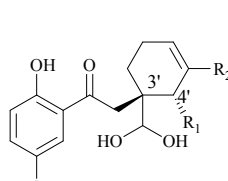
1015 (-) R=CH₃



1016 (+)

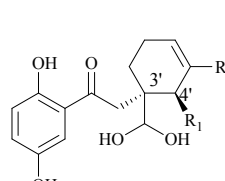


1017 (-)



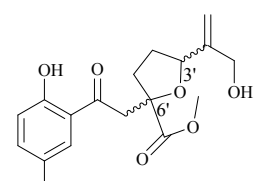
1018 (+) R₁=OH, R₂=CH₂OH

1021 (-) R₁=H, R₂=CHO



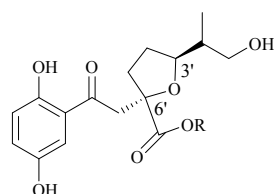
1019 (-) R₁=OH, R₂=CH₂OH

1020 (+) R₁=H, R₂=CHO



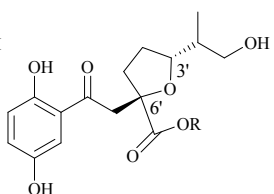
1022 (+) 3'S, 6'S

1023 (-) 3'R, 6'R



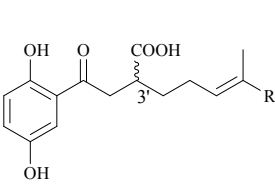
1024 (+) R=CH₃

1026 (+) R=H



1025 (-) R=CH₃

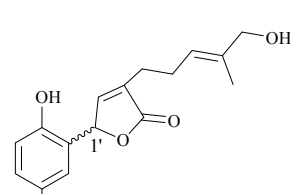
1027 (-) R=H



1028 (+) R=CH₂OH 3'R

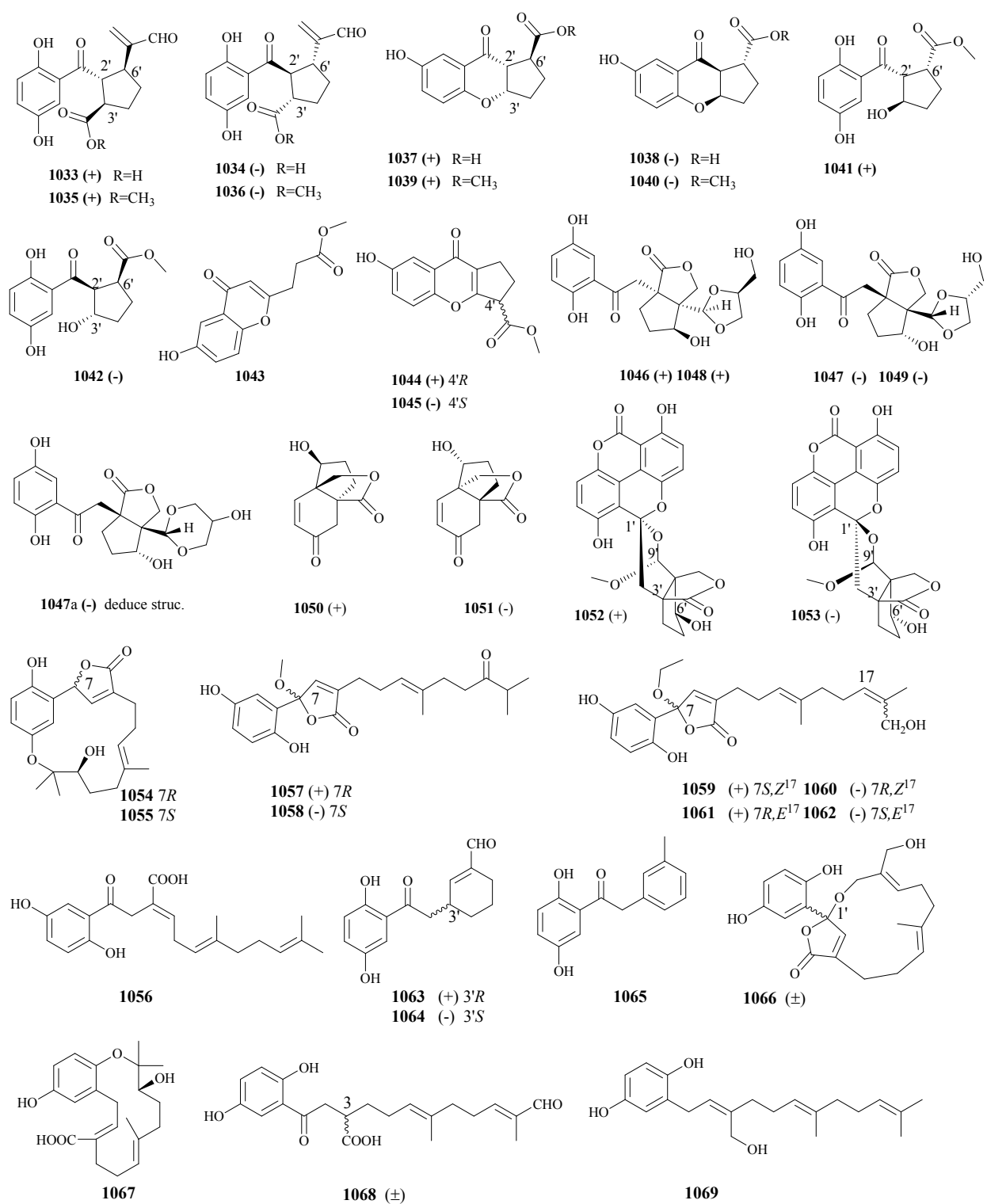
1029 (-) R=CH₂OH 3'S

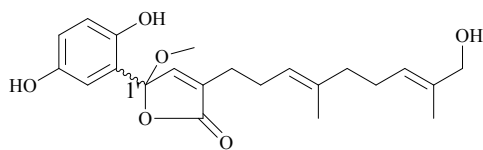
1030 (+) R=COOH 3'R



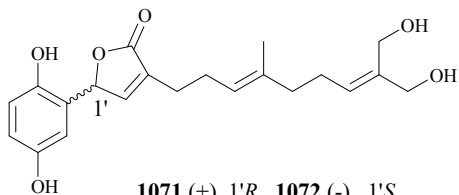
1031 (+) 1'R

1032 (-) 1'S

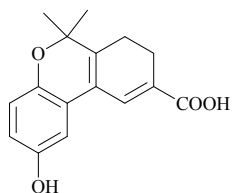




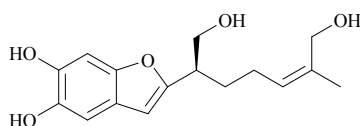
1070 ($\pm$)



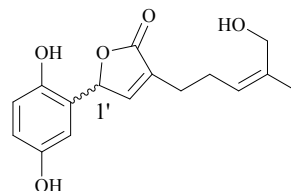
1071 (+) 1'R **1072** (-) 1'S



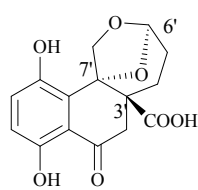
1073



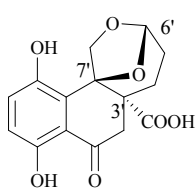
1074



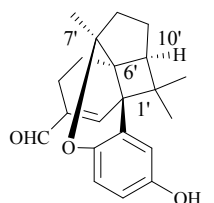
1075 (+) 1'R **1076** (-) 1'S



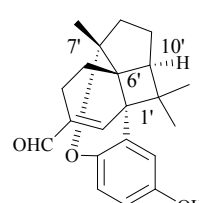
1077 (+)



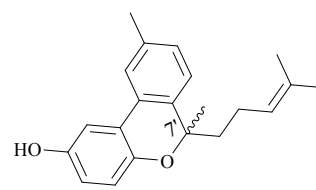
1078 (-)



1079 (+)

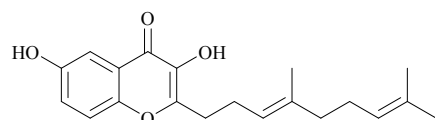


1080 (-)

