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# Dictionary of Natural Products

VOLUME FIVE

R-Z



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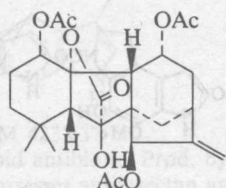
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W.C. Chan, S. Newlands, C. Williams, J. Wilson

# R

## Rabdoepigibberellolide

[81398-21-2]



$C_{26}H_{34}O_9$  M 490.549

Constit. of *Rabdosia shikokiana*. Cryst. Mp 255.5-256.5°.

$[\alpha]_D^{25} - 89^\circ$  (c, 0.28 in  $CHCl_3$ ).

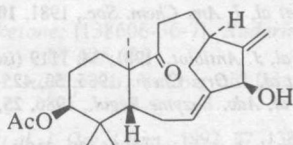
Ochi, M. et al, *J. Chem. Soc., Chem. Commun.*, 1982, 810.

## Rabdohakusin

R-00002

ent-8,9-Seco-3 $\alpha$ -acetoxy-15 $\alpha$ -hydroxy-7,16-kauradien-9-one

[93740-31-9]



$C_{22}H_{32}O_4$  M 360.492

Stereochem. wrongly given in the lit. Constit. of *Rabdosia umbrosa* var. *hakusanensis*. Cryst. Mp 97-98°.

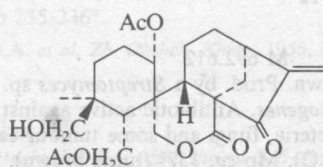
$[\alpha]_D^{25} + 78.7^\circ$  (c, 0.127 in  $CHCl_3$ ).

Kubo, I. et al, *Chem. Lett.*, 1984, 1613.

## Rabdokaurin B

R-00003

[142465-71-2]



$C_{24}H_{32}O_8$  M 448.512

Constit. of *Rabdosia longituba*. Amorph. powder.  $[\alpha]_D^{26}$

$+ 57.6^\circ$  (c, 0.81 in MeOH).

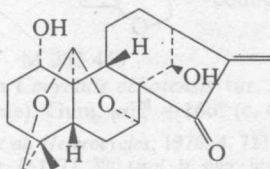
Takeda, Y. et al, *Phytochemistry*, 1992, 31, 1687 (isol, pmr, cmr)

## Rabdoserrin A

R-00004

ent-7,20:18,20-Diepoxy-1 $\beta$ ,14 $\alpha$ -dihydroxy-16-kauren-15-one

[96685-01-7]



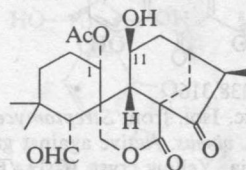
$C_{20}H_{26}O_5$  M 346.422

Isol. from *Rabdosia serra*. Shows antitumour props.

Jin, R. et al, *CA*, 1985, 103, 175377a.

## Rabdosichuanin A

R-00005



$C_{22}H_{30}O_7$  M 406.475

Constit. of *Rabdosia setschwanensis*. Cryst. Mp 225-227°.

$[\alpha]_D^{25} + 107.3^\circ$  (c, 0.55 in MeOH).

1-Deacetoxy, 11-epimer: *Rabdosichuanin B*

$C_{20}H_{28}O_5$  M 348.438

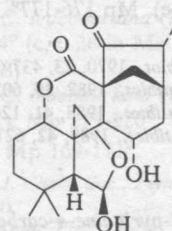
Constit. of *R. setschwanensis*. Cryst. Mp 241-243°.

$[\alpha]_D^{24} - 58.16^\circ$  (c, 0.576 in MeOH).

Hao, H. et al, *Phytochemistry*, 1990, 29, 2591 (isol, pmr, cmr)

## Rabdosichuanin C

R-00006



$C_{20}H_{28}O_6$  M 364.438

Constit. of *Rabdosia setschwanensis*. Cryst. Mp 231-233°.

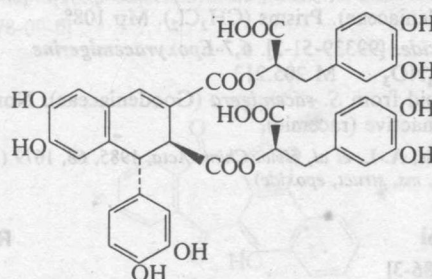
$[\alpha]_D^{25} - 120.94^\circ$  (c, 0.55 in MeOH).

Hao, H. et al, *Phytochemistry*, 1990, 29, 2591 (isol, pmr, cmr)

## Rabdoisin

R-00007

[119152-54-4]



$C_{36}H_{30}O_{16}$  M 718.623

Constit. of *Rabdosia japonica*. Used for treatment of gastrointestinal disorders. Tanning agent. Light brown amorph. powder.  $[\alpha]_D^{25} - 78^\circ$  (c, 3.5 in MeOH).

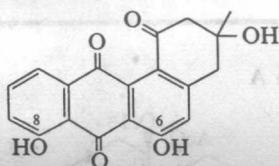
Agata, I. et al, *Chem. Pharm. Bull.*, 1988, 36, 3223 (isol, struct)

Agata, I. et al, *Phytochemistry*, 1989, 28, 2447 (isol, pmr, cmr)

**Rabelomycin**

R-00008

3,4-Dihydro-3,6,8-trihydroxy-3-methylbenz[a]anthracene-1,7,12(2H)trione, 9CI  
[28399-50-0]



$C_{19}H_{14}O_6$  M 338.316

Quinone antibiotic. Isol. from *Streptomyces olivaceus* and *S. matensis* ssp. *vineus*. Active against gram-positive and negative bacteria. Yellow cryst. ( $C_6H_6$ /EtOH). Mp 193° dec.  $[\alpha]_D^{25} - 102^\circ$  (c, 1 in  $CHCl_3$ ).

8-Me ether: [117620-88-9]. 8-O-Methylrabelomycin

$C_{20}H_{16}O_6$  M 352.343

From *S. tsusimaensis*. Active against gram-positive bacteria. Yellow needles. Mp 191-193°.  $[\alpha]_D^{25} - 118^\circ$  (c, 0.5 in MeOH).

6-Deoxy, 8-Me ether: [117620-87-8]. 6-Deoxy-8-O-methylrabelomycin. MM 47755. Antibiotic MM 47755

$C_{20}H_{16}O_5$  M 336.343

From *S. tsusimaensis* and another *S. sp.* Weakly active against gram-positive bacteria. Yellow prisms ( $CHCl_3$ /cyclohexane). Mp 176-177°.  $[\alpha]_D^{20} - 136^\circ$  (c, 0.04 in MeOH).

Liu, W.-C. et al, *J. Antibiot.*, 1970, **23**, 437 (isol)

Imamura, N. et al, *J. Antibiot.*, 1982, **35**, 602 (isol)

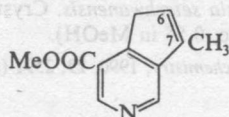
Shigihara, Y. et al, *J. Antibiot.*, 1988, **41**, 1260 (deriv)

Gilpin, M.L. et al, *J. Antibiot.*, 1989, **42**, 627 (deriv)

**Racemigerine**

R-00009

Methyl 7-methyl-5H-2-pyridine-4-carboxylate, 9CI  
[76655-39-5]



$C_{11}H_{11}NO_2$  M 189.213

Alkaloid from the aerial parts of *Scaevola racemigera* (Goodeniaceae). Prisms ( $CH_2Cl_2$ ). Mp 108°.

6,7-Epoxy: [99339-51-2]. 6,7-Epoxyracemigerine

$C_{11}H_{11}NO_3$  M 205.213

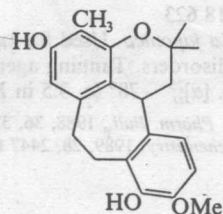
Alkaloid from *S. racemigera* (Goodeniaceae). Noncryst. Opt. inactive (racemic).

Skaltsounis, A.-L. et al, *Helv. Chim. Acta*, 1985, **68**, 1679 (isol, uv, ir, pmr, ms, struct, epoxide)

**Racemosol**

R-00010

[103805-86-3]



$C_{21}H_{24}O_4$  M 340.418

Constit. of *Bauhinia racemosa*. Deep-red prisms ( $CHCl_3$ ).

Mp 202°.  $[\alpha]_D^{25} + 64.45^\circ$  (c, 1.0 in  $CHCl_3$ ).

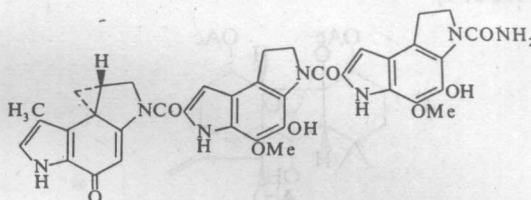
Anjaneyulu, A.S.R. et al, *Tetrahedron*, 1986, **42**, 2417 (cryst struct)

**Rachelmycin**

R-00011

CC 1065. ML 1065. NSC 298223, 106/78. Antibiotic CC 1065

[69866-21-3]



$C_{37}H_{33}N_7O_8$  M 703.710

Obt. from *Streptomyces zelensis* and *S. canulus*.

Antitumour agent, active against gram-positive and -negative bacteria and yeasts. Clustered needles or amorph. amber-coloured foam.  $\lambda_{max}$  236 (sh), 258 (sh) and 364 (end absorption) nm.

▷ Toxic. DF4936780.

