Supplemental Materials

for

Terpenoid compositions of resins from *Callitris* species (Cupressaceae)

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Contents

Figure S1. Mass spectra of derivatives and related or isomeric terpenoids:

(a) Callitrisin (I),

(b) columellarin (II),

(c) dihydrocolumellarin (III),

(d) dehydroabietane (IV, standard),

(e) callitrisol (V),

(f) ferruginol (VI, standard),

(g) sandaracopimara-8(14),15-dien-3β-ol (VII),

(h) callitrisic acid (X, standard),

(i) dehydroabietic acid (XI, standard),

(j) veadeiroic acid (standard),

(k) 19-hydroxyferruginol (XVIII, standard),

(l) hinokiol (XIX),

(m) 3β,18-dihydroxypimara-8(14),15-diene (XX),

(n) methyl abieta-6,8,11,13-tetraen-18-oate (standard),

(o) methyl cleistantha-6,8,11,13-tetraen-19-oate (standard),

(p) methyl abieta-8,11,13,15-tetraen-18-oate,

(q) methyl cleistantha-8,11,13,15-tetraen-19-oate (standard),

(r) methyl dehydroabietate (XI, standard),

(s) methyl veadeiroate (standard),

(t) methyl 5β-callitrisate (standard),

(u) methyl 5β-dehydroabietate (standard),

(v) methyl 10β(H)-9,10-*seco*-callitrisate,

(w) methyl 10α(H)-9,10-*seco*-callitrisate,

(x) methyl 10β(H)-9,10-*seco*-dehydroabietate,

(y) methyl 10α(H)-9,10-*seco*-dehydroabietate,

(z) methyl 4,5,9,10-bis-*seco*-dehydroabietate,

(aa) callitrisic acid-TMS (X, standard),

(bb) dehydroabietic acid-TMS (XI, standard),

(cc) veadeiroic acid-TMS (standard),

(dd) 4-*epi*-pimaric acid-TMS (XII),

(ee) *iso*-communic acid-TMS (XIII),

(ff) 12Z-communic acid-TMS (XIV),

(gg) 12E-communic acid-TMS (XV),

(hh) sandaracopimaric acid-TMS (XVI, standard),

(ii) ozic acid-TMS (XVIII),

(jj) 10β(H)-9,10-*seco*-dehydroabietic acid-TMS,

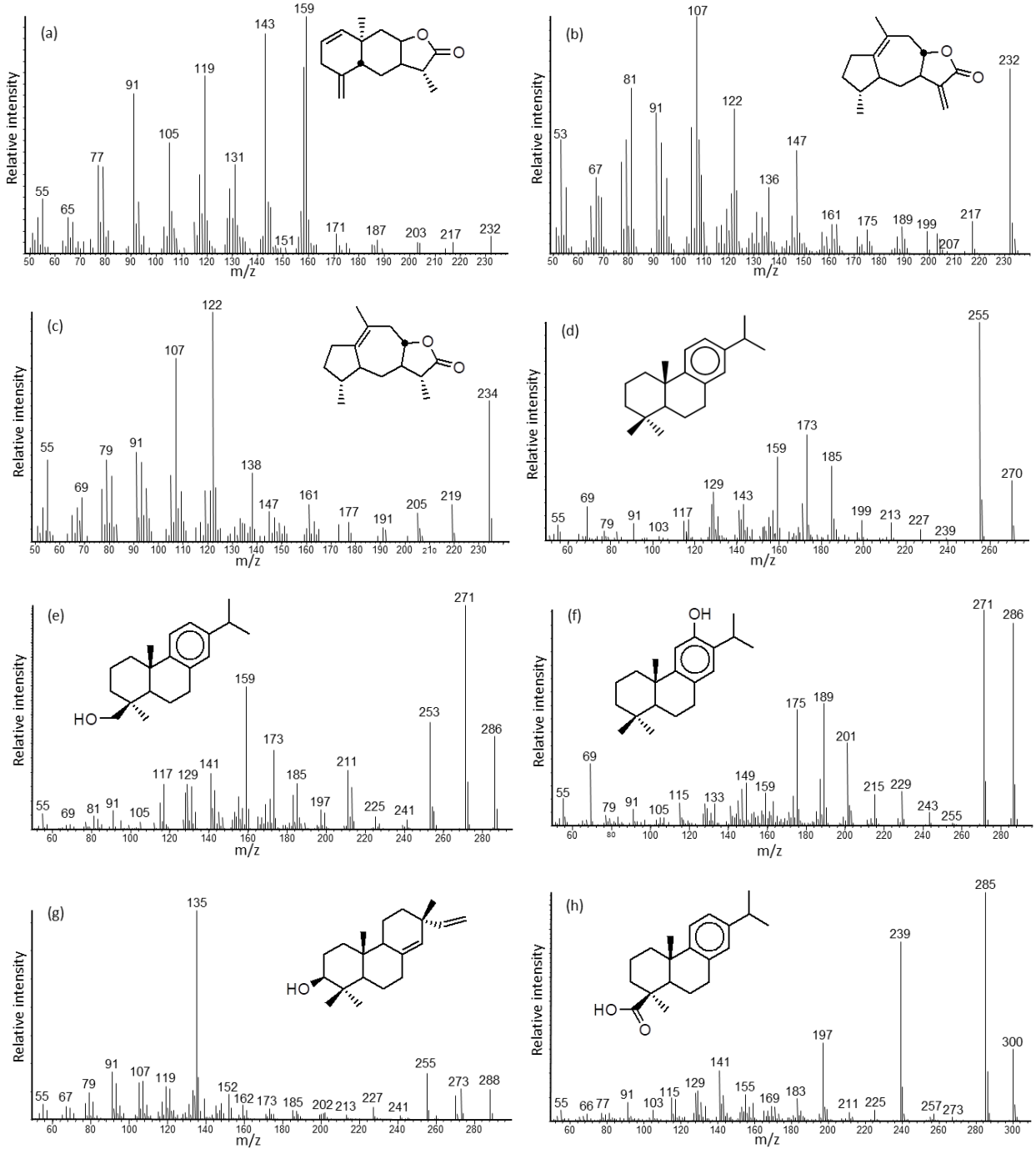
(kk) 4,5,9,10-bis-*seco*-dehydroabietic acid-TMS,