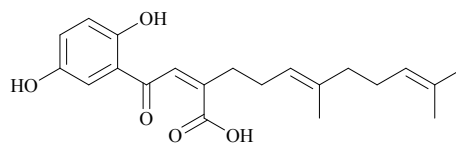


1084 (+) 7'R

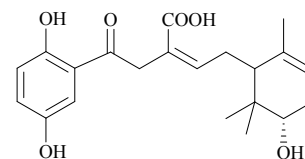
1085 (-) 7'S



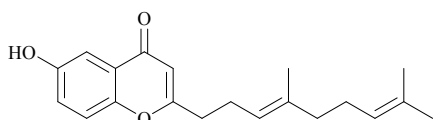
1081



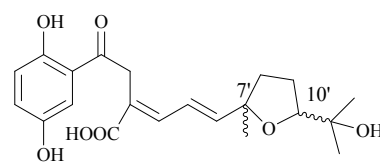
1082



1089

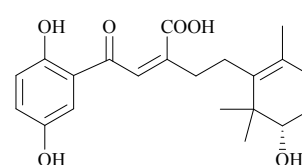


1083

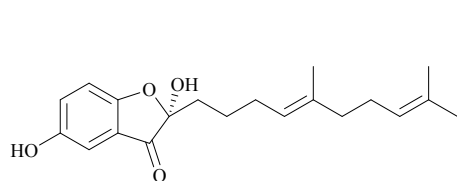


1086 (+) 7'R,10'S **1087** (-) 7'S,10'R

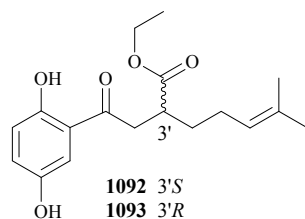
1088* 7'R,10'R



1090* 10'R

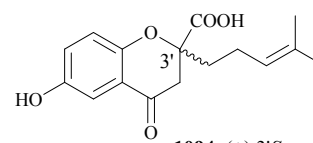


1091* 2'S



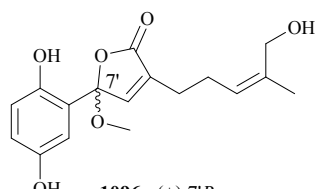
1092 3'S

1093 3'R



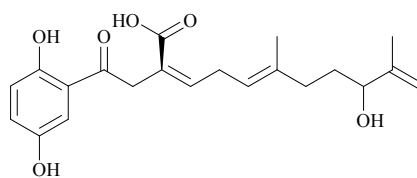
1094 (+) 3'S

1095 (-) 3'R

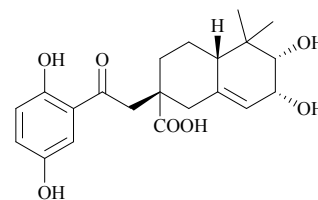


1096 (+) 7'R

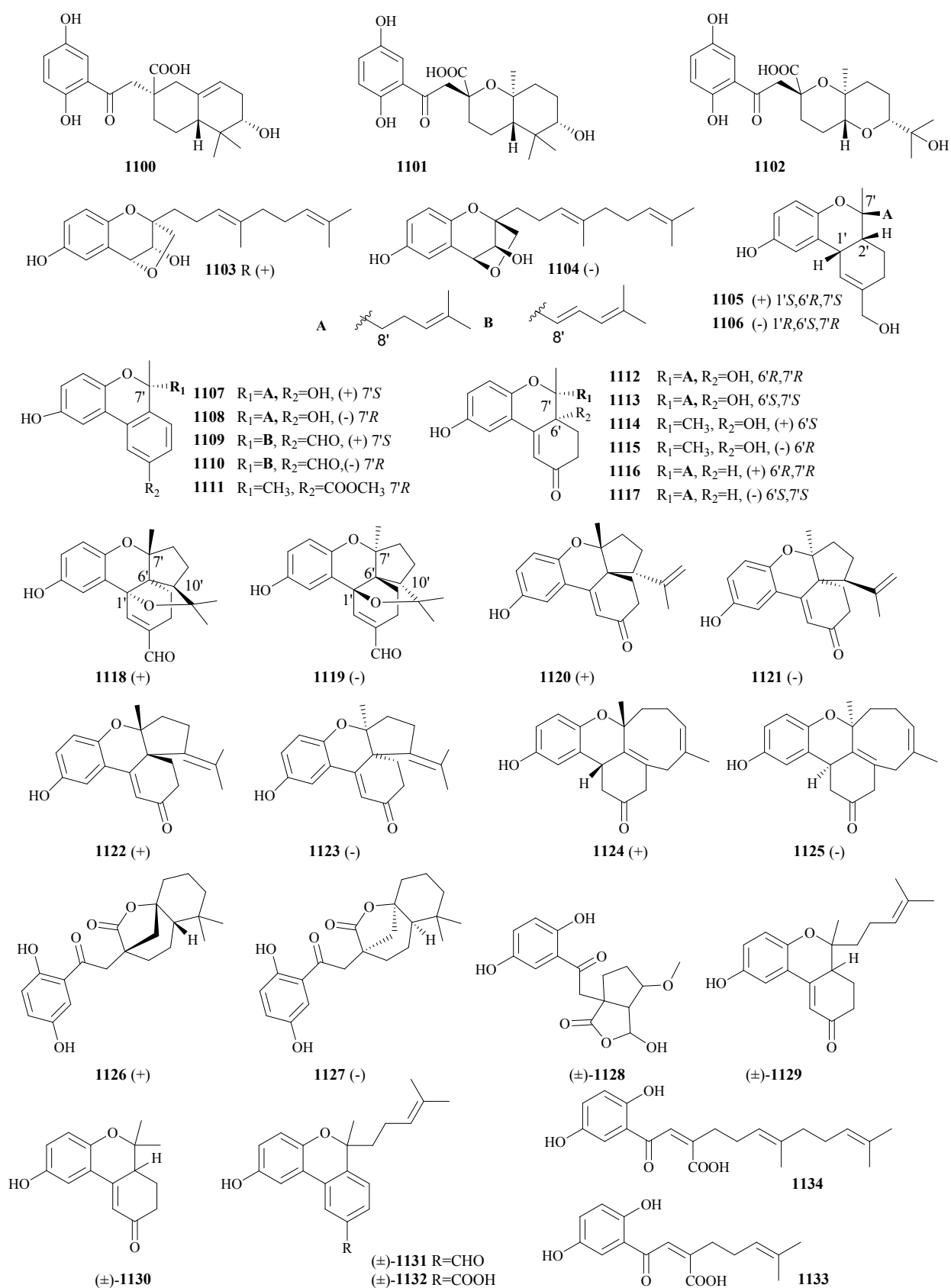
1097 (-) 7'S

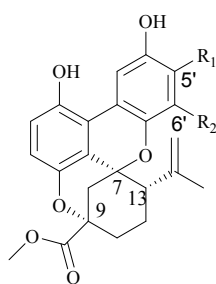


1098



1099





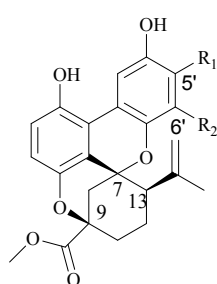
$R_1=H, R_2=A_1, R_3=OAc$ **1135**(+)

$R_1=H, R_2=A_1, R_3=H$ **1137** (+)

$R_1=H, R_2=B_1, R_3=OAc$ **1139** (+)

$R_1=B_1, R_2=H, R_3=OH$ **1141**(+)

$R_1=A_1, R_2=H, R_3=OAc$ **1143** (+)



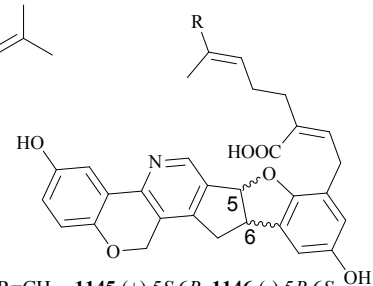
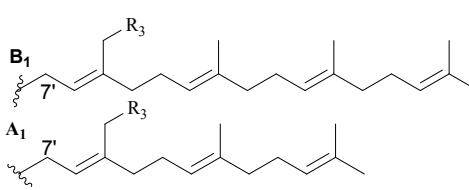
$R_1=H, R_2=A_1, R_3=OAc$ **1136** (-)

$R_1=H, R_2=A_1, R_3=H$ **1138** (-)

$R_1=H, R_2=B_1, R_3=OAc$ **1140** (-)

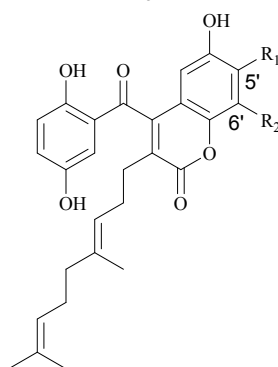
$R_1=B_1, R_2=H, R_3=OH$ **1142** (-)

$R_1=A_1, R_2=H, R_3=OAc$ **1144** (-)



$R=CH_3$ **1145** (+) $5S,6R$ **1146** (-) $5R,6S$

$R=CHO$ **1147** (+) $5S,6R$ **1148** (-) $5R,6S$



$R_1=H, R_2=A_2, R_3=COOH$ **1151**

$R_1=H, R_2=A_2, R_3=CH_2OH$ **1152**

$R_1=A_2, R_2=H, R_3=CH_2OH$ **1153**

$R_1=B_2, R_2=H$ **1154**

$R_1=H, R_2=A_2, R_3=CH_2OH$ **1149**

$R_1=H, R_2=A_2, R_3=CH_3$ **1155**

$R_1=A_2, R_2=H, R_3=CH_2OH$ **1150**

$R_1=A_2, R_2=H, R_3=COOH$ **1156**

$R_1=H, R_2=A_2, R_3=CH_2OH$ **1157** (+)

