

INTERNATIONAL  
STANDARD

ISO  
1750

First edition  
1981-12-01

AMENDMENT 7  
2021-12

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**Pesticides and other agrochemicals —  
Common names**

**AMENDMENT 7**

*Produits phytosanitaires et assimilés — Noms communs*

*AMENDEMENT 7*

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[ISO 1750:1981/Amd 7:2021](https://standards.iteh.ai/catalog/standards/sist/7ef6811d-b949-47d5-9b84-e6af2387b7de/iso-1750-1981-amd-7-2021)

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Reference number  
ISO 1750:1981/Amd.7:2021(E)

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Phone: +41 22 749 01 11  
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The procedures used to develop this document and those intended for its further maintenance are described in the ISO/IEC Directives, Part 1. In particular, the different approval criteria needed for the different types of ISO documents should be noted. This document was drafted in accordance with the editorial rules of the ISO/IEC Directives, Part 2 (see [www.iso.org/directives](http://www.iso.org/directives)).

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This document was prepared by Technical Committee ISO/TC 81, *Common names for pesticides and other agrochemicals*.

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<https://standards.iteh.ai/catalog/standards/sist/7ef6811d-b949-47d5-9b84-e6af2387b7de/iso-1750-1981-amd-7-2021>

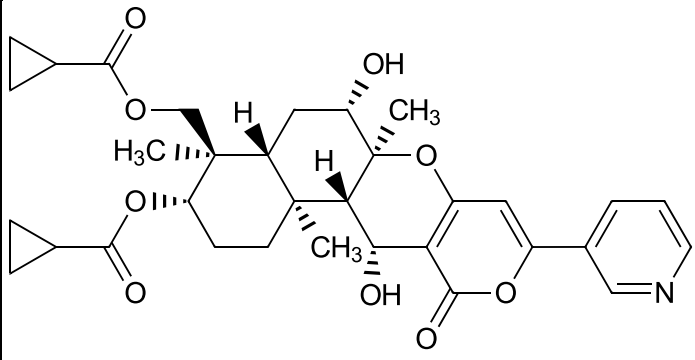
## Clause 4, Approved common names

Add these new common names approved since Amd.5.

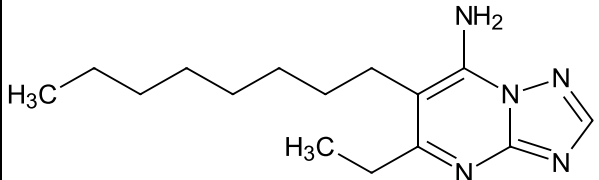
The common names in this section are those that have been approved since Amendment 5 to ISO 1750:1981.

<b>Common name</b>	<b>acynonapyr</b>
IUPAC PIN	(1 <i>R</i> ,3 <i>r</i> ,5 <i>S</i> )-3-[2-propoxy-4-(trifluoromethyl)phenoxy]-9-[[5-(trifluoromethyl)pyridin-2-yl]oxy]-9-azabicyclo[3.3.1]nonane
IUPAC name	9-[[5-(trifluoromethyl)-2-pyridyl]oxy]-3- <i>endo</i> -[[ $\alpha,\alpha,\alpha$ -trifluoro-2-propoxy- <i>p</i> -tolyl]oxy]-9-azabicyclo[3.3.1]nonane
CAS name	(3- <i>endo</i> )-3-[2-propoxy-4-(trifluoromethyl)phenoxy]-9-[[5-(trifluoromethyl)-2-pyridinyl]oxy]-9-azabicyclo[3.3.1]nonane
Structure	 <p style="text-align: center; color: red; font-weight: bold;">iTeh STANDARD PREVIEW (standards.iteh.ai)</p>
CAS Reg. No.	1332838-17-1 <a href="https://standards.iteh.ai/catalog/standards/sist/7ef6811d-b949-47d5-9b84-36af2387b7de/iso-1750-1981-amd-7-2021">ISO 1750:1981/Amd 7:2021</a>
Formula	C <sub>24</sub> H <sub>26</sub> F <sub>6</sub> N <sub>2</sub> O <sub>3</sub> <a href="https://standards.iteh.ai/catalog/standards/sist/7ef6811d-b949-47d5-9b84-36af2387b7de/iso-1750-1981-amd-7-2021">6af2387b7de/iso-1750-1981-amd-7-2021</a>
Use	acaricide
InChIKey	GIDAJLLAARKRMS-BWTSREIZSA-N
InChI	InChI=1S/C24H26F6N2O3/c1-2-10-33-21-11-15(23(25,26)27)6-8-20(21)34-19-12-17-4-3-5-18(13-19)32(17)35-22-9-7-16(14-31-22)24(28,29)30/h6-9,11,14,17-19H,2-5,10,12-13H2,1H3/t17-,18+,19+