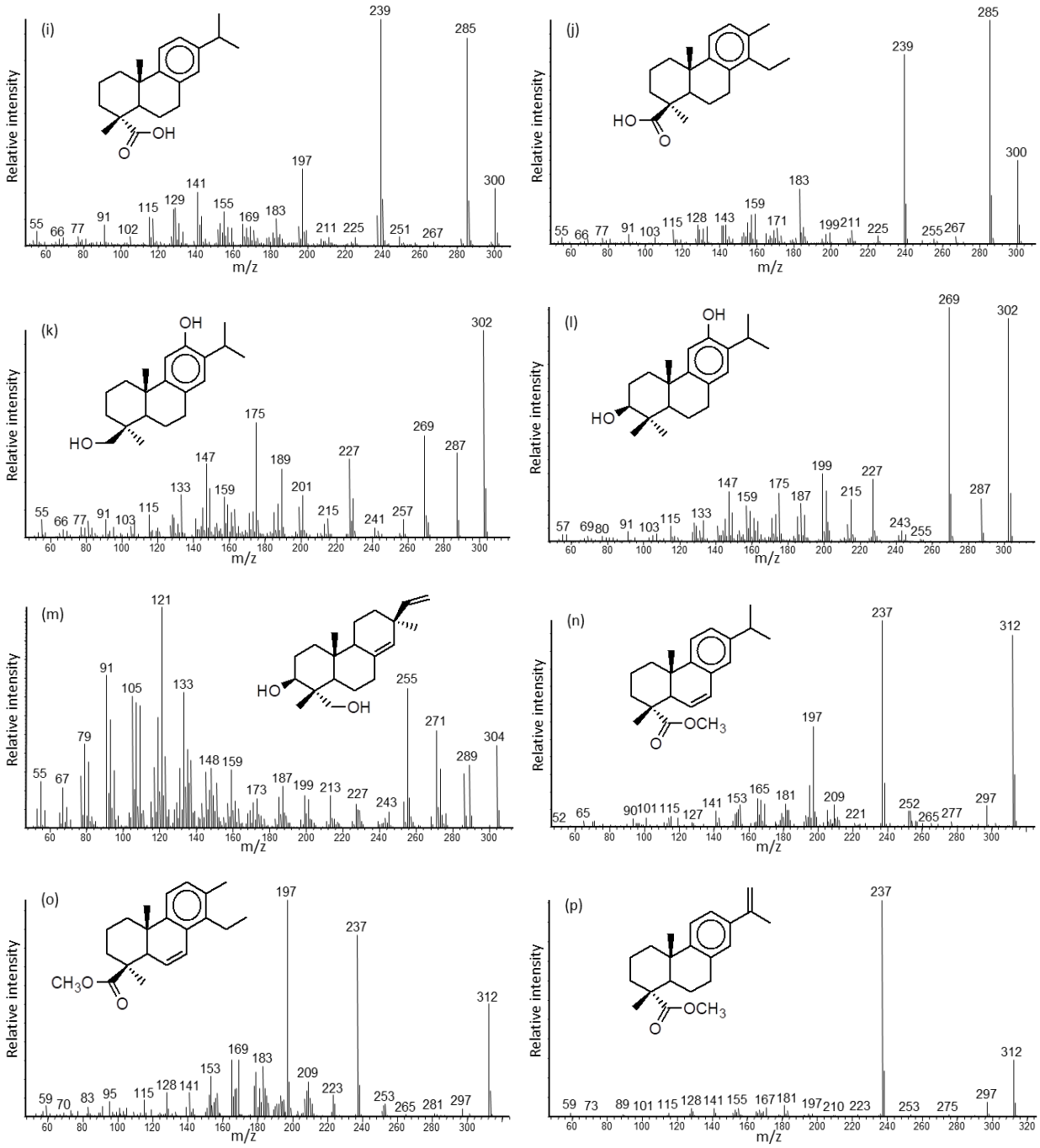
(ll) 10α(H)-9,10-*seco*-dehydroabietic acid-TMS,