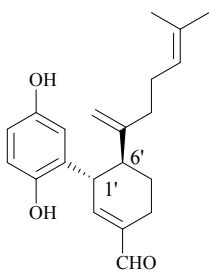
$R_1=H, R_2=A_2, R_3=COOH$ **1159** (+)

$R_1=A_2, R_2=H, R_3=COOH$ **1161** (-)

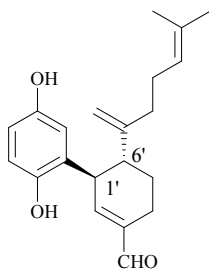
$R_1=H, R_2=A_2, R_3=CH_2OH$ **1158** (-)

$R_1=H, R_2=A_2, R_3=COOH$ **1160** (-)

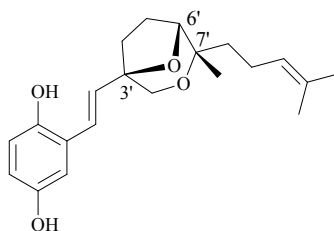
$R_1=A_2, R_2=H, R_3=COOH$ **1162** (+)



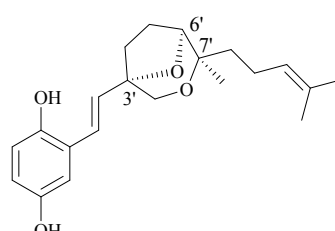
1163 (+)



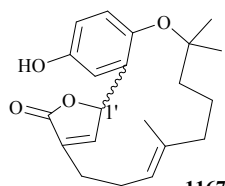
1164 (-)



1165 (+)

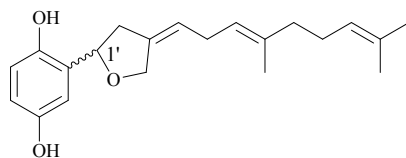


1166 (-)



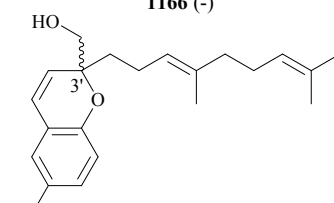
1167 1'R

1168 1'S



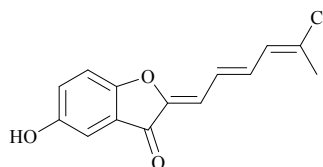
1169 1'R

1170 1'S

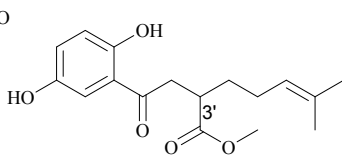


1171 3'R

1172 3'S

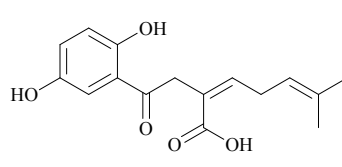


1173

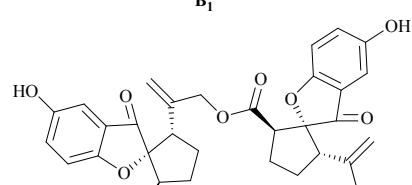
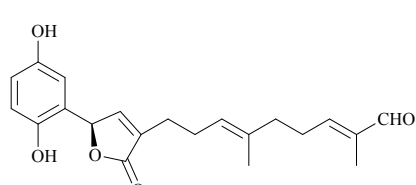
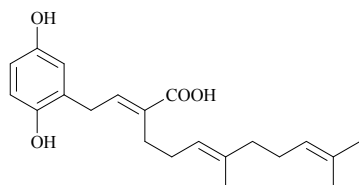
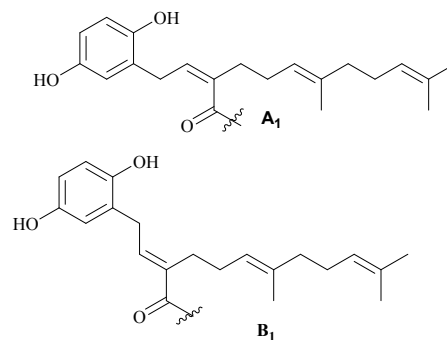
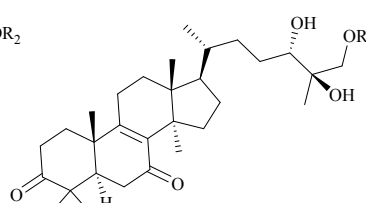
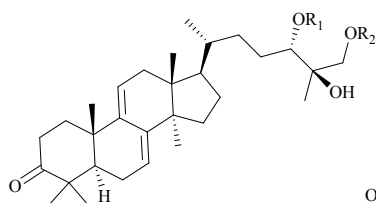
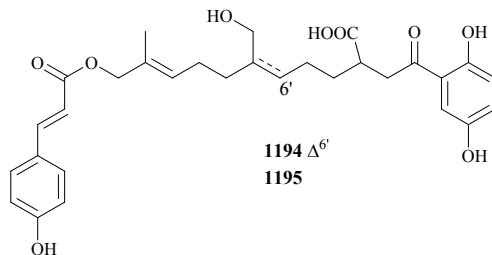
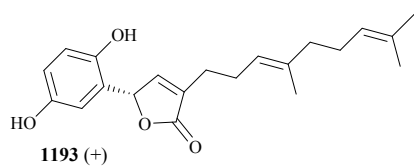
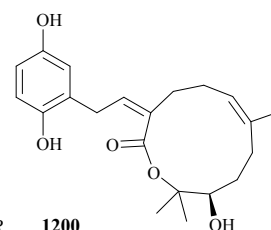
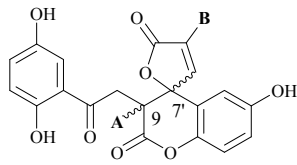
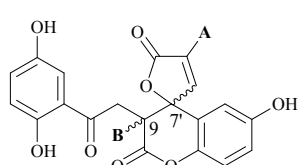
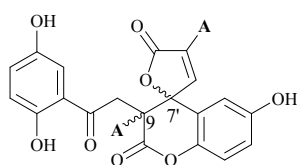
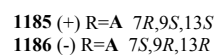
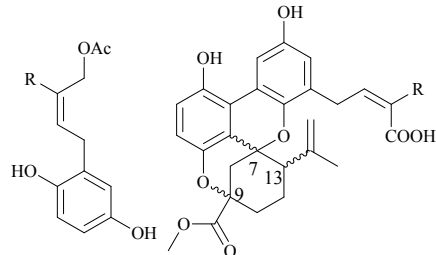
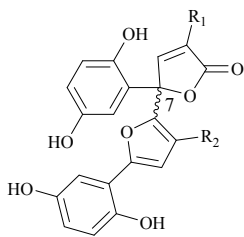
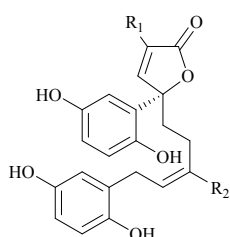
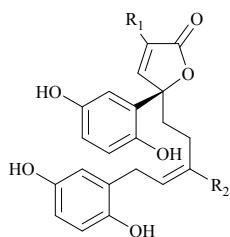
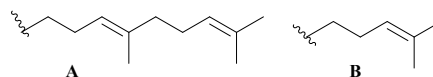
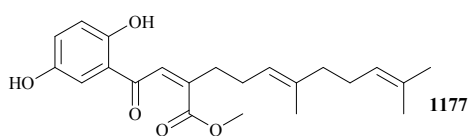


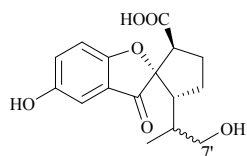
1174 (+) 3'R

1175 (-) 3'S



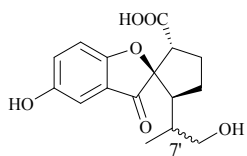
1176





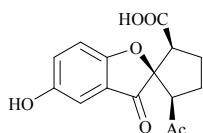
1206 (+) 7'S

1208 (+) 7'R

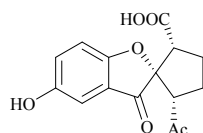


1207 (-) 7'R

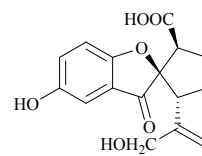
1209 (-) 7'S



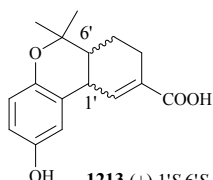
1210 (+)



1211 (-)