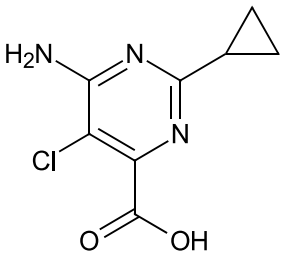
<b>Common name</b>	<b>afidopyropen</b>
IUPAC PIN	{{(3 <i>S</i> ,4 <i>R</i> ,4 <i>aR</i> ,6 <i>S</i> ,6 <i>aS</i> ,12 <i>R</i> ,12 <i>aS</i> ,12 <i>bS</i> )-3-[(cyclopropanecarbonyl)oxy]-6,12-dihydroxy-4,6 <i>a</i> ,12 <i>b</i> -trimethyl-11-oxo-9-(pyridin-3-yl)-1,3,4,4 <i>a</i> ,5,6,6 <i>a</i> ,12,12 <i>a</i> ,12 <i>b</i> -decahydro-2 <i>H</i> ,11 <i>H</i> -benzo[ <i>f</i> ]pyrano[4,3- <i>b</i> ][1]benzopyran-4-yl)methyl cyclopropanecarboxylate
IUPAC name	{{(3 <i>S</i> ,4 <i>R</i> ,4 <i>aR</i> ,6 <i>S</i> ,6 <i>aS</i> ,12 <i>R</i> ,12 <i>aS</i> ,12 <i>bS</i> )-3-[(cyclopropylcarbonyl)oxy]-1,3,4,4 <i>a</i> ,5,6,6 <i>a</i> ,12,12 <i>a</i> ,12 <i>b</i> -decahydro-6,12-dihydroxy-4,6 <i>a</i> ,12 <i>b</i> -trimethyl-11-oxo-9-(3-pyridyl)-2 <i>H</i> ,11 <i>H</i> -benzo[ <i>f</i> ]pyrano[4,3- <i>b</i> ]chromen-4-yl)methyl cyclopropanecarboxylate
CAS name	[(3 <i>S</i> ,4 <i>R</i> ,4 <i>aR</i> ,6 <i>S</i> ,6 <i>aS</i> ,12 <i>R</i> ,12 <i>aS</i> ,12 <i>bS</i> )-3-[(cyclopropylcarbonyl)oxy]-1,3,4,4 <i>a</i> ,5,6,6 <i>a</i> ,12,12 <i>a</i> ,12 <i>b</i> -decahydro-6,12-dihydroxy-4,6 <i>a</i> ,12 <i>b</i> -trimethyl-11-oxo-9-(3-pyridinyl)-2 <i>H</i> ,11 <i>H</i> -naphtho[2,1- <i>b</i> ]pyrano[3,4- <i>e</i> ]pyran-4-yl)methyl cyclopropanecarboxylate

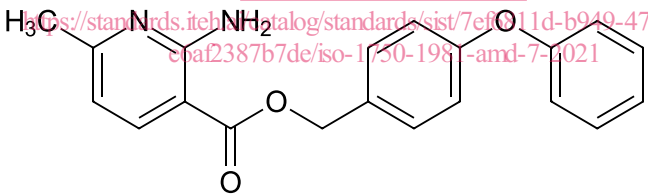
<b>Common name</b>	<b>afidopyropen</b>
<b>Structure</b>	
<b>CAS Reg. No.</b>	915972-17-7
<b>Formula</b>	C <sub>33</sub> H <sub>39</sub> NO <sub>9</sub>
<b>Use</b>	insecticide
<b>InChIKey</b>	LRZWFURXIMFONG-HRSIRGMGSA-N
<b>InChI</b>	InChI=1S/C33H39NO9/c1-31-11-10-24(42-29(38)18-8-9-18)32(2,16-40-28(37)17-6-7-17)22(31)14-23(35)33(3)27(31)26(36)25-21(43-33)13-20(41-30(25)39)19-5-4-12-34-15-19/h4-5,12-13,15,17-18,22-24,26-27,35-36H,6-11,14,16H2,1-3H3/t22-,23+,24+,26+,27-,31+,32+,33-/m1/s1

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<b>Common name</b>	<b>ametotradin</b> (standards.itech.ai)
<b>IUPAC PIN</b>	5-ethyl-6-octyl[1,2,4]triazolo[1,5-a]pyrimidin-7-amine
<b>IUPAC name</b>	5-ethyl-6-octyl[1,2,4]triazolo[1,5-a]pyrimidin-7-amine
<b>CAS name</b>	5-ethyl-6-octyl[1,2,4]triazolo[1,5-a]pyrimidin-7-amine
<b>Structure</b>	
<b>CAS Reg. No.</b>	865318-97-4
<b>Formula</b>	C <sub>15</sub> H <sub>25</sub> N <sub>5</sub>
<b>Use</b>	fungicide
<b>InChIKey</b>	GKGQIOFASHYUJZ-UHFFFAOYSA-N
<b>InChI</b>	InChI=1S/C15H25N5/c1-3-5-6-7-8-9-10-12-13(4-2)19-15-17-11-18-20(15)14(12)16/h11H,3-10,16H2,1-2H3