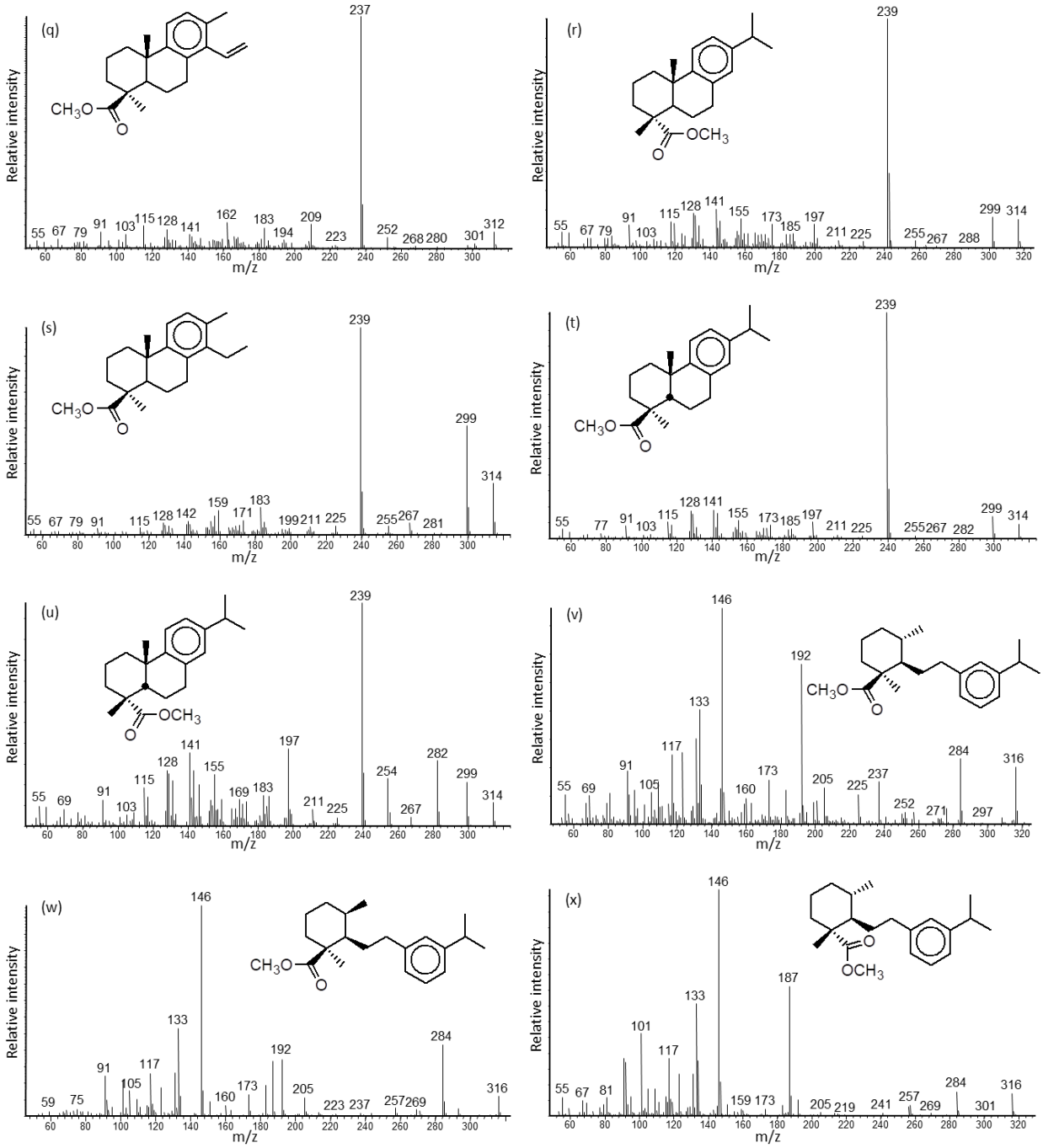
(mm) 7-oxodehydroabietic acid-TMS (standard),