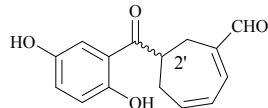


1212



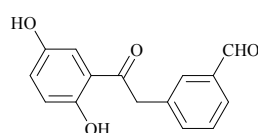
1213 (+) 1'S,6'S

1214 (-) 1'R,6'R

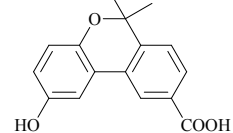


1215 (+) 2'S

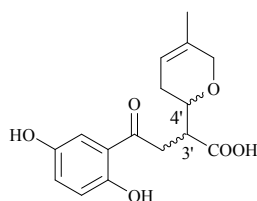
1216 (-) 2'R



1217



1218

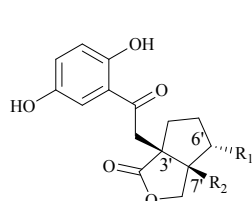


1219 (+) 3'R,4'R

1220 (-) 3'S,4'S

1221 (+) 3'R,4'S

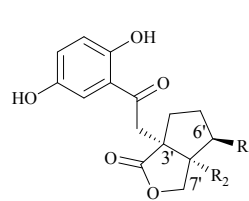
1222 (-) 3'S,4'R



1223 R₁=OEt, R₂=CHO (+)

1225 R₁=OH, R₂=CHO (+)

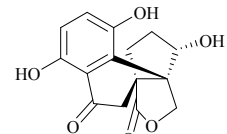
1227 R₁=OH, R₂=CH(OEt)₂ (+)



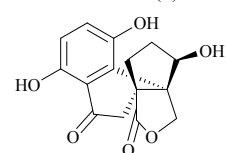
1224 R₁=OEt, R₂=CHO (-)

1226 R₁=OH, R₂=CHO (-)

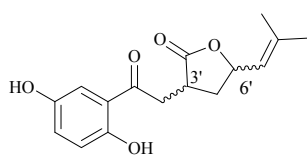
1228 R₁=OH, R₂=CH(OEt)₂ (-)



1229 (+)



1230 (-)

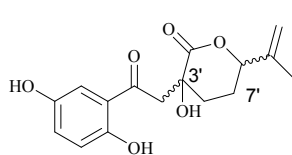


1231 (+) 3'S,6'S

1232 (-) 3'R,6'R

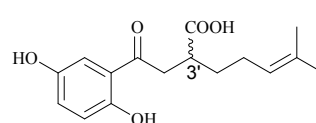
1233 (+) 3'R,6'S

1234 (-) 3'S,6'R



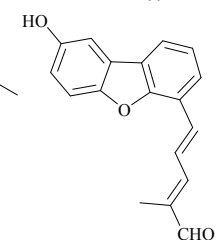
1235 (+) 3'R,7'R

1236 (-) 3'S,7'S

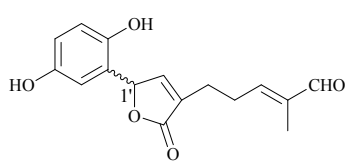


1237 (+) 3'R

1238 (-) 3'S

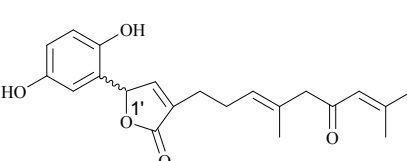


1243



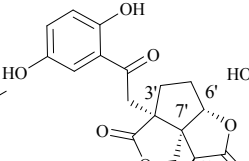
1239 (+) 1'R

1240 (-) 1'S

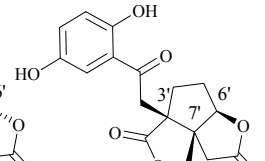


1241 (+) 1'R

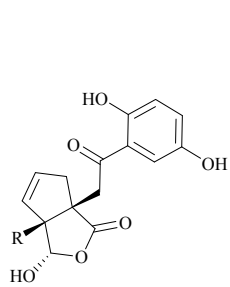
1242 (-) 1'S



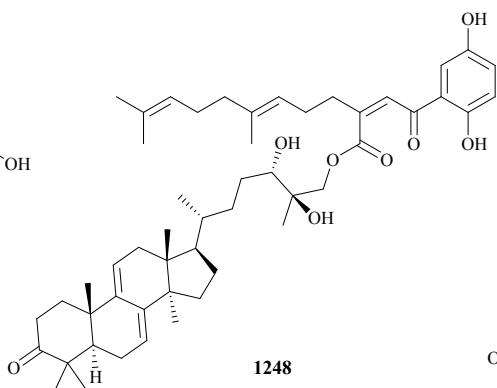
1244 (+)



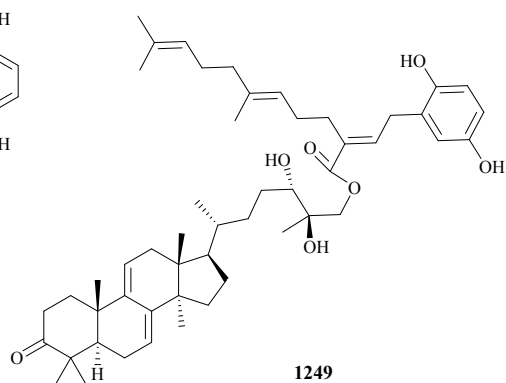
1245 (-)



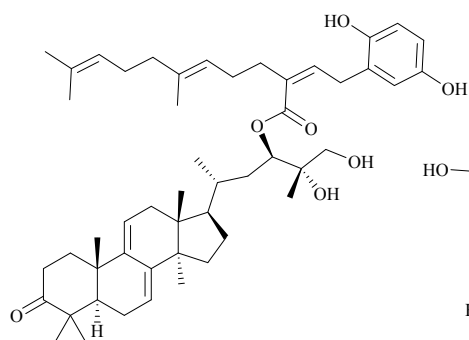
1246 R=CH₃
1247 R=H



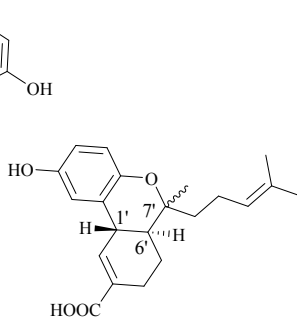
1248



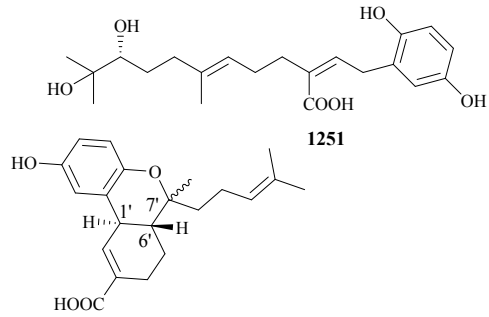
1249



1250

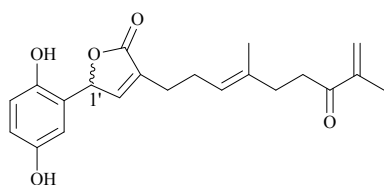


1252 (+) 1'S,6'S,7'R
1254 (+) 1'S,6'S,7'S

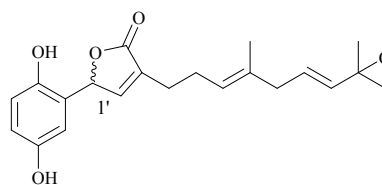


1253 (-) 1'R,6'R,7'S
1255 (-) 1'R,6'R,7'R

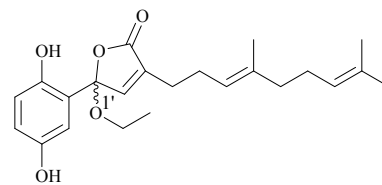
1251



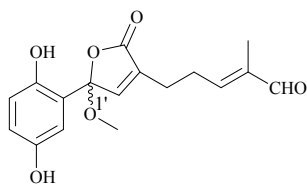
1256 (+) 1'R
1257 (-) 1'S



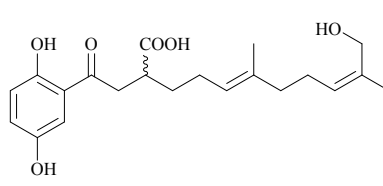
1258 (+) 1'R
1259 (-) 1'S



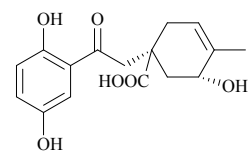
1260 (+) 1'R
1261 (-) 1'S



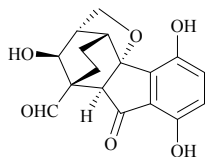
1264 (+) 1'R
1265 (-) 1'S



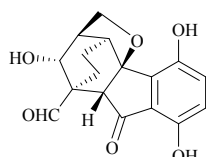
1262 (+) 3'R **1263** (-) 3'S



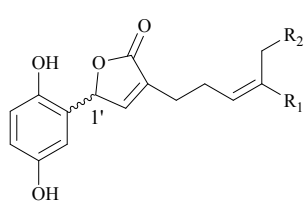
1266



1267 (+)

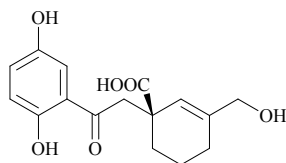


1268 (-)