Hanka, L.J. et al, *J. Antibiot.*, 1978, **31**, 1211 (isol)

Martin, D.G. et al, *J. Antibiot.*, 1980, **33**, 902 (struct)

Chidester, C.G. et al, *J. Am. Chem. Soc.*, 1981, **103**, 7629 (cryst struct)

Martin, D.G. et al, *J. Antibiot.*, 1981, **34**, 1119 (isol)

Sundberg, R.I. et al, *J. Org. Chem.*, 1985, **50**, 425 (bibl)

Wierenga, W. et al, *Adv. Enzyme Regul.*, 1986, **25**, 141 (rev, pharmacol)

Reynolds, V.L. et al, *J. Antibiot.*, 1986, **39**, 319 (rev, pharmacol)

Rawal, V.H. et al, *Heterocycles*, 1987, **25**, 701 (rev, synth)

Kelly, R.C. et al, *J. Am. Chem. Soc.*, 1987, **109**, 6837 (synth)

Watt, W. et al, *Acta Crystallogr., Sect. C*, 1988, **44**, 1675 (cryst struct)

Bolton, R.E. et al, *J. Chem. Soc., Perkin Trans. 1*, 1988, 2491 (synth, bibl)

Lewis, R.J., *Sax's Dangerous Properties of Industrial Materials*, 8th Ed., Van Nostrand-Reinhold, 1992, APT375.

**Ractinomycin A**

R-00012

[1401-11-2]

$C_{33}H_{30}N_3O_{14}$  M 692.612

Struct. unknown. Prod. by a *Streptomyces* sp. related to *S. phaeochromogenes*. Antibiotic active against gram-positive bacteria, fungi and some tumour cells. Orange needles ( $Et_2O$ ). Mp ca. 157° (turns brown), blackens at 208°. The same organism also produces Ractinomycin B.

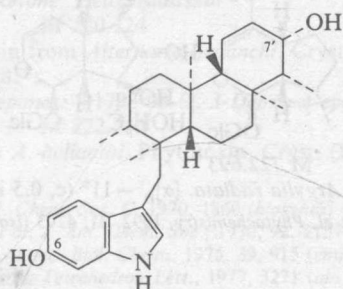
▷ VE2800000.

Utahara, R. et al, *J. Antibiot., Ser. A*, 1955, **8**, 132; 1957, **10**, 115 (isol, uv)

Wakiki, H. et al, *Antibiot. Chemother. (Washington, D.C.)*, 1958, **8**, 228 (isol)

## Radarin B

[138629-07-9]

 $C_{28}H_{41}NO_2$  M 423.637

Indole diterpenoid antibiotic. Prod. by *Aspergillus sulphureus*. Possesses antiinsectan and cytotoxic props. Yellow solid. Mp 115-118° dec.  $[\alpha]_D + 39.4^\circ$  (c, 0.003 in  $CHCl_3$ ).

6-Deoxy: [138606-37-8]. **Radarin D** $C_{28}H_{41}NO$  M 407.638

Prod. by *A. sulphureus*. Yellow oil.  $[\alpha]_D + 31.8^\circ$  (c, 0.003 in  $CHCl_3$ ).

7'-Ketone: [138606-35-6]. **Radarin A** $C_{28}H_{39}NO_2$  M 421.622

Prod. by *A. sulphureus*. Pink oil.  $[\alpha]_D + 11.1^\circ$  (c, 0.005 in  $CHCl_3$ ).

6-Deoxy, 7'-ketone: [138606-36-7]. **Radarin C** $C_{28}H_{39}NO$  M 405.622

Prod. by *A. sulphureus*. Yellow oil.  $[\alpha]_D + 6.7^\circ$  (c, 0.002 in  $CHCl_3$ ).

Laakso, J.A. et al, *J. Org. Chem.*, 1992, 57, 138 (isol, pmr, cmr, struct)

## Raddeamine

R-00014

Struct. unknown

 $C_{23}H_{37}NO_2$  M 359.551

Prob. a steroidal alkaloid. Alkaloid from *Fritillaria raddeana* (Liliaceae). Mp 271-272°. Occurs with Alvanidine, A-01152 and Alvanine, A-01153.

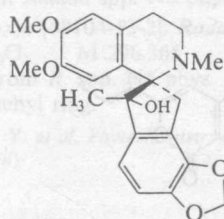
B, HCl: Mp 235-236°.

Aslanov, Kh.A. et al, *Zh. Obshch. Khim.*, 1956, 20, 579; *CA*, 50, 13971.

## Raddeanamine

R-00015

3',4',6,8-Tetrahydro-6',7'-dimethoxy-2',6-dimethylspiro[7H-indeno[4,5-d]-1,3-dioxole-7,1'(2'H)-isoquinolin]-6-ol, 9CI [59614-35-6]



Absolute configuration

 $C_{23}H_{25}NO_5$  M 383.443

Alkaloid from *Corydalis ochotensis* var. *raddeana* (Fumariaceae). Gum.  $[\alpha]_D^{20} + 166^\circ$  (c, 0.68 in MeOH).

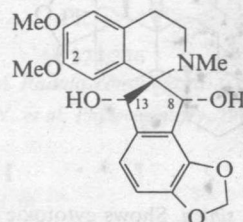
Kametani, T. et al, *Heterocycles*, 1976, 4, 723; *J. Chem. Soc., Perkin Trans. 1*, 1977, 390 (isol, ir, pmr, ms, struct)

## Raddeanine†

R-00016

3',4',6,8-Tetrahydro-6',7'-dimethoxy-2'-methylspiro[7H-indeno[4,5-d]-1,3-dioxole-7,1'(2'H)-isoquinoline]-6,8-diol, 9CI. 13-epi-Yenusomine

[59654-07-8]



(+) -form

 $C_{21}H_{23}NO_6$  M 385.416

Diastereoisomeric with Yenusomine, Y-00020.

(+) -form

Alkaloid from *Corydalis ochotensis* var. *raddeana* and *C. govaniana* (Fumariaceae). Prisms (Me<sub>2</sub>CO). Mp 208-209° (200-202°).  $[\alpha]_D + 79.4^\circ$  (c, 0.11 in MeOH).

O<sup>8</sup>-Ac: [59614-36-7]. **Raddeanidine** $C_{23}H_{25}NO_7$  M 427.453

Alkaloid from *C. ochotensis* var. *raddeana* (Fumariaceae). Gum.  $[\alpha]_D^{20} + 82.7^\circ$  (c, 0.52 in MeOH).

O<sup>2</sup>-De-Me: [64191-01-1]. **Ledeboridine. Ledebouridine** $C_{20}H_{21}NO_6$  M 371.389

Alkaloid from *C. ledebouriana* (Fumariaceae). Mp 140-141°.  $[\alpha]_D + 114^\circ$  (c, 0.28 in MeOH).

(±) -form [64234-40-8]

Alkaloid from *C. ledebouriana* (Fumariaceae). Cryst. (MeOH/ $CHCl_3$ ). Mp 219-220°, Mp 201-202° (synthetic).

O<sup>8</sup>-Ac: Synthetic. Mp 169-171°.

Kametani, T. et al, *J. Chem. Soc., Perkin Trans. 1*, 1977, 390 (isol, ir, pmr, ms, struct)

Israilov, I.A. et al, *Khim. Prir. Soedin.*, 1977, 428; *CA*, 87, 148667q (isol, ir, pmr, ms, struct)

Mukhopadhyay, S. et al, *J. Nat. Prod. (Lloydia)*, 1987, 50, 270

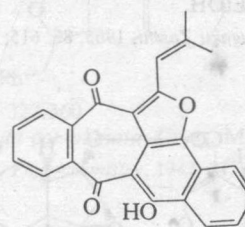
(isol, uv, ir, pmr, cmr, ms)

Hanaoka, M. et al, *Chem. Pharm. Bull.*, 1988, 36, 4248 (synth, ir, pmr, Raddeanine, Raddeanidine)

## Radermachol

R-00017

9-Hydroxy-2-(2-methyl-1-propenyl)benzo[g]benzo[5,6]cyclohepta[1,2,3-cd]benzofuran-3,8-dione, 9CI [95378-00-0]

 $C_{24}H_{16}O_4$  M 368.388Red pigment from roots of *Radermachera xylocarpa*.

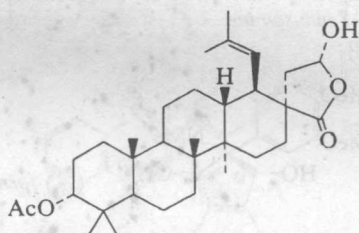
Cryst. (MeOH/hexane). Mp 217-218°.

Joshi, B.S. et al, *Tetrahedron Lett.*, 1984, 25, 5847 (cryst struct)

Jiang, Q. et al, *Tetrahedron Lett.*, 1991, 32, 5283 (synth)

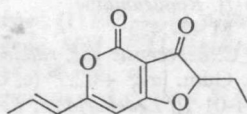
## Radermasinin

[105531-92-8]

 $C_{32}H_{50}O_5$  M 514.744Constit. of *Radermachia sinica*. Shows cytotoxic activity.Cryst. Mp 239-242°.  $[\alpha]_D^{20} -27.0^\circ$  (c, 0.58 in  $CHCl_3$ ).Rice, G.K. et al, *J. Chem. Soc., Chem. Commun.*, 1986, 1397 (cryst struct)

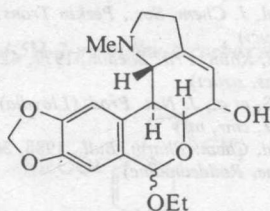
## Radianthin

[97399-68-3]

 $C_{12}H_{12}O_4$  M 220.224Phytotoxin from *Alternaria helianthi*. Low melting solid.Mp 5°.  $[\alpha]_D^{20} 0^\circ$  (c, 0.15 in  $CHCl_3$ ).Tal, B. et al, *Phytochemistry*, 1985, 24, 729.