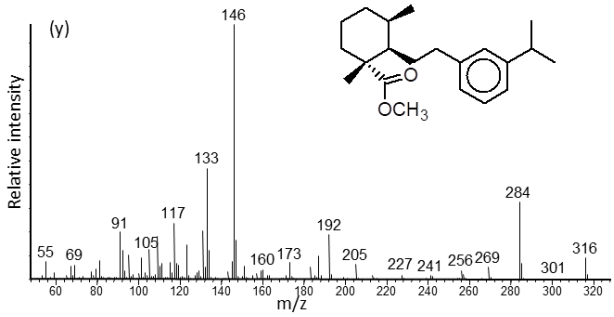
(nn) lambertianic acid-TMS (XXII, standard).

Note - structures shown on mass spectra and listed above are a reference standard, or are tentative based on a literature citation and interpretation of the fragmentation pattern with consideration of the relative GC retention index.


(a)
(b)
(c)
(d)
(e)
(f)
(g)
(h)
Relative intensity
m/z
HO
OH
O


(i)
(j)
(k)
(l)
(m)
(n)
(o)
(p)
Relative intensity
m/z
OH
HO
O
CH3O
OCH3


(q)
(r)
(s)
(t)
(u)
(v)
(w)
(x)
Relative intensity
m/z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Relative intensity
m/z
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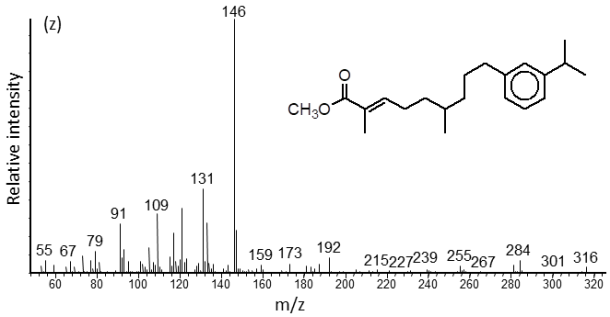
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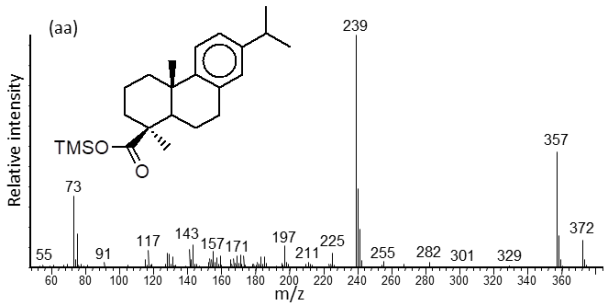
Relative intensity
m/z
CH3O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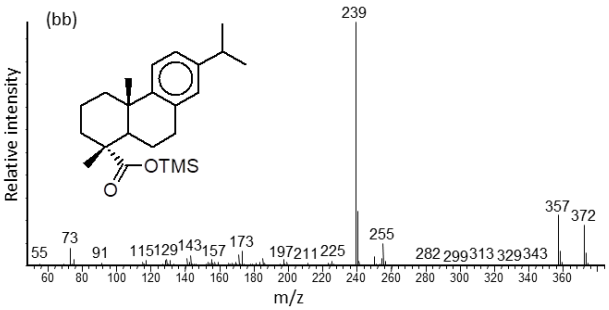
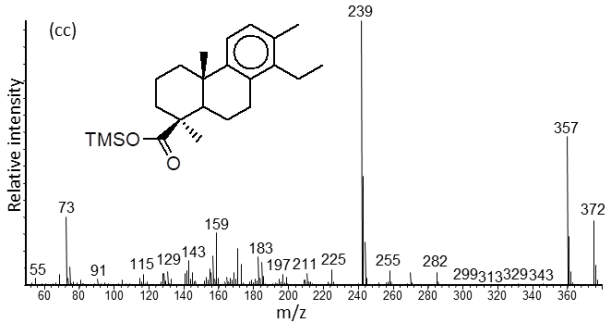
Relative intensity
m/z
TMSO