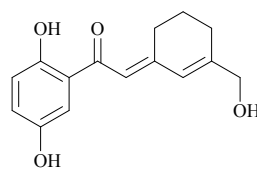


$R_1=\text{COOH}, R_2=\text{H}$ **1269** (+) 1'*R*, **1270** (-) 1'*S*

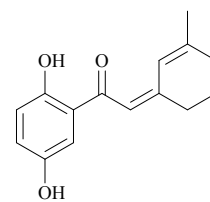
$R_1=\text{CHO}, R_2=\text{CH}_3$ **1271** ($\pm$)



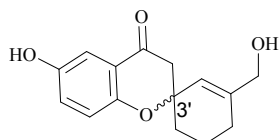
1272



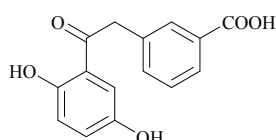
1273



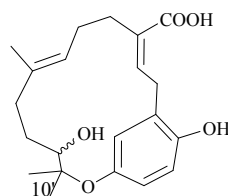
1274



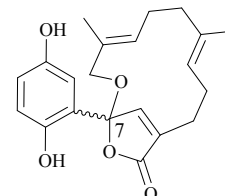
1275 (+) 3'*R* **1276** (-) 3'*S*



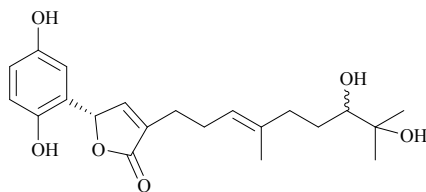
1277



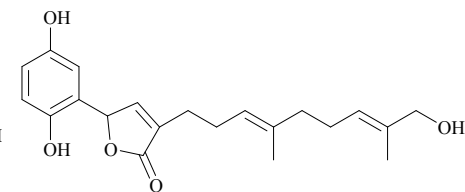
1278 (+) 10'*S* **1279** (-) 10'*R*



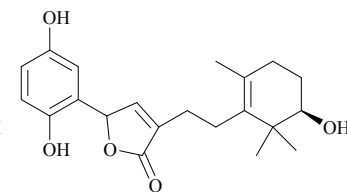
1281 7*R* **1282** 7*S*



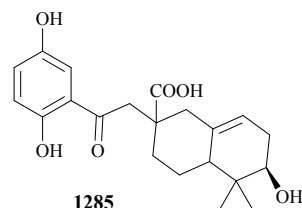
1280



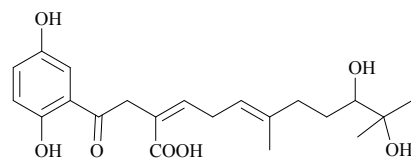
1283



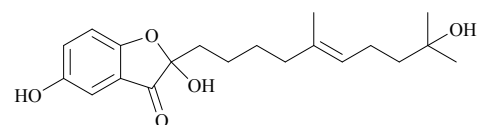
1284



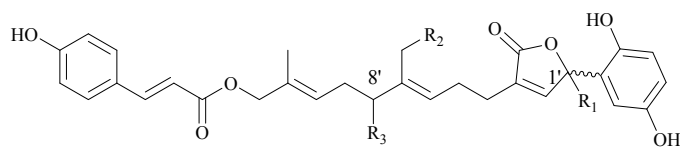
1285



1286



1287



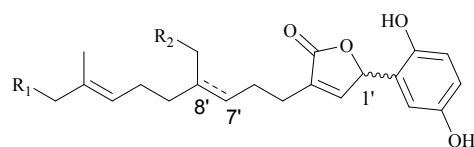
$R_1=R_3=\text{H}, R_2=\text{OH}$ **1288** 1'*R*, **1289** 1'*S*

$R_1=\text{OCH}_3, R_2=\text{OH}, R_3=\text{H}$ **1290** 1'*R*, **1291** 1'*S*

$R_1=\text{OCH}_3, R_2=R_3=\text{H}$ **1298** ($\pm$)

$R_1=R_2=R_3=\text{H}$ **1311** (+) 1'*R* **1312** (-) 1'*S*

$R_1=\text{H}, R_2=R_3=\text{OH}$ **1313** 1'*S* **1314** 1'*R*



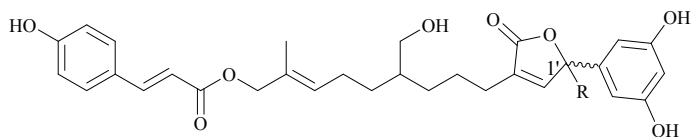
$R_1=\text{OH}, R_2=\text{H}, \Delta^7$ **1299** (+) 1'*R* **1300** (-) 1'*S*

$R_1=\text{OH}, R_2=\text{OAc}, \Delta^7$ **1301** (+) 1'*R* **1302** (-) 1'*S*

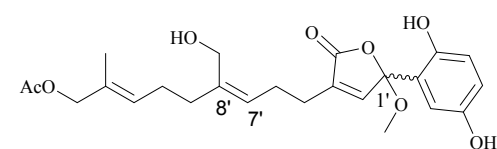
$R_1=\text{OH}, R_2=\text{OAc}$ **1303** (+) 1'*R* **1304** (-) 1'*S*

$R_1=\text{OAc}, R_2=\text{OH}, \Delta^7$ **1305** (+) 1'*R* **1306** (-) 1'*S*

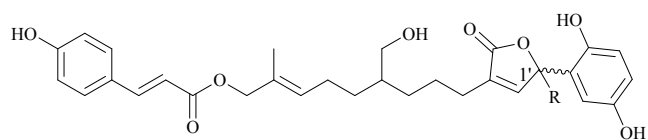
$R_1=\text{OAc}, R_2=\text{OH}$ **1307** (+) 1'*R* **1308** (-) 1'*S*



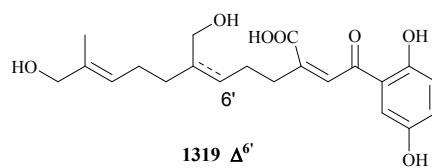
$R=\text{OCH}_3$ **1292** 1'*R*, **1293** 1'*S*



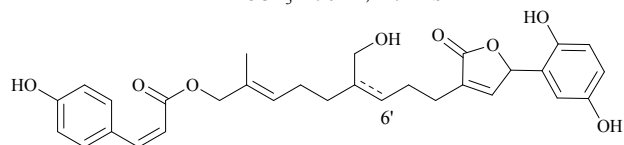
1309 (+) **1310** (-)



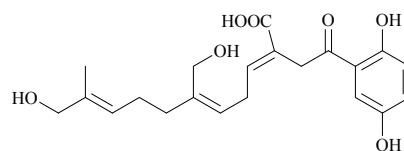
R=H **1294** 1'*R*, **1295** 1'*S*
R=OCH₃ **1296** 1'*R*, **1297** 1'*S*



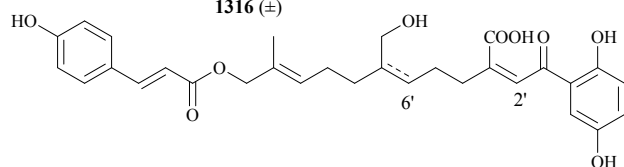
1319 Δ^{6'}
1321 (±)



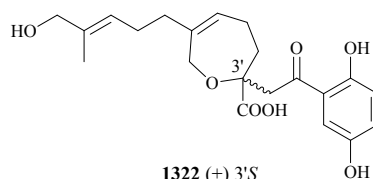
1315 (±) Δ^{6'}
1316 (±)



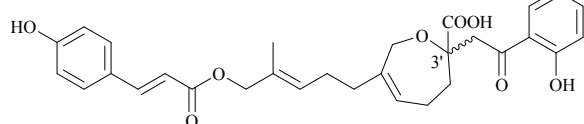
1320



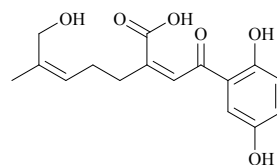
1317 Δ^{6'}
1318 (±) racemates



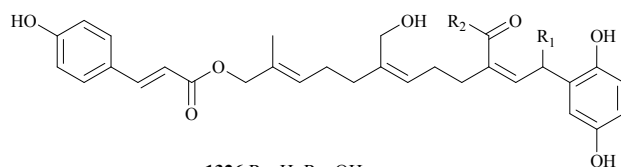
1322 (+) 3'*S*
1323 (-) 3'*R*



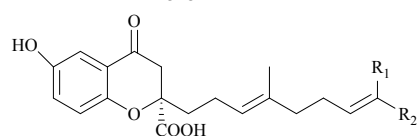
1324 (+) 3'*S*
1325 (-) 3'*R*



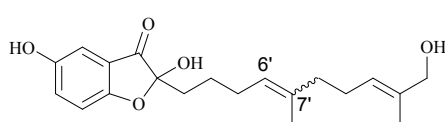
1328



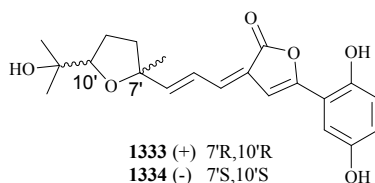
1326 R₁=H, R₂=OH
1327 R₁=ketone, R₂=OCH₃