## Radiatorine†

R-00020

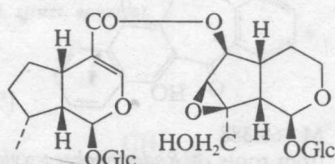
 $C_{19}H_{23}NO_5$  M 345.394Alkaloid from the rhizomes of *Lycoris radiata*(Amaryllidaceae). Cryst. ( $Me_2CO$ ). Mp 171-172.5°.

Assumed to be an artifact arising from the processing of the plant with EtOH.

Uyeo, S. et al, *Yakugaku Zasshi*, 1965, 85, 615; *CA*, 63, 11632e.

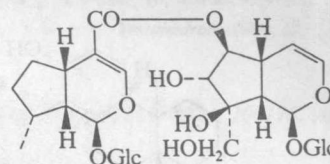
## Radiatoride E

R-00021

 $C_{31}H_{46}O_{18}$  M 706.694Constit. of *Argylia radiata*.  $[\alpha]_D -113^\circ$  (c, 1.5 in MeOH).Bianco, A. et al, *Phytochemistry*, 1992, 31, 4203 (isol, pmr, cmr)

## Radiatoride F

R-00022

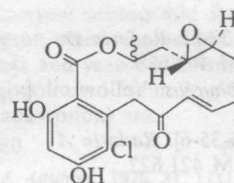
 $C_{31}H_{46}O_{19}$  M 722.693Constit. of *Argylia radiata*.  $[\alpha]_D -11^\circ$  (c, 0.5 in MeOH).Bianco, A. et al, *Phytochemistry*, 1992, 31, 4203 (isol, pmr, cmr)

## Radicicoll

R-00023

8-Chloro-1a,14,15,15a-tetrahydro-9,11-dihydroxy-14-methyl-6H-oxireno[e][2]benzoxacetyl tetradecin-6,12(7H)-dione, 9Cl. Monorden

[12772-57-5]

 $C_{18}H_{17}ClO_6$  M 364.781Macrolide antibiotic. Metab. of *Nectria radiculicola*, *Monosporium* spp., *Penicillium luteoaurantium* and *Monocillium nordinii*. Active against gram-positive and gram-negative bacteria, fungi and tumours. Potent tranquilliser of low toxicity. Prisms ( $Et_2O$ ). Mp 195°.  $[\alpha]_D^{20} +216^\circ$  (c, 1 in  $CHCl_3$ ).

▷ RR1105000.

Di-Ac: Cryst. Mp 189-190°.

Di-Me ether: [75207-16-8].

Shows antitumour and nematocidal props. Mp 186-187°.

 $[\alpha]_D -58^\circ$  (c, 1 in  $CHCl_3$ ). Other Dialkyl ethers show similar props.

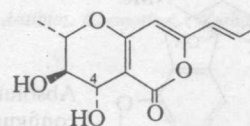
[11021-61-7]

McCapra, F. et al, *Tetrahedron Lett.*, 1964, 869 (struct, nmr, ir, ms)Mirrington, R.N. et al, *Aust. J. Chem.*, 1966, 19, 1265 (struct, ir, uv, nmr)Betina, V. et al, *Z. Allg. Mikrobiol.*, 1972, 12, 355; 1973, 13, 287 (props)Nozawa, K. et al, *J. Nat. Prod. (Lloydia)*, 1979, 42, 374.Ayer, W.A. et al, *Can. J. Microbiol.*, 1980, 26, 766 (isol, derivs)U.S. Pat., 4 228 079, (1980); *CA*, 94, 121499 (derivs)Lampilas, M. et al, *Tetrahedron Lett.*, 1992, 33, 773, 777 (synth)

## Radicinol

R-00024

[65647-66-7]

 $C_{12}H_{14}O_5$  M 238.240Metab. of *Cochliobolus lunata*. Oil.  $[\alpha]_D^{21} -175^\circ$  (c, 1 in  $CHCl_3$ ).4-Ketone: [10088-95-6]. *Radicinin*. *Stemphyllone* $C_{12}H_{12}O_5$  M 236.224Metab. of *Stemphyllium radicinum*, *Alternaria**chrysanthemi* and *Curvularia lunata*. Active againstgram-positive bacteria and fungi. Phytotoxin. Needles ( $EtOH$ ). Mp 220° dec., 236-240° dec.  $[\alpha]_D^{27} -208^\circ$  (c, 1.25 in  $EtOH$ ).

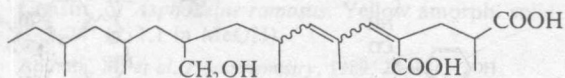
▷ UQ1360000.

4-Ketone, Ac: Cryst. Mp 197°. [ $\alpha$ ]<sub>D</sub><sup>27</sup> –267° (c, 0.7 in Py).3-Deoxy, 4-ketone: **Deoxyradicinin**C<sub>12</sub>H<sub>12</sub>O<sub>4</sub> M 220.224Phytoalexin from *Alternaria helianthi*. Cryst. (Me<sub>2</sub>CO). Mp 183-185°.

3-Deoxy, 4-epimer: [91793-98-5]. 3-Deoxy-4-epiradicinol

C<sub>12</sub>H<sub>14</sub>O<sub>4</sub> M 222.240Isol. from *A. helianthi*. Phytotoxin. Cryst. (EtOH). Dec. <100°.Grove, J.F., *J. Chem. Soc. C*, 1970, 1860 (biosynth)Tanabe, M. et al, *J. Am. Chem. Soc.*, 1970, **92**, 2157 (biosynth)Seto, H. et al, *Agric. Biol. Chem.*, 1975, **39**, 915 (cmr)Nukina, M. et al, *Tetrahedron Lett.*, 1977, 3271 (abs config)Robeson, D.J. et al, *Phytochemistry*, 1982, **21**, 1821, 2359; 1984,**23**, 767 (isol, cryst struct, abs config, derivs)Strijewski, A. et al, *Biotechnol. Lett.*, 1982, **4**, 495 (biosynth)Tal, B. et al, *J. Chem. Soc., Perkin Trans. 1*, 1988, 1283 (biosynth)**Radicionic acid****R-00025**

4-Carboxy-10-hydroxymethyl-2,6,8,12,14-pentamethyl-4,6-hexadecadienoic acid. 2-(6-Hydroxymethyl-2,4,8,10-tetramethyl-2-dodecenylidene)-4-methylpentanedioic acid [49620-14-6]

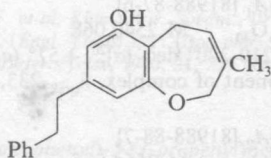
C<sub>23</sub>H<sub>40</sub>O<sub>5</sub> M 396.566Produced by an unidentified fungus. Isol. from a *Penicillium* sp. Plant growth regulator. Oil.

Di-Me ester: [42998-57-2].

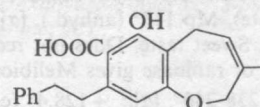
[ $\alpha$ ]<sub>D</sub><sup>24</sup> +15° (c, 0.92 in CHCl<sub>3</sub>).Sassa, T. et al, *Agric. Biol. Chem.*, 1973, **37**, 1221 (isol)Sassa, T. et al, *Tetrahedron Lett.*, 1973, 2333 (isol, struct)Japan. Pat., 75 81 863, (1975); *CA*, **83**, 159157 (props)Seto, H. et al, *Tetrahedron Lett.*, 1977, 4083 (biosynth, cmr)**Radulanin A****R-00026**

2,5-Dihydro-3-methyl-8-(2-phenylethyl)-1-benzoxepin-6-ol, 9CI

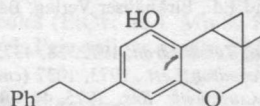
[68104-12-1]

C<sub>19</sub>H<sub>20</sub>O<sub>2</sub> M 280.366Isol. from *Radula* spp. No phys. props. reported.4'-Hydroxy: [68104-13-2]. **Radulanin C**C<sub>19</sub>H<sub>20</sub>O<sub>3</sub> M 296.365Isol. from *R. spp.* No phys. props. reported. Substd. in the phenyl ring.Asakawa, Y. et al, *Phytochemistry*, 1978, **17**, 2005, 2115; 1981, **20**, 858 (isol)**Radulanin H****R-00027**

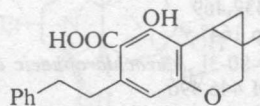
[85526-70-1]

C<sub>20</sub>H<sub>20</sub>O<sub>4</sub> M 324.376Constit. of *Radula complanata*. Cryst. Mp 122-123°.Asakawa, Y. et al, *Phytochemistry*, 1982, **21**, 2481.**Radulanin I****R-00028**

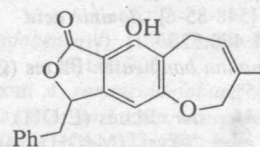
[132839-41-9]

C<sub>19</sub>H<sub>20</sub>O<sub>2</sub> M 280.366Constit. of the liverwort, *Radula javanica*. Oil. [ $\alpha$ ]<sub>D</sub> –12.8° (c, 0.56 in MeOH).Me ether: [132839-42-0]. **Radulanin J**C<sub>20</sub>H<sub>22</sub>O<sub>2</sub> M 294.393Constit. of *R. javanica*. Oil. [ $\alpha$ ]<sub>D</sub> +47.1° (c, 0.15 in MeOH).Asakawa, Y. et al, *Phytochemistry*, 1991, **30**, 325 (isol, pmr, cmr)**Radulanin K****R-00029**