(bb)
Relative intensity
m/z
OTMS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Relative intensity
m/z
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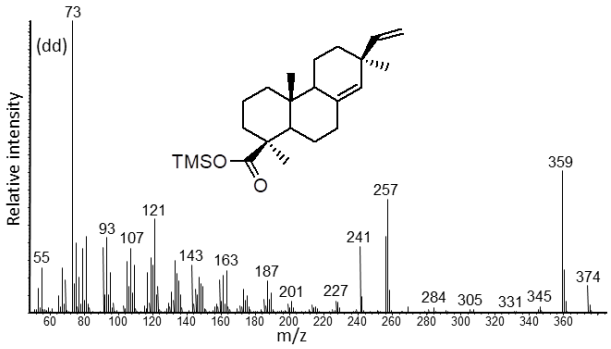
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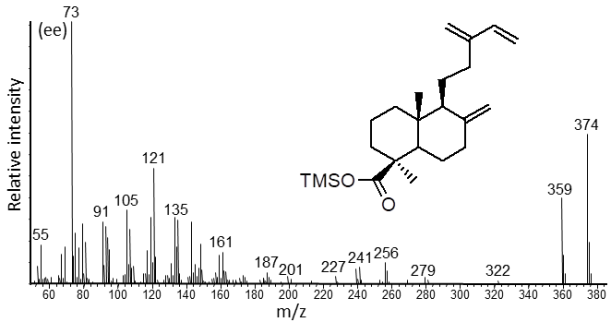
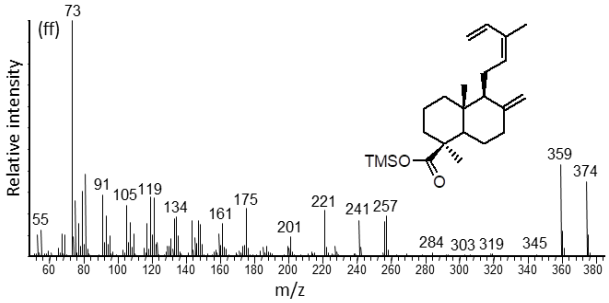
Relative intensity
m/z
TMSO