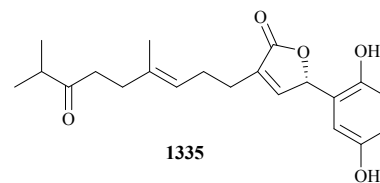
1329 R₁=CH₃, R₂=CH₂OH
1330 R₁=CH₂OH, R₂=CH₃



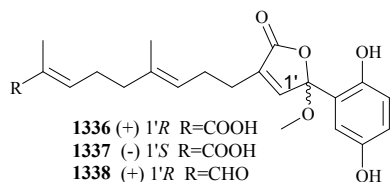
1331 *E* (Δ⁶⁽⁷⁾)
1332 *Z* (Δ⁶⁽⁷⁾)



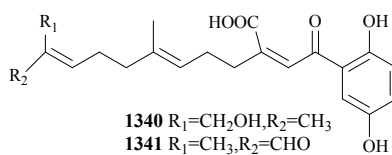
1333 (+) 7'*R*, 10'*R*
1334 (-) 7'*S*, 10'*S*



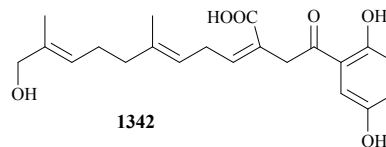
1335



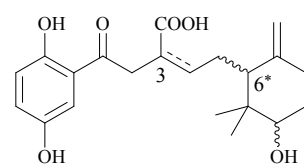
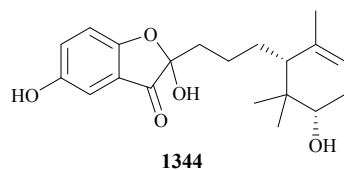
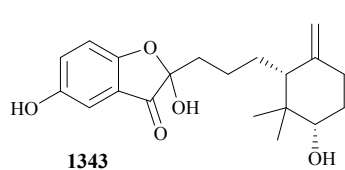
1336 (+) 1'*R* R=COOH
1337 (-) 1'*S* R=COOH
1338 (+) 1'*R* R=CHO
1339 (-) 1'*S* R=CHO



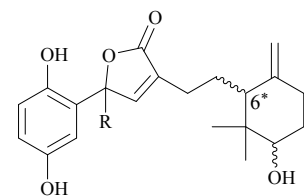
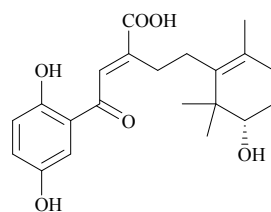
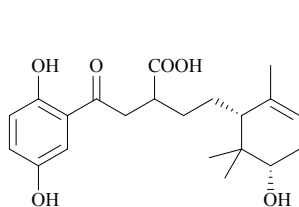
1340 R₁=CH₂OH, R₂=CH₃
1341 R₁=CH₃, R₂=CHO



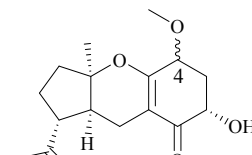
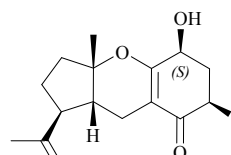
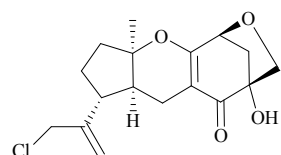
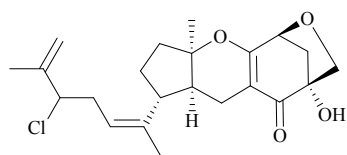
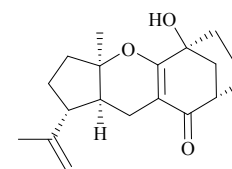
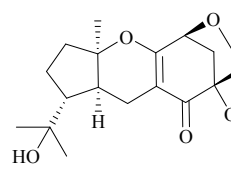
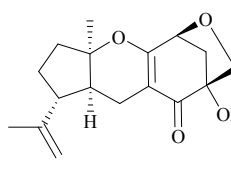
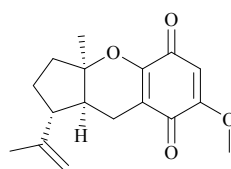
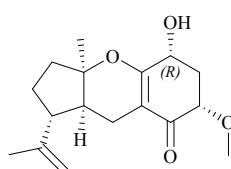
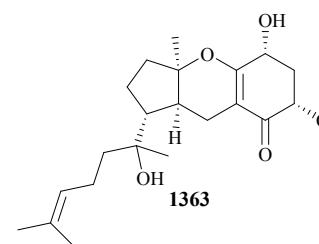
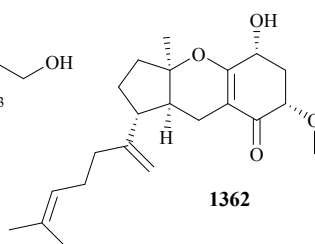
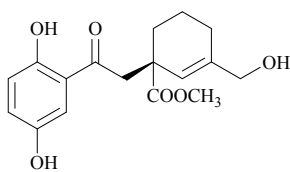
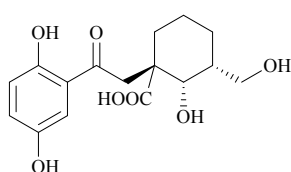
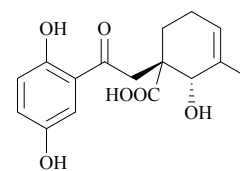
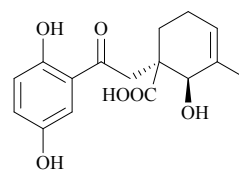
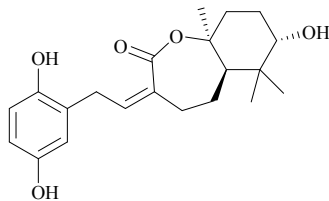
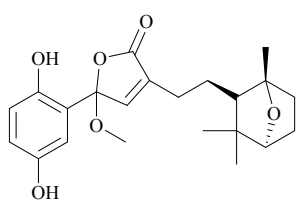
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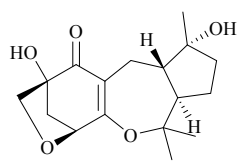


1347 (-) Δ^3 (6R)

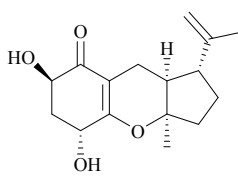


R=H 1350 (+), 1351 (-) / 1352 (+), 1353 (-)
R=OH 1354 (+) 1355 (-)

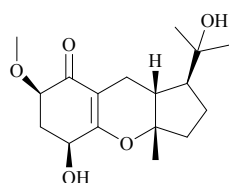




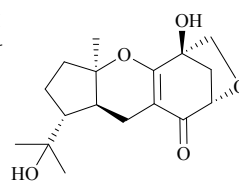
1374



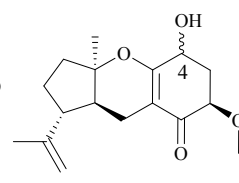
1375



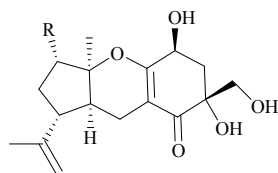
1376 (-)