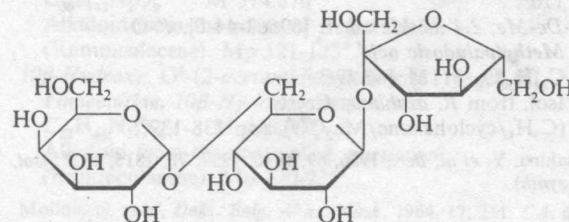
[132839-43-1]

C<sub>20</sub>H<sub>20</sub>O<sub>4</sub> M 324.376Constit. of the liverwort, *Radula javanica*. Cryst. (hexane).Mp 147-149°. [ $\alpha$ ]<sub>D</sub> +7.6° (c, 0.26 in MeOH).Asakawa, Y. et al, *Phytochemistry*, 1991, **30**, 325 (isol, pmr, cmr)**Radulanolide****R-00030**

[85526-71-2]

C<sub>20</sub>H<sub>18</sub>O<sub>4</sub> M 322.360Constit. of *Radula complanata*. Cryst. Mp 103-104°.Asakawa, Y. et al, *Phytochemistry*, 1982, **21**, 2481.**Raffinose, 8CI****R-00031**

$\beta$ -D-Fructofuranosyl O- $\alpha$ -D-galactopyranosyl-(1 $\rightarrow$ 6)- $\alpha$ -D-glucopyranoside, 9CI. Melitriose. Gossypose. Melitose [512-69-6]



$C_{18}H_{32}O_{16}$  M 504.441

Occurs in sugar beet, cotton seeds, manna. Widely distributed in plants, esp. in the seeds. Prisms +  $5H_2O$ . Mp 78° (hydrate), Mp 118° (anhyd.).  $[\alpha]_D^{20} + 123^\circ$  (c, 2 in  $H_2O$ ) (anhyd.). Sweet taste. Does not reduce Fehling's soln. Invertase or raffinase gives Melibiose and Fructose.

**Undeca-Me:** Bp<sub>2</sub> 238-240°.  $[\alpha]_D^{17} + 128.4^\circ$  (c, 1 in  $H_2O$ ).

**Undeca-Ac:** Mp 101°.  $[\alpha]_D^{20} + 100^\circ$  (c, 8 in EtOH).

Tanret, G., *Bull. Soc. Chim. Fr.*, 1895, 261.

Haworth, N.W., *J. Chem. Soc.*, 1923, 3125.

French, D., *Adv. Carbohydr. Chem. Biochem.*, 1954, 9, 149 (rev)

Duperon, R., *C. R. Hebd. Seances Acad. Sci.*, 1955, 241, 1817 (occur)

Bourne, E.J. et al, *Biochem. J.*, 1965, 97, 802 (biosynth)

Karrer, W. et al, *Konstitution und Vorkommen der Organischen Pflanzenstoffe*, 2nd Ed., Birkhäuser Verlag, Basel, 1972-1985, no. 671 (occur)

Kamerling, J.P. et al, *Tetrahedron*, 1972, 28, 4375 (ms)

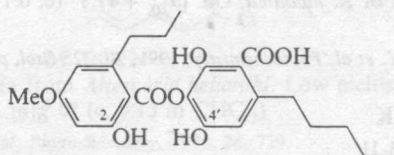
Bock, K. et al, *Tetrahedron Lett.*, 1973, 1037 (cmr)

Anteunis, M. et al, *Carbohydr. Res.*, 1975, 44, 101 (pmr)

## Ramalinolic acid

R-00032

2,4-Dihydroxy-3-[(2-hydroxy-4-methoxy-6-propylbenzoyl)oxy]-6-pentylbenzoic acid, 9CI  
[552-56-7]



$C_{23}H_{28}O_8$  M 432.469

Cryst. ( $C_6H_6$ ). Mp 164°.

2-Me ether: [2879-80-3]. **Merochlorophaeic acid**

$C_{24}H_{30}O_8$  M 446.496

Isol. from *Cladonia merochlorophaea*. Strong inhibitor of prostaglandin biosynth. Plates (MeOH aq.), needles ( $C_6H_6$ /cyclohexane). Mp 164-166°.

4'-Me ether: [486-36-2]. **Homosekikaic acid**. Nemoxyinic acid

$C_{24}H_{30}O_8$  M 446.496

Constit. of *Cladonia subpityrea* and *Cenomyce fimbriata*. Plates (pet. ether). Mp 133-134°.

4'-Me ether, Me ester: Mp 106°.

2,4'-Di-Me ether: [548-85-6]. **Boninic acid**

$C_{25}H_{32}O_8$  M 460.523

Constit. of *Ramalina boninensis*. Plates ( $C_6H_6$ /pet. ether). Mp 134.5°.

2,4'-Di-Me ether, Me ester: Plates (EtOH). Mp 86°.

Tri-Me ether, Me ester: Cryst. (MeOH). Mp 75°.

O-De-Me, 2-Me ether: [19833-81-9]. **Paludosic acid**

$C_{23}H_{28}O_8$  M 432.469

Isol. from *R. paludosa*. Needles (cyclohexane/ $C_6H_6$ /EtOAc). Mp 170-171° (158.5-159.5°). Mp. depends on solv. of recryst.

O-De-Me, 4'-Me ether: [103538-13-2]. **4'-O-Methylnorhomosekikaic acid**

$C_{23}H_{28}O_8$  M 432.469

Trace metab. of *R. luciae*. Cryst. ( $CHCl_3/CCl_4$ ). Mp 176°.

O-De-Me, 2,4'-di-Me ether: [69563-44-6]. **4'-O-Methylpaludosic acid**

$C_{24}H_{30}O_8$  M 446.496

Isol. from *R. asahinae*. Cryst. ( $C_6H_6$ /cyclohexane/ $Me_2CO$ ). Mp 138-139°.

Asahina, Y. et al, *Ber.*, 1936, 69, 1896; 1937, 70, 1815, 1821 (isol, synth)

Shibata, S. et al, *Phytochemistry*, 1965, 4, 133 (*Merochlorophaeic acid*)

Culberson, C.F., *Bryologist*, 1967, 70, 397 (isol, *Paludosic acid*)

Huneck, S. et al, *Z. Naturforsch.*, B, 1968, 23, 717 (pmr, *Merochlorophaeic acid*)

Elix, J.A. et al, *Aust. J. Chem.*, 1975, 28, 399 (synth, pmr, ms)

Chester, D.O. et al, *Aust. J. Chem.*, 1978, 31, 2745 (4'-O-Methylpaludosic acid, 4'-O-Methylnorhomosekikaic acid)

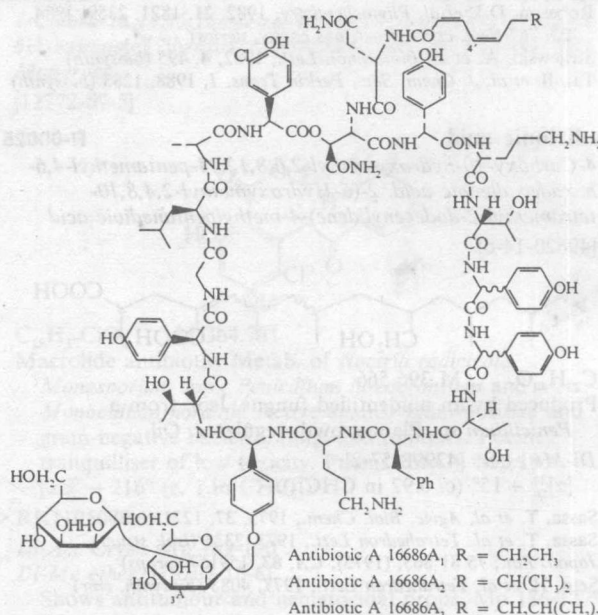
Shibuya, M. et al, *Chem. Pharm. Bull.*, 1983, 31, 407 (pharmacol, *Merochlorophaeic acid*)

## Ramoplanin, INN

R-00033

A 16686. Antibiotic A 16686

[76168-82-6]



Glycolipodepsipeptide antibiotic complex. Prod. by *Actinoplanes* sp. ATCC 33076. Active against gram-positive bacteria. Shows antiplaque activity. Inhibits cell wall biosynthesis.

**Antibiotic A 16686A<sub>1</sub>** [81988-87-6]

$C_{118}H_{152}ClN_{21}O_{40}$  M 2540.068

Powder. Mp 210-220° dec.  $[\alpha]_D^{20} + 57^\circ$  (c, 0.51 in  $H_2O$ ). Minor component of complex.  $\lambda_{max}$  233, 266 nm (MeOH).

**Antibiotic A 16686A<sub>2</sub>** [81988-88-7]

$C_{119}H_{154}ClN_{21}O_{40}$  M 2554.095

Powder. Mp 210-220° dec.  $[\alpha]_D^{20} + 73^\circ$  (c, 0.49 in  $H_2O$ ). Major component of complex.

B.HCl: Mp 250°.

4'E-Isomer, 3'A-O- $\alpha$ -D-mannopyranoside: [135789-09-2].

**Ramoplanose**. UK 71903. Antibiotic UK 71903

$C_{125}H_{164}ClN_{21}O_{45}$  M 2716.237

Prod. by A. sp.

**Antibiotic A 16686A<sub>3</sub>** [81988-89-8]

$C_{120}H_{156}ClN_{21}O_{40}$  M 2568.122

Powder. Mp 220°.  $[\alpha]_D^{20} + 50^\circ$  (c, 0.48 in 0.1M HCl). Minor component of complex.  $\lambda_{max}$  234, 267 nm (MeOH).