(ee)
Relative intensity
m/z
TMSO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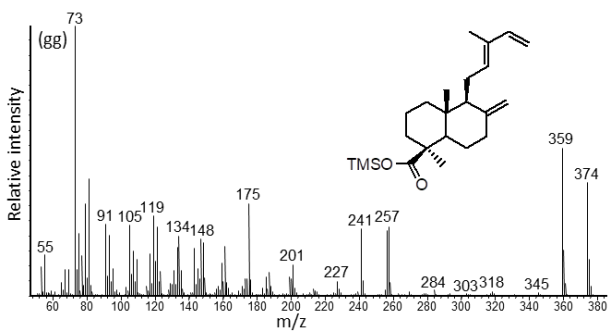
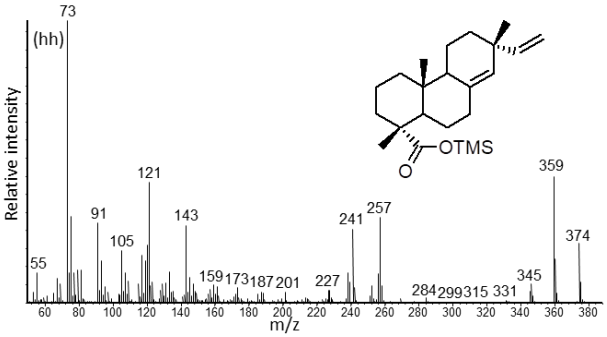
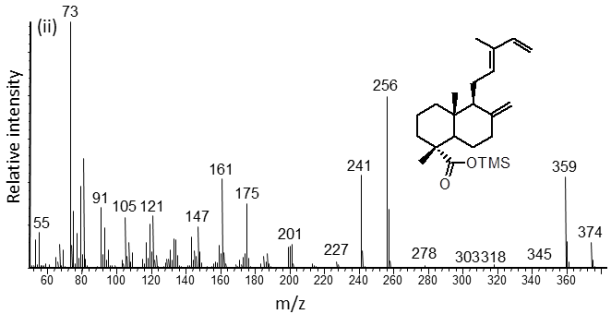
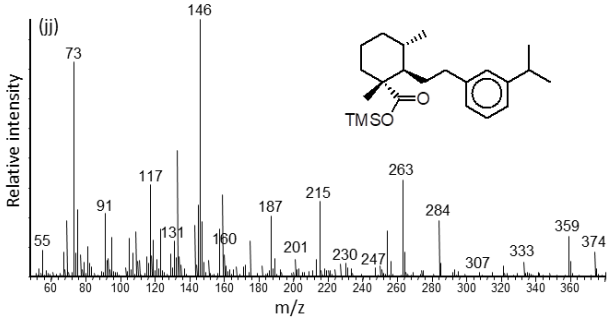
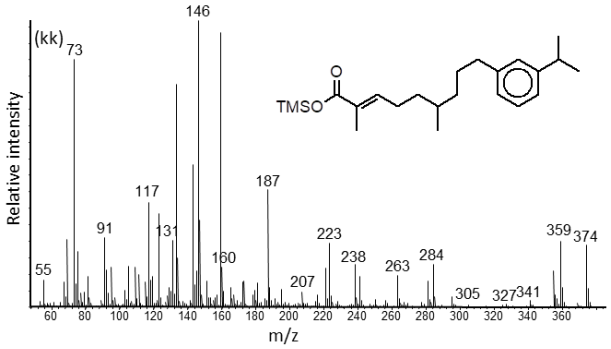
Relative intensity
m/z
TMSO


(gg)
Relative intensity
m/z
TMSO
O


(hh)
Relative intensity
m/z
OTMS
O


(ii)
Relative intensity
m/z
OTMS
O


(jj)
Relative intensity
m/z
TMSO
O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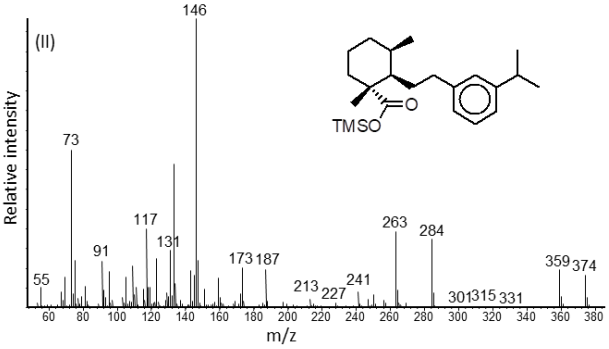
Relative intensity
m/z
TMSO
O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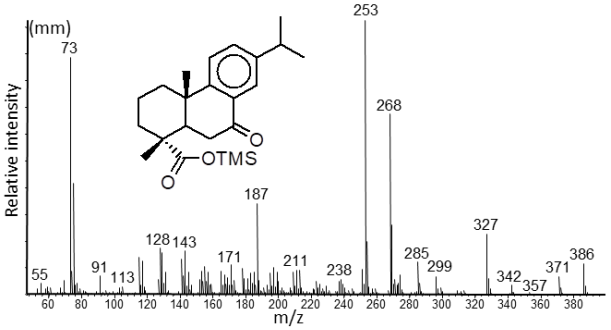
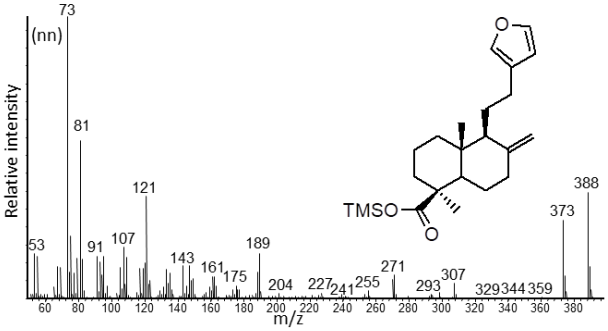
Relative intensity
m/z
TMSO
O


(mm)
Relative intensity
m/z
OTMS
O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Relative intensity
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TMSO
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