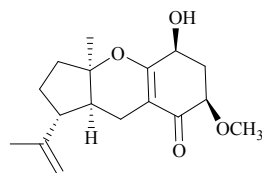
1377



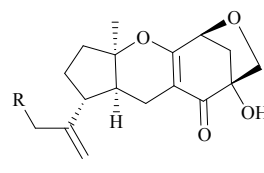
1378 4S 1379 4R



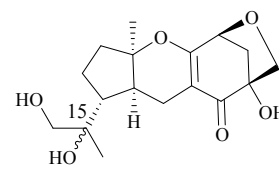
1380R=H 1381 R=OH



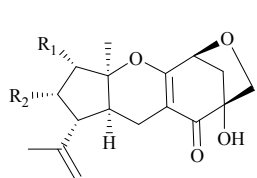
1382



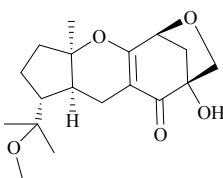
1383 R=OAc 1388 R=OH



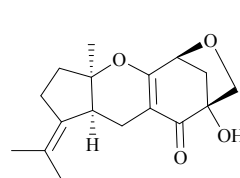
1384 15S 1385 15R



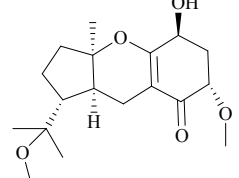
1386 R₁=H, R₂=OH
1387 R₁=OH, R₂=H



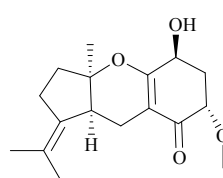
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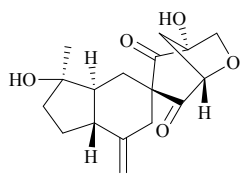
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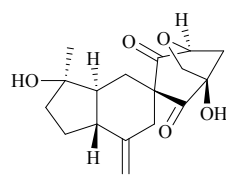
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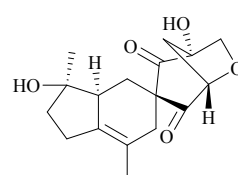
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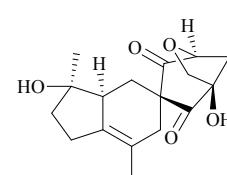
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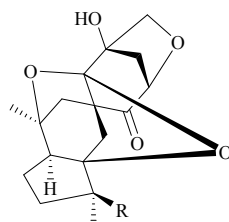
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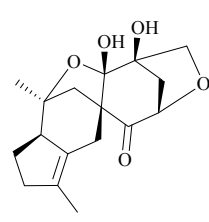
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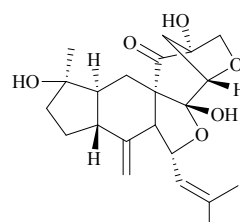
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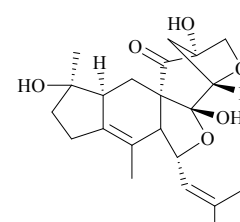
1397 R=H
1398 R=OH



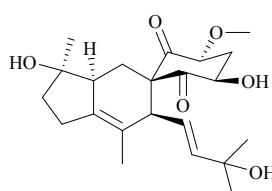
1399



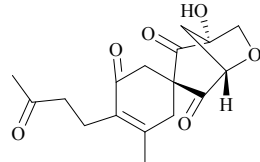
1400



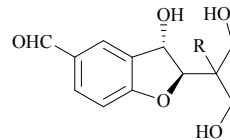
1401



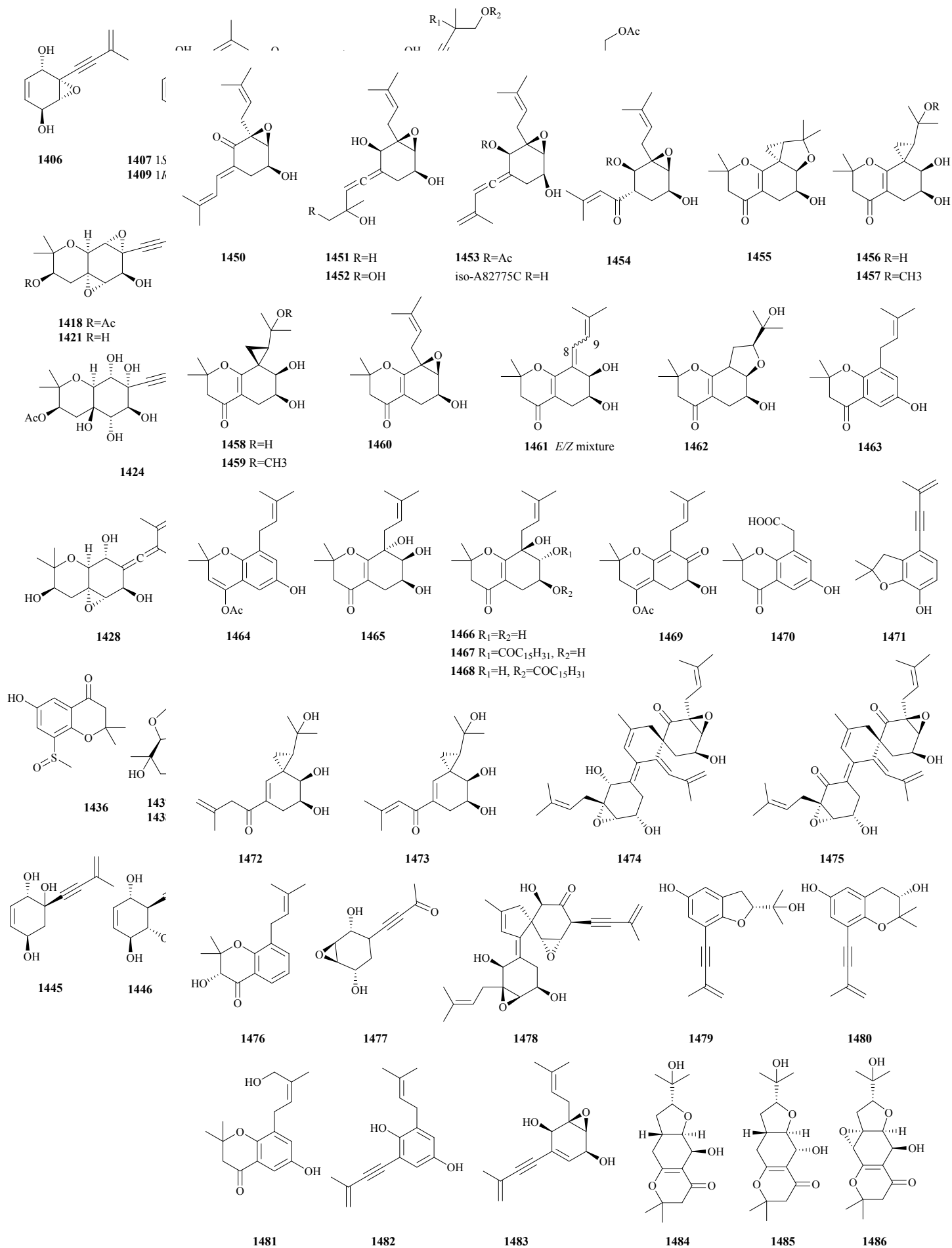
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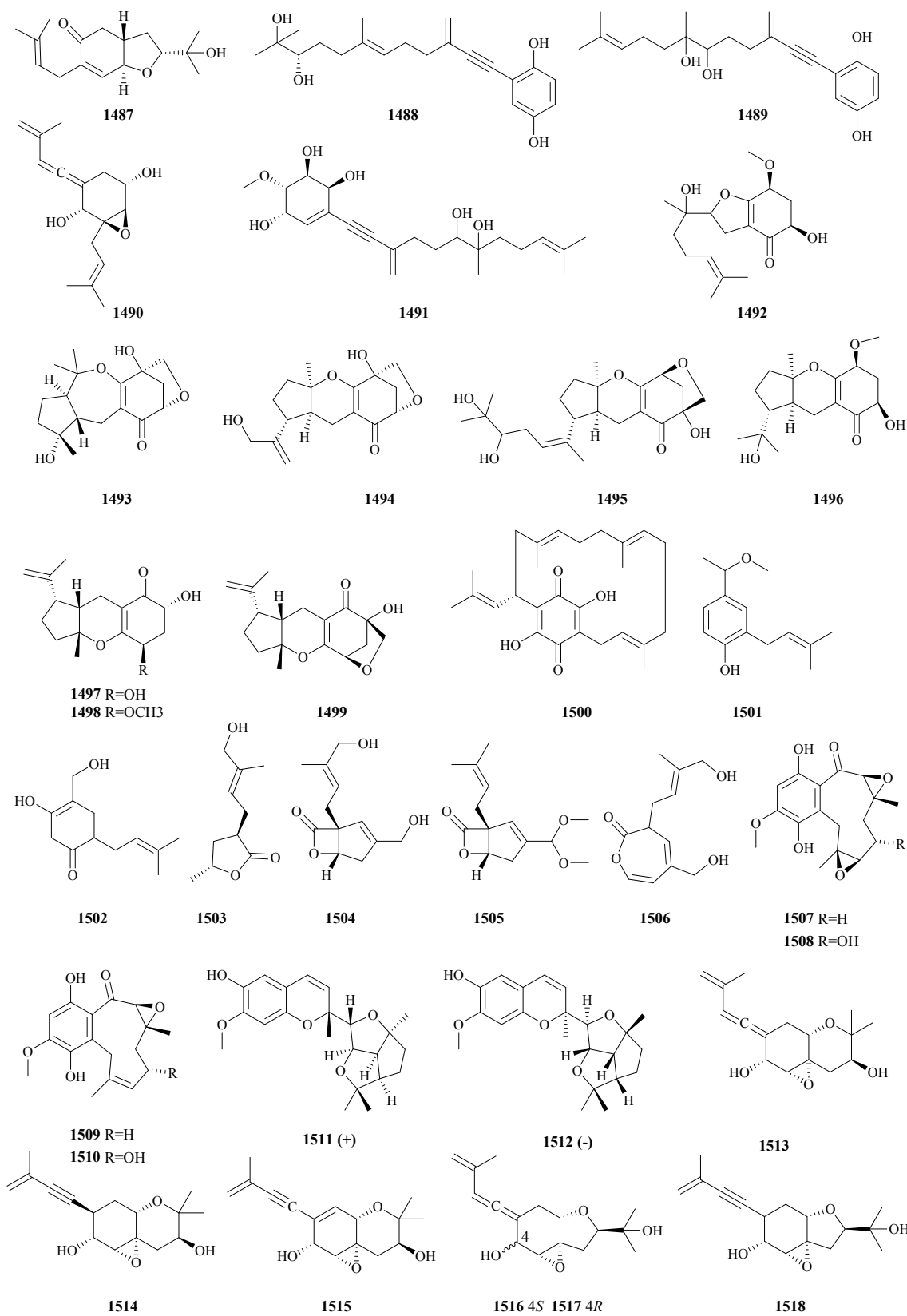