Ger. Pat., 3 013 246, (1980); CA, 94, 63774 (isol, props)

Eur. Pat., 46 201, (1982); CA, 97, 4672 (isol, props, nmr)

Pallanza, R. et al, *Antimicrob. Agents Chemother.*, 1984, 26, 462 (props)

Cavalleri, B. et al, *J. Antibiot.*, 1984, 37, 309 (isol)

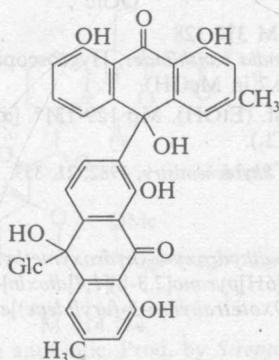
Pallanza, R. et al, *J. Antibiot.*, 1984, 37, 318 (isol, props)

- Ciabatti, R. *et al*, *J. Antibiot.*, 1989, **42**, 254, 268 (pmr, cmr, ms, struct)  
 Skelton, N.J. *et al*, *J. Am. Chem. Soc.*, 1991, **113**, 7522 (Ramoplanose)

**Ramosin†**

[120163-26-0]

R-00034

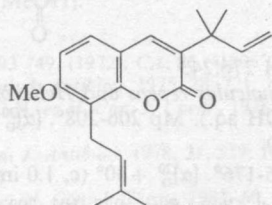
 $C_{36}H_{32}O_{13}$  M 672.641

Constit. of *Asphodelus ramosus*. Yellow amorph. solid.  $[\alpha]_D^{25}$  -74° (c, 1.1 in MeOH).

Adinolfi, M. *et al*, *Phytochemistry*, 1989, **28**, 284.**Ramosinin**

[70894-25-6]

R-00035

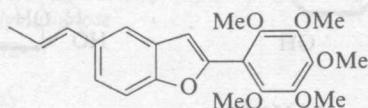
 $C_{20}H_{24}O_3$  M 312.408

Constit. of *Haplophyllum ramosissimum*. Cryst. (EtOAc). Mp 85-86°.

Gashimov, N.F. *et al*, *Khim. Prir. Soedin.*, 1979, **15**, 15; *Chem. Nat. Compd. (Engl. Transl.)*, 11 (isol, pmr)Salvá, J. *et al*, *Heterocycles*, 1990, **31**, 255 (synth)**Ramosissin**

2-(Pentamethoxyphenyl)-5-(1-propenyl)benzofuran, 9CI  
 [110784-16-2]

R-00036

 $C_{22}H_{24}O_6$  M 384.428

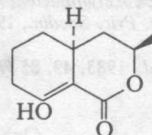
Constit. of *Krameria ramosissima*. Cryst. Mp 92-94°.

Achenbach, H. *et al*, *Phytochemistry*, 1987, **26**, 2041.**Ramulosin**

R-00037

3,4,4a,5,6,7-Hexahydro-8-hydroxy-3-methyl-1H-2-benzopyran-1-one, 9CI

[29914-01-0]

 $C_{10}H_{14}O_3$  M 182.219

Metab. of *Pestalotia ramulosa*. Plates (Me<sub>2</sub>CO aq.). Mp 120-121°.  $[\alpha]_D^{25}$  +18° (c, 2.8 in EtOH).

Benzoyl: Needles (EtOH aq.). Mp 82-83°.

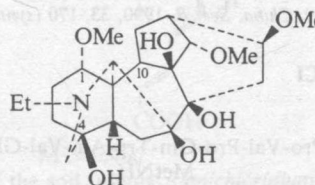
4-Nitrobenzoyl: Pale-yellow needles (Me<sub>2</sub>CO aq.). Mp 135-136°.

Stodola, F.H. *et al*, *Biochem. J.*, 1964, **93**, 92 (isol, struct)Tanenbaum, S.W. *et al*, *Tetrahedron Lett.*, 1970, 2377 (pmr)Findlay, J.A. *et al*, *Can. J. Chem.*, 1976, **54**, 3419 (abs config)Mori, K. *et al*, *Tetrahedron*, 1985, **41**, 5295 (synth)Takano, S. *et al*, *Heterocycles*, 1989, **29**, 2101 (synth, bibl)Asaoka, M. *et al*, *Tetrahedron*, 1990, **46**, 1541 (synth)**Ranaconine**

R-00038

Hydroxylappaconine

[70474-32-7]

 $C_{23}H_{37}NO_7$  M 439.548Alkaloid detected by glc-ms in *Aconitum leucostomum*.

Alkaline hydrol. product of Ranaconitine

(Ranunculaceae). Mp 107-109°.

O<sup>4</sup>-(2-Aminobenzoyl): [82872-80-8]. N-Deacetylranaconitine $C_{30}H_{42}N_2O_8$  M 558.670Alkaloid from the roots of *A. finetianum*

(Ranunculaceae). Shows good analgesic activity. Mp 125-127°.

O<sup>4</sup>-(2-Acetamidobenzoyl): [1360-76-5]. Ranaconitine $C_{32}H_{44}N_2O_9$  M 600.708Alkaloid from *A. ranunculaefolium*, *A. sinomontanum*, *A.**finetianum* and *A. barbatum* var. *puberulum*(Ranunculaceae). Mp 132-134°.  $[\alpha]_D^{27}$  +33.2° (c, 0.5 in CHCl<sub>3</sub>).O<sup>4</sup>-(2-Acetamidobenzoyl), 1-epimer: [83685-20-5].

Puberanine

 $C_{32}H_{44}N_2O_9$  M 600.708Alkaloid from the roots of *A. barbatum* var. *puberulum*(Ranunculaceae). Amorph.  $[\alpha]_D^{20}$  +16.6° (c, 0.6 in CHCl<sub>3</sub>).10β-Hydroxy, O<sup>4</sup>-(2-aminobenzoyl): [82872-81-9]. N-

Deacetylfinaconitine

 $C_{30}H_{42}N_2O_9$  M 574.670Alkaloid from the roots of *A. finetianum*

(Ranunculaceae). Mp 121-123°.

10β-Hydroxy, O<sup>4</sup>-(2-acetamidobenzoyl): [81161-27-5].

Finaconitine. 10β-Hydroxyranaconitine

 $C_{32}H_{44}N_2O_{10}$  M 616.707Alkaloid from the roots of *A. finetianum*

(Ranunculaceae). Mp 220-221°.

Mollov, N. *et al*, *Dokl. Bolg. Akad. Nauk*, 1964, **17**, 251; *CA*, **61**, 12324h (isol, ir, Ranaconitine)

- Pelletier, S.W. *et al*, *Tetrahedron Lett.*, 1978, 5045 (ir, pmr, cmr, struct, Ranaconitine)  
 Pelletier, S.W. *et al*, *Can. J. Chem.*, 1979, 57, 1652 (cmr)  
 Jiang, S. *et al*, *Yaoxue Xuebao*, 1982, 17, 283; 1983, 18, 440; *CA*, 97, 20736a; 100, 20505e (Finaconitine, Deacetylranaconitine, Deacetylfinaconitine)  
 Plugar, V.N. *et al*, *Khim. Prir. Soedin.*, 1982, 80; *CA*, 97, 39185s (glc, ms)  
 Yu, D. *et al*, *Planta Med.*, 1983, 49, 85 (Puberanine)

**Ranakinin R**

R-00039

[60972-29-4]

H-Arg-Pro-Pro-Gly-Phe-Thr-Pro-Phe-Arg-Ile-Ala-Pro-Glu-Ile-Val-OH

Bradykinin analogue. Isol. from *Rana rugosa* skin.Yanaihara, C. *et al*, *Adv. Exp. Med. Biol.*, Sect. A, 1978, 120, 185 (synth)Yasuhara, T. *et al*, *Chem. Pharm. Bull.*, 1979, 27, 486 (isol, struct)**Ranamargarin**

R-00040

[132151-82-7]

H-Asp-Asp-Ala-Ser-Asp-Arg-Ala-Lys-Lys-Phe-Tyr-Gly-Leu-Met-NH<sub>2</sub>

Isol. from frog skin.

Tong, Y. *et al*, *Regul. Pept.*, 1988, 22, 182 (isol)Lu, Y. *et al*, *Sci. China, Ser. B*, 1990, 33, 170 (synth)**Ranatensin, 9CI**

R-00041

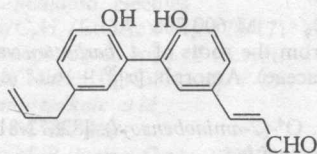
[29451-71-6]

H-5-OxoPro-Val-Pro-Gln-Trp-Ala-Val-Gly-His-Phe-MetNH<sub>2</sub>C<sub>61</sub>H<sub>84</sub>N<sub>16</sub>O<sub>13</sub>S M 1281.500Isol. from frog skin (*Rana pipiens*). Vasoactive peptide related to Bombesin.Nakajima, T. *et al*, *Fed. Proc.*, 1970, 23, 282 (isol)Rivier, J.E. *et al*, *Biochemistry*, 1978, 17, 1766 (synth, props)Hernandez, O. *et al*, *J. Labelled Compd.*, 1984, 7, 893 (hplc)Howard, J.M. *et al*, *Aust. J. Pharm.*, 1985, 248, G196 (biol activity)Krane, I.M. *et al*, *J. Biol. Chem.*, 1988, 263, 13317; 1990, 265, 7091 (cloning)Ersparmer, G.F. *et al*, *Regul. Pept.*, 1988, 21, 1 (biol activity)**Randainal**

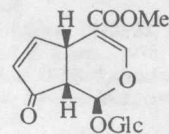
R-00042

3-[2',6'-Dihydroxy-5'-(2-propenyl)[1,1'-biphenyl-3-yl]]-2-propenal, 9CI

[87562-13-8]

C<sub>18</sub>H<sub>16</sub>O<sub>3</sub> M 280.323Constit. of *Sassafras randaiense*. Cryst. (EtOAc/hexane). Mp 137-139°.Alcohol: [93753-25-4]. **Randaianol**C<sub>18</sub>H<sub>18</sub>O<sub>3</sub> M 282.338Isol. from *S. randaiense*. Oil.Chen, F.-C. *et al*, *Phytochemistry*, 1983, 22, 616.El-Ferally, F.S., *Phytochemistry*, 1984, 23, 2329 (Randainol)**Randioside**

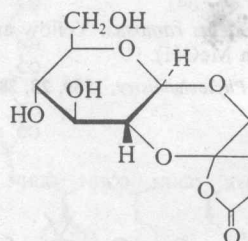
[82345-53-7]

C<sub>16</sub>H<sub>20</sub>O<sub>10</sub> M 372.328Constit. of *Randia canthioides*. Hygroscopic powder. [α]<sub>D</sub><sup>17</sup> -29.4° (c, 0.2 in MeOH).Tetra-Ac: Cryst. (EtOH). Mp 129-131°. [α]<sub>D</sub><sup>30</sup> -56.8° (c, 0.56 in CHCl<sub>3</sub>).Uesato, S. *et al*, *Phytochemistry*, 1982, 21, 353.**Ranuncoside**

R-00044

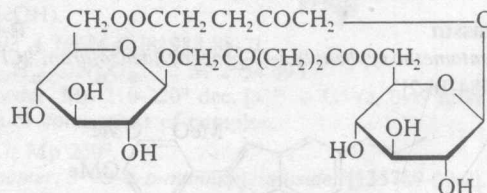
Hexahydro-7,8-dihydroxy-6-(hydroxymethyl)spiro[furan-2(5H),2'(3'H)-[6H]pyrano[2,3-b][1,4]dioxin]-5-one, 9CI. 1,2-O-[2-(S)-2-(2-Oxotetrahydro-5-furylidene)]ethylene-α-D-glucopyranose

[35879-55-1]

C<sub>11</sub>H<sub>16</sub>O<sub>8</sub> M 276.243Constit. of *Ranunculus repens* and *Helleborus foetidus*.Needles (EtOH aq.). Mp 206-208°. [α]<sub>D</sub><sup>20</sup> +40.2° (c, 0.5 in EtOH aq.).Tri-Ac: Mp 175-176°. [α]<sub>D</sub><sup>20</sup> +30° (c, 1.0 in CHCl<sub>3</sub>).Tschesche, R. *et al*, *Chem. Ber.*, 1972, 105, 290 (isol, struct, pmr, ms)Mariezcurrena, R.A. *et al*, *Acta Crystallogr., Sect. B*, 1973, 29, 1030.**Ranunculoside, 9CI**

R-00045

[35879-56-2]

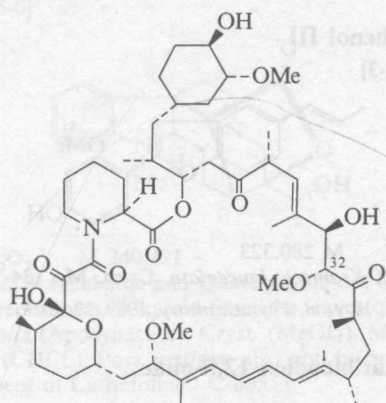
C<sub>22</sub>H<sub>32</sub>O<sub>16</sub> M 552.485Extracted from *Ranunculus repens* and *Helleborus foetidus*.Mp 153-154°. [α]<sub>D</sub><sup>20</sup> -4° (c, 1.0 in MeOH).Hexa-Ac: Mp 233-234°. [α]<sub>D</sub><sup>20</sup> -3.0° (c, 0.94 in CHCl<sub>3</sub>).Tschesche, R. *et al*, *Chem. Ber.*, 1972, 105, 290 (synth, pmr)

**Rapamycin**

**R-00046**

AY 22989. Antibiotic AY 22989

[53123-88-9]



$C_{51}H_{79}NO_{13}$  M 914.184

Polyene-type antibiotic. Prod. by *Streptomyces hygroscopicus*. Antifungal agent. Cryst. (Et<sub>2</sub>O). Mp 183-185°.  $[\alpha]_D^{25} - 58.2^\circ$  (MeOH).

▷ VE6250000.

32-Demethoxy: [83482-58-0]. **Demethoxyrapamycin**. AY 24668. Antibiotic AY 24668

$C_{50}H_{77}NO_{12}$  M 884.158

From *S. hygroscopicus*. Antifungal agent with v. slight antitumour activity. Cryst. (Et<sub>2</sub>O). Mp 122-124°.  $[\alpha]_D^{25} - 124.4^\circ$  (MeOH).

[85537-35-5]

U.S. Pat., 3 993 749, (1972); CA, 86, 41806 (synth, ir, pmr)  
Vézina, C. et al. J. Antibiot., 1975, 28, 721; 1983, 36, 351 (isol)  
Swindells, D.C.N. et al. Can. J. Chem., 1978, 56, 2491 (cryst struct)

Baker, H. et al. J. Antibiot., 1978, 31, 539; 1979, 32, 630 (pharmacol)

Findlay, J.A. et al. Can. J. Chem., 1982, 60, 2046 (uv, ir, pmr, cmr, ms, cd)

U.S. Pat., 4 375 464, (1983); CA, 98, 177506 (isol)

McAlpine, J.B. et al. J. Antibiot., 1991, 44, 688 (pmr, cmr)

Paiva, N.L. et al. J. Nat. Prod. (Lloydia), 1991, 54, 167 (cmr, biosynth)

Curran, D.P. et al. Tetrahedron Lett., 1992, 33, 2295 (synth)

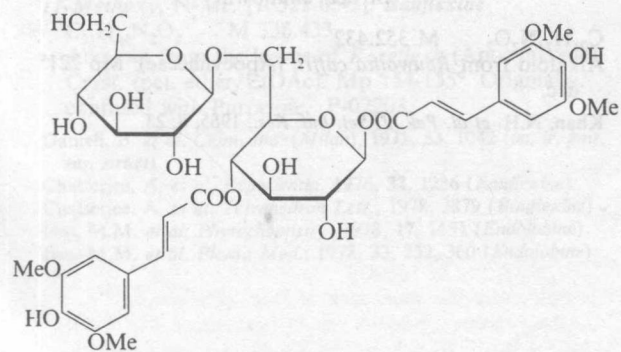
Lewis, R.J., Sax's Dangerous Properties of Industrial Materials, 8th Ed., Van Nostrand-Reinhold, 1992, RBK000.

**Raphanusol A**

**R-00047**

6-O-β-D-glucopyranosyl β-D-glucopyranose 1,4-bis[3-(4-hydroxy-3,5-dimethoxyphenyl)-2-propenoate], 9CI. 1,4-Di-O-sinapoylgentiobiose

[74565-72-3]



$C_{34}H_{42}O_{19}$  M 754.694

Isol. from *Raphanus sativus* (Sakura jima radish) seedling.

Endogenous hypocotyl growth inhibitor. Powder

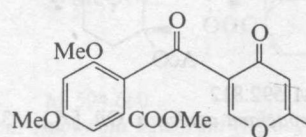
(Me<sub>2</sub>CO/C<sub>6</sub>H<sub>6</sub>). Mp 137-138°.  $[\alpha]_D^{22} + 79.4^\circ$  (c, 0.23 in MeOH).

Hase, T. et al, Phytochemistry, 1982, 21, 1021 (isol, struct)

Hase, T. et al, CA, 1985, 103, 138518b.

**Rapicone**

**R-00048**



$C_{17}H_{16}O_7$  M 332.309

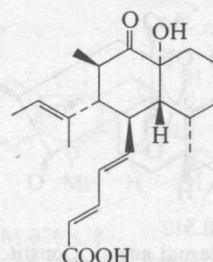
Metab. of *Ramichloridium apiculatum*. Needles (C<sub>6</sub>H<sub>6</sub>). Mp 162-163°.

Nozawa, K. et al, Phytochemistry, 1992, 31, 4178 (isol, pmr, cmr)

**Rapiculine**

**R-00049**

[133961-70-3]



$C_{21}H_{30}O_4$  M 346.466

Metab. of the soil fungus *Ramichloridium apiculatum*.

Needles (C<sub>6</sub>H<sub>6</sub>). Mp 199°.  $[\alpha]_D^{23} - 73^\circ$  (c, 0.35 in MeOH).

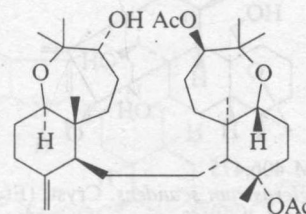
Nozawa, K. et al, J. Chem. Soc., Perkin Trans. 1, 1991, 537 (isol, pmr, cmr)

Nozawa, K. et al, Phytochemistry, 1992, 31, 4177 (abs config)

**Raspacionin**

**R-00050**

[132210-64-1]

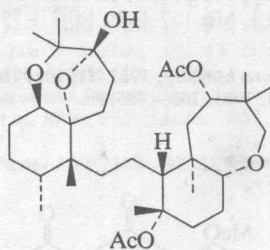


$C_{34}H_{56}O_7$  M 576.812

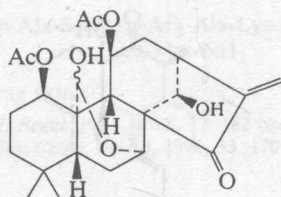
Constit. of sponge *Raspaciona aculeata*. Prisms (heptane).