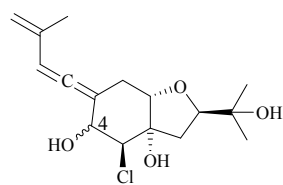
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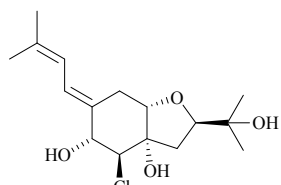
1404 R=H
1405 R=OH



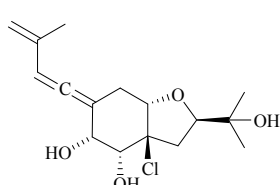




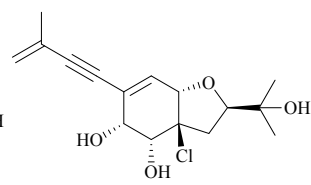
1519 4*R* 1520 4*S*



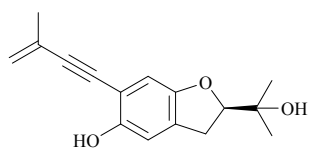
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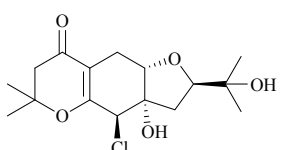
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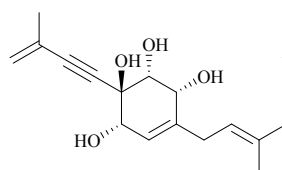
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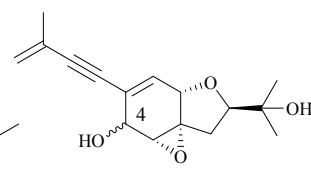
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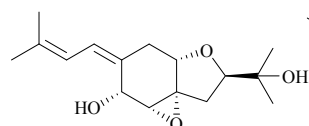
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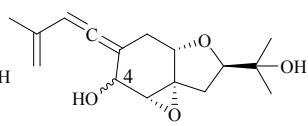
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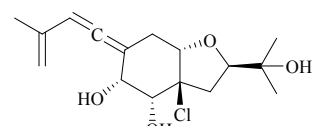
1527 4*R* 1528 4*S*



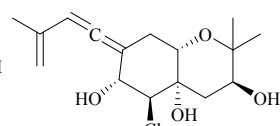
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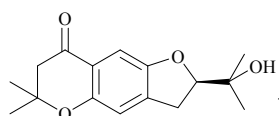
1530 4*R* 1531 4*S*



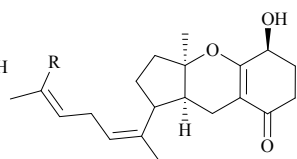
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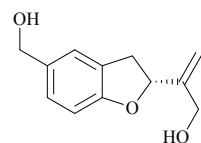
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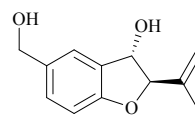
1534



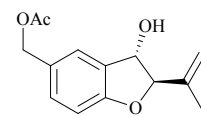
1535 R=CH₃
1536 R=CH₂OH
1537 R=COOH



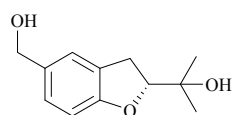
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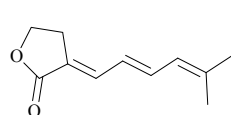
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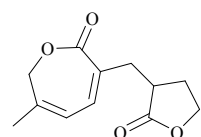
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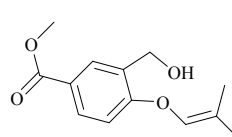
1541



1542

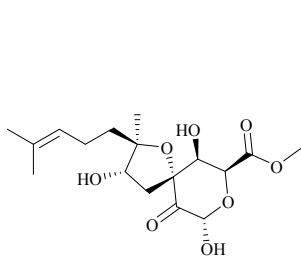


1543

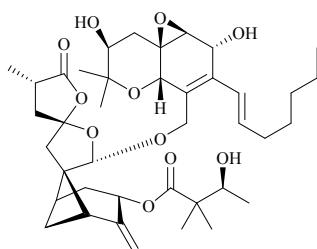


1544

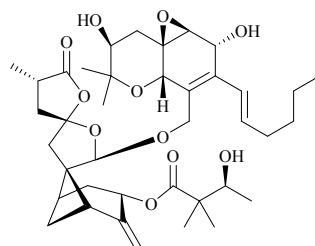
6. Miscellaneous meroterpenoids (1545-1585)



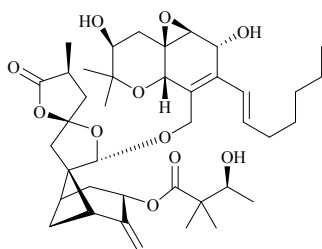
1545



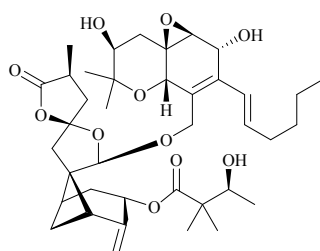
1546



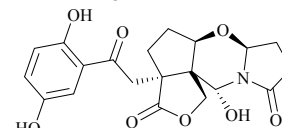
1547



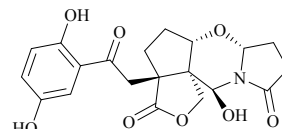
1548



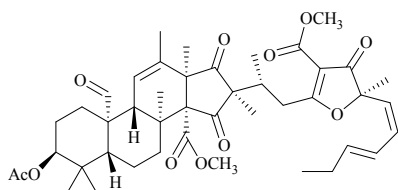
1549



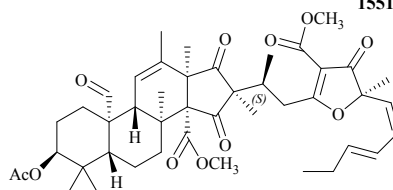
1550 (+)



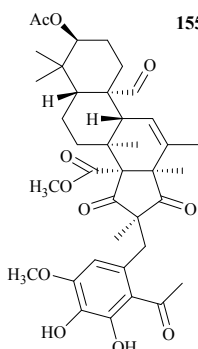
1551 (-)



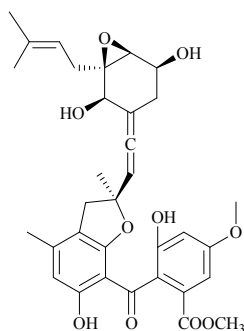
1552



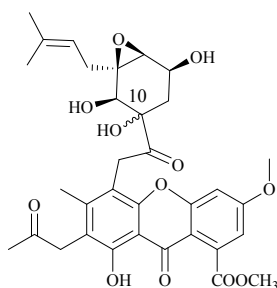
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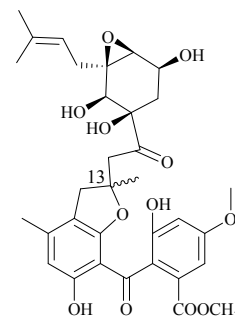
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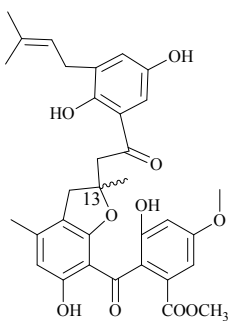
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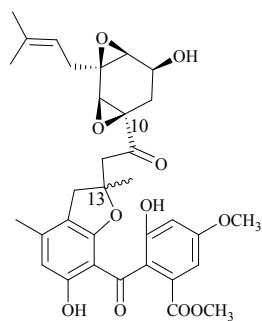
1556 10R 1557 10S



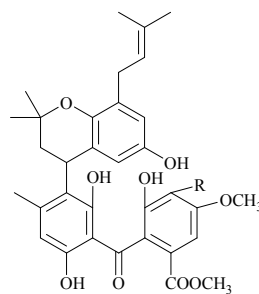
1558 13R/S mixture



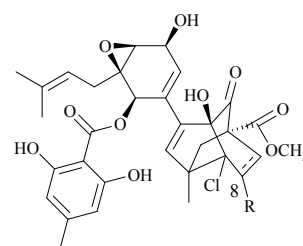
1559 1 13R/S (Racemate)



1560 13R 1561 13S



1562 R=H 1563 R=Cl



1564 R=OCH₃, Δ⁸
1565 R=ketone