Mp 188-189°.  $[\alpha]_D + 31.4^\circ$  (c, 1.5 in CHCl<sub>3</sub>).

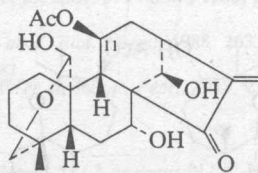
Cimino, G. et al, Tetrahedron Lett., 1990, 31, 6565 (cryst struct)

**Raspacionin A****R-00051** $C_{34}H_{56}O_8$  M 592.812Constit. of *Raspaciona aculeata*. Oil.  $[\alpha]_D -3.95^\circ$  (c, 1.17 in  $CHCl_3$ ).Cimino, G. et al, *Tetrahedron*, 1992, 48, 9013 (isol, pmr, cmr, crystal structure)**Rastronol D****R-00052**ent-1 $\alpha$ ,11 $\alpha$ -Diacetoxy-7 $\beta$ ,20-epoxy-14 $\alpha$ ,20-dihydroxy-16-kauren-15-one

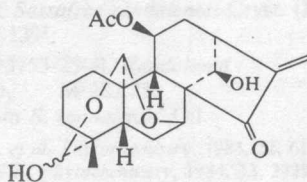
[59859-96-0]

 $C_{24}H_{32}O_8$  M 448.512A cyclic (20 $\rightarrow$ 7) internal acetal. Constit. of *Englerastrum scandens*. Cryst. (Me<sub>2</sub>CO). Mp 198.1-199.2°.  $[\alpha]_D -80.5^\circ$ .Nomoto, K. et al, *Helv. Chim. Acta*, 1976, 59, 772.**Rastronol G****R-00053**ent-11 $\alpha$ -Acetoxy-19,20-epoxy-7 $\beta$ ,14 $\alpha$ ,20-trihydroxy-16-kauren-15-one

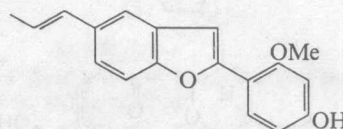
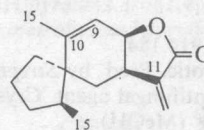
[59859-99-3]

 $C_{22}H_{30}O_7$  M 406.475Constit. of *Englerastrum scandens*. Cryst. (Et<sub>2</sub>O). Mp 228.4-228.9°.  $[\alpha]_D -146^\circ$ .Nomoto, K. et al, *Helv. Chim. Acta*, 1976, 59, 772.**Rastronol H****R-00054**ent-11 $\alpha$ -Acetoxy-7 $\beta$ ,20:19,20-diepoxy-14 $\alpha$ ,19-dihydroxy-16-kauren-15-one

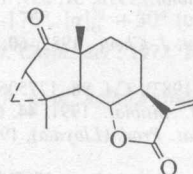
[59860-00-3]

 $C_{22}H_{28}O_7$  M 404.459Constit. of *Englerastrum scandens*. Cryst. (MeOH). Mp 173.5-174.5°.  $[\alpha]_D -77.2^\circ$ .Nomoto, K. et al, *Helv. Chim. Acta*, 1976, 59, 772.**Ratanhiaphenol III****R-00055**

[91432-06-3]

 $C_{18}H_{16}O_3$  M 280.323Constit. of *Krameria lanceolata*. Cryst. Mp 134-135°.Achenbach, H. et al, *Phytochemistry*, 1989, 28, 1959.**9,11(13)-Ratibidadien-12,8-olide****R-00056** $C_{15}H_{20}O_2$  M 232.322(4*S*,5*S*,7*R*,8*R*)-form [94137-83-4]Constit. of *Ratibida columnifera*. Gum. $\Delta^{10,15}$ -Isomer: [94137-79-8]. 10(15),11(13)-Ratibidadien-12,8-olide $C_{15}H_{20}O_2$  M 232.322Constit. of *R. columnifera*. Gum.Herz, W. et al, *J. Org. Chem.*, 1985, 50, 610.**Ratibinolide****R-00057**

[130170-06-8]

 $C_{15}H_{18}O_3$  M 246.305Constit. of *Ratibida latipalearis*. Cryst. Mp 142-144°.  $[\alpha]_D^{25} +94^\circ$  ( $CHCl_3$ ).Mata, R. et al, *Heterocycles*, 1990, 31, 1111 (isol, pmr, cmr, crystal structure)**Raucaffridine****R-00058**

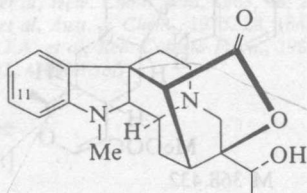
Struct. unknown

 $C_{21}H_{24}N_2O_3$  M 352.432Alkaloid from *Rauwolfia caffra* (Apocynaceae). Mp 221° dec.Khan, N.H. et al, *Pak. J. Sci. Ind. Res.*, 1965, 8, 23.

**Raucubaine**

R-00059

1,2,19,20-Tetrahydro-19,20-dihydroxy-1-methylakuummolan-17-oic acid  $\gamma$ -lactone, 9CI. Quaternoline  
[75418-95-0]



$C_{20}H_{24}N_2O_3$  M 340.421

Identity of Raucubaine and Quaternoline not certain, may be stereoisomers. Alkaloid from the leaves of *Rauwolfia salicifolia* (Apocynaceae). Cryst. (MeOH). Mp 224°.  $[\alpha]_D^{20}$  –18° (CHCl<sub>3</sub>). Poss. artifact, also obt. by acid treatment of Cathafoline, C-00553.

11-Methoxy: [56283-51-3]. **Caberoline**

$C_{21}H_{26}N_2O_4$  M 370.447

Alkaloid from *Cabucala* spp. and *Alstonia plumosa* (Apocynaceae). Poss. artifact; also obt. by acid treatment of Caberine.

[57498-99-4]

Mamatas-Kalamaras, S. et al, *Phytochemistry*, 1975, **14**, 1849 (Quaternoline)

Kutney, J.P. et al, *Heterocycles*, 1980, **14**, 1309 (uv, ir, pmr, cd, ms, cryst struct)

Pauptit, R.A. et al, *Can. J. Chem.*, 1981, **59**, 1007 (cryst struct)

Jacquier, M.J. et al, *Phytochemistry*, 1982, **21**, 2973 (isol)

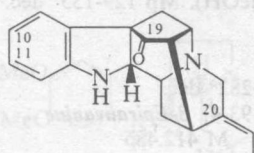
Sierra, P. et al, *Collect. Czech. Chem. Commun.*, 1982, **47**, 2912 (isol, spectra, abs config)

**Rauflorine**

R-00060

19,20-Didehydro-1-demethylajmalan-17-one, 9CI

[36063-54-4]



$C_{19}H_{20}N_2O$  M 292.380

Alkaloid from the root bark of *Rauwolfia confertiflora* (Apocynaceae). Cryst. (MeOH). Mp 221°.  $[\alpha]_D^{20}$  +312° (c, 1.54 in CHCl<sub>3</sub>).

10-Methoxy: [67627-71-8]. **Endolobine**

$C_{20}H_{22}N_2O_2$  M 322.406

Alkaloid from the leaves and stem bark of *R. cumminsii*, and the leaves of *R. mombasiana* (Apocynaceae).

11-Methoxy, N-Me: [70522-05-3]. **Rauflexine**

$C_{21}H_{24}N_2O_2$  M 336.433

Alkaloid from the leaves of *R. reflexa* (Apocynaceae). Cryst. (pet. ether/EtOAc). Mp 154-155°. Originally confused with Purpeline, P-02203.

Danieli, B. et al, *Chim. Ind. (Milan)*, 1971, **53**, 1042 (uv, ir, pmr, ms, struct)

Chatterjee, A. et al, *Experientia*, 1976, **32**, 1236 (Rauflexine)

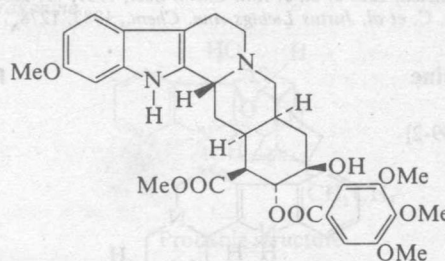
Chatterjee, A. et al, *Tetrahedron Lett.*, 1978, 3879 (Rauflexine)

Iwu, M.M. et al, *Phytochemistry*, 1978, **17**, 1651 (Endolobine)

Iwu, M.M. et al, *Planta Med.*, 1978, **33**, 232, 360 (Endolobine)

**Raugustine**

R-00061



$C_{32}H_{38}N_2O_9$  M 594.660

Alkaloid from *Rauwolfia ligustrina* (Apocynaceae). Cryst. + 1H<sub>2</sub>O. Mp 160-170° dec.  $[\alpha]_D^{24}$  –50° (c, 0.609 in CHCl<sub>3</sub>).

B, HNO<sub>3</sub>: Mp 262-263° dec. (in vacuo).

Mono-Ac: Mp 232-234° (block).

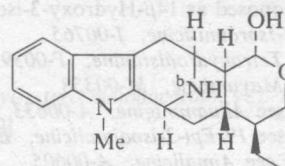
Müller, J.M., *Experientia*, 1957, **12**, 479 (isol, uv)

**Raumacline**

R-00062

4-Demethyl-20,21-dihydroalstophyllan-17-ol, 9CI

[132923-02-5]



$C_{20}H_{26}N_2O_2$  M 326.438

Alkaloid from cell cultures of *Rauwolfia serpentina* (Apocynaceae). Mp 197-200°.

N<sup>b</sup>-Me: [132943-65-8]. N<sup>b</sup>-Methylraumacline

$C_{21}H_{28}N_2O_2$  M 340.464

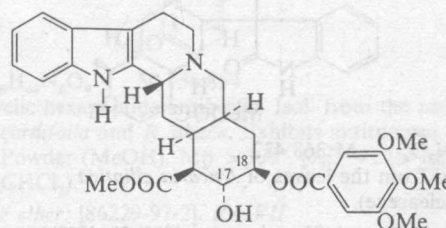
Alkaloid from cell cultures of *R. serpentina* (Apocynaceae). Amorph. solid.

Polz, L. et al, *Tetrahedron Lett.*, 1990, **31**, 6693 (isol, uv, cd, pmr, cmr, struct)

**Raunescine†**

R-00063

[117-73-7]



$C_{31}H_{36}N_2O_8$  M 564.634

Alkaloid from *Rauwolfia canescens* and *R. tetraphylla*, also detected in *R. ligustrina* (Apocynaceae). Cryst. (MeOH aq.). Mp 160-170°.  $[\alpha]_D$  –74°.

B, HNO<sub>3</sub>: Mp 223-225°.  $[\alpha]_D$  –80° (c, 1 in 5M AcOH).

Mono-Ac: Mp 157-162°.  $[\alpha]_D$  –157° (c, 1 in CHCl<sub>3</sub>).

O<sup>17</sup>-(3,4,5-Trimethoxybenzoyl), O<sup>18</sup>-de-(3,4,5-trimethoxybenzoyl): [483-07-8]. **Isoraunescine**

$C_{31}H_{36}N_2O_8$  M 564.634

Alkaloid from *R. canescens*, also detected in *R. ligustrina* (Apocynaceae). Cryst. (MeOH aq.). Mp 241-242°.  $[\alpha]_D$  –70°.

O<sup>17</sup>-Me: see Deserpidine, D-00554

Hosanky, N. et al, *J. Am. Pharm. Assoc.*, 1955, **44**, 635 (isol, uv)

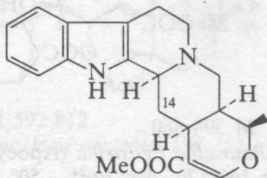
Huebner, C.F. *et al*, *J. Am. Chem. Soc.*, 1957, **79**, 250 (*struct*)  
 van Tamelen, E.E. *et al*, *J. Am. Chem. Soc.*, 1957, **79**, 5256 (*struct*)  
 Szántay, C. *et al*, *Justus Liebigs Ann. Chem.*, 1983, 1278.

**Rauniticine**

Ervine

[5299-09-2]

R-00064

C<sub>21</sub>H<sub>24</sub>N<sub>2</sub>O<sub>3</sub> M 352.432

Alkaloid from *Rauwolfia nitida*, *Uncaria elliptica*, *U. attenuata*, *Vinca libanotica*, *V. major* and *Veratrum album lobelianum* (Apocynaceae, Naucleaceae, Liliaceae).  
 Cryst. (MeOH). Mp 233-235°. [ $\alpha$ ]<sub>D</sub> -6.6° (Py), [ $\alpha$ ]<sub>D</sub> -38.4° (CHCl<sub>3</sub>).

14 $\alpha$ -Hydroxy: 14 $\alpha$ -HydroxyrauniticineC<sub>21</sub>H<sub>24</sub>N<sub>2</sub>O<sub>4</sub> M 368.432

Alkaloid from the leaves of *U. attenuata* (Naucleaceae).  
 Originally proposed as 14 $\beta$ -Hydroxy-3-isorauniticine.

3-Epimer: see 3-Isorauniticine, I-00765

19-Epimer: see Tetrahydroalstonine, T-00391

20-Epimer: see Mayumbine, M-00358

3,19-Diepimer: see Akuammigine, A-00633

3,20-Diepimer: see 19-Epi-3-isoajmalicine, E-00332

19,20-Diepimer: see Ajmalicine, A-00605

Salkin, R. *et al*, *J. Pharm. Sci.*, 1961, **50**, 1038 (*isol*, *uv*, *ir*, *struct*)  
 Shamma, M. *et al*, *J. Am. Chem. Soc.*, 1963, **85**, 2507 (*stereochem*)  
 Finch, N. *et al*, *Tetrahedron*, 1966, **22**, 1327 (*uv*, *ord*, *stereochem*)  
 Aynilian, G.H. *et al*, *J. Nat. Prod. (Lloydia)*, 1974, **37**, 299 (*isol*, *w*, *ir*, *ms*)

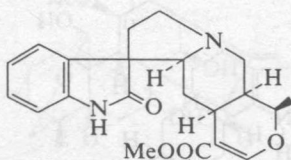
Ponglux, D. *et al*, *Phytochemistry*, 1980, **19**, 2013 (*isol*, *uv*, *pmr*, *ms*)

Phillipson, J.D. *et al*, *Phytochemistry*, 1983, **22**, 1809 (*isol*, *pmr*)

Yamanaka, E. *et al*, *Chem. Pharm. Bull.*, 1986, **34**, 3713 (14 $\alpha$ -Hydroxyrauniticine)

**Rauniticine oxindole A**

R-00065

C<sub>21</sub>H<sub>24</sub>N<sub>2</sub>O<sub>4</sub> M 368.432

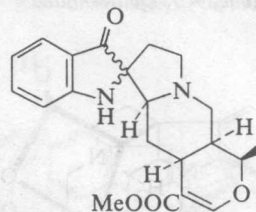
Alkaloid from the leaves of *Uncaria elliptica* (Naucleaceae).

Phillipson, J.D. *et al*, *Phytochemistry*, 1983, **22**, 1809 (*uv*, *pmr*, *ms*, *struct*)

**Rauniticine pseudoindoxyl**

R-00066

10,11-Didemethoxy-3-isoreserpine pseudoindoxyl, 9CI

C<sub>21</sub>H<sub>24</sub>N<sub>2</sub>O<sub>4</sub> M 368.432

Alkaloid from the leaves of *Uncaria elliptica* (Naucleaceae).

3-Epimer: 3-Isorauniticine pseudoindoxyl

C<sub>21</sub>H<sub>24</sub>N<sub>2</sub>O<sub>4</sub> M 368.432

Alkaloid from leaves of *U. elliptica* (Naucleaceae).

3,19-Diepimer: Akuammigine pseudoindoxyl

C<sub>21</sub>H<sub>24</sub>N<sub>2</sub>O<sub>4</sub> M 368.432

Alkaloid from leaves of *U. elliptica* (Naucleaceae).

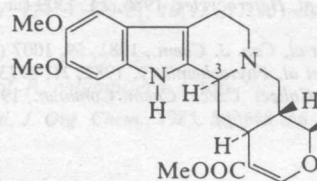
[5802-00-6, 88335-34-6, 88335-35-7, 88375-63-7]

Phillipson, J.D. *et al*, *Phytochemistry*, 1983, **22**, 1809 (*uv*, *pmr*, *ms*, *struct*, *synth*)

**Rauvanine**

R-00067

[3148-42-3]

C<sub>23</sub>H<sub>28</sub>N<sub>2</sub>O<sub>5</sub> M 412.485

Alkaloid from *Rauwolfia vomitoria* (Apocynaceae). Leaflets + 0.5H<sub>2</sub>O (MeOH). Mp 129-135° dec. [ $\alpha$ ]<sub>D</sub> +32.5° (c, 1 in CHCl<sub>3</sub>).

▷ RP5776500.

B,HCl: Mp 280-282° dec.

3-Epimer: [6835-93-4]. 3-Epirauvanine

C<sub>23</sub>H<sub>28</sub>N<sub>2</sub>O<sub>5</sub> M 412.485

Alkaloid from the bark of *Neisosperma oppositifolia* (Apocynaceae). Amorph. [ $\alpha$ ]<sub>D</sub> +111° (CHCl<sub>3</sub>).

Goutarel, R. *et al*, C. R. Hebd. Seances Acad. Sci., 1961, **253**, 2589 (*uv*, *ir*, *pmr*, *struct*)

Poisson, J. *et al*, *Bull. Soc. Chim. Fr.*, 1964, 2853 (*epimer*, *synth*, *ir*, *pmr*)

Finch, N. *et al*, *Tetrahedron*, 1966, **22**, 1327 (*uv*, *ord*, *stereochem*)

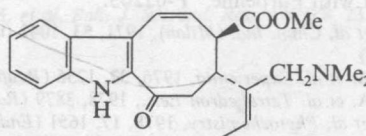
Amarasekara, A.S. *et al*, *Fitoterapia*, 1986, **57**, 55; *CA*, **105**, 39383p (*epimer*, *isol*)

**Rauviridine**

R-00068

Vobasine methine

[57567-12-1]

C<sub>22</sub>H<sub>26</sub>N<sub>2</sub>O<sub>3</sub> M 366.459

Alkaloid from the stem bark of *Rauwolfia viridis* (Apocynaceae). Needles (Et<sub>2</sub>O/pet. ether). Mp 145-146°. [ $\alpha$ ]<sub>D</sub> -103.7° (c, 0.9 in CHCl<sub>3</sub>). Data given relates to synthetic prod.

The Dictionary of Natural Products  
 is also available in a fully  
 substructure-searchable CD-ROM version

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 Marketing Department (SDD),  
 Chapman & Hall, for details