<b>Common name</b>	<b>aminocyclopyrachlor</b>
<b>IUPAC PIN</b>	6-amino-5-chloro-2-cyclopropylpyrimidine-4-carboxylic acid
<b>IUPAC name</b>	6-amino-5-chloro-2-cyclopropylpyrimidine-4-carboxylic acid
<b>CAS name</b>	6-amino-5-chloro-2-cyclopropyl-4-pyrimidinecarboxylic acid

<b>Common name</b>	<b>aminocyclopyrachlor</b>
Structure	
CAS Reg. No.	858956-08-8
Formula	C <sub>8</sub> H <sub>8</sub> ClN <sub>3</sub> O <sub>2</sub>
Use	herbicide
InChIKey	KWAIHLIXESXTJL-UHFFFAOYSA-N
InChI	InChI=1S/C8H8ClN3O2/c9-4-5(8(13)14)11-7(3-1-2-3)12-6(4)10/h3H,1-2H2,(H,13,14)(H2,10,11,12)
Notes	It should be stated which ester or salt is present, for example aminocyclopyrachlor-methyl [858954-83-3] or aminocyclopyrachlor-potassium [858956-35-1].

<b>Common name</b>	<b>aminopyrifen</b>
IUPAC PIN	(4-phenoxyphenyl)methyl 2-amino-6-methylpyridine-3-carboxylate
IUPAC name	4-phenoxybenzyl 2-amino-6-methylnicotinate
CAS name	(4-phenoxyphenyl)methyl 2-amino-6-methyl-3-pyridinecarboxylate
Structure	
CAS Reg. No.	1531626-08-0
Formula	C <sub>20</sub> H <sub>18</sub> N <sub>2</sub> O <sub>3</sub>
Use	fungicide
InChIKey	PWWPULQZEAPTTB-UHFFFAOYSA-N
InChI	InChI=1S/C20H18N2O3/c1-14-7-12-18(19(21)22-14)20(23)24-13-15-8-10-17(11-9-15)25-16-5-3-2-4-6-16/h2-12H,13H2,1H3,(H2,21,22)

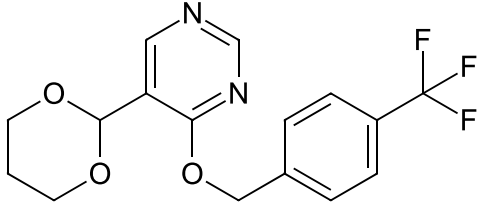
<b>Common name</b>	<b>beflubutamid-M</b>
IUPAC PIN	(2 <i>S</i> )- <i>N</i> -benzyl-2-[4-fluoro-3-(trifluoromethyl)phenoxy]butanamide
IUPAC name	( <i>S</i> )- <i>N</i> -benzyl-2-[( $\alpha,\alpha,\alpha,4$ -tetrafluoro- <i>m</i> -tolyl)oxy]butyramide
CAS name	(2 <i>S</i> )-2-[4-fluoro-3-(trifluoromethyl)phenoxy]- <i>N</i> -(phenylmethyl)butanamide

<b>Common name</b>	<b>beflubutamid-M</b>
<b>Structure</b>	
<b>CAS Reg. No.</b>	113614-09-8
<b>Formula</b>	C <sub>18</sub> H <sub>17</sub> F <sub>4</sub> NO <sub>2</sub>
<b>Use</b>	herbicide
<b>InChIKey</b>	FFQPZWRNXXKPNPX-INIZCTEOSA-N
<b>InChI</b>	InChI=1S/C18H17F4NO2/c1-2-16(17(24)23-11-12-6-4-3-5-7-12)25-13-8-9-15(19)14(10-13)18(20,21)22/h3-10,16H,2,11H2,1H3,(H,23,24)/t16-/m0/s1
<b>Notes</b>	The unresolved enantiomeric mixture has the common name beflubutamid [113614-08-7].

<b>Common name</b>	<b>benzovindiflupyr</b>
<b>IUPAC PIN</b>	<i>rac-N-[(1R,4S)-9-(dichloromethylidene)-1,2,3,4-tetrahydro-1,4-methanonaphthalen-5-yl]-3-(difluoromethyl)-1-methyl-1H-pyrazole-4-carboxamide</i>
<b>IUPAC name</b>	<i>N-[(1RS,4SR)-9-(dichloromethylene)-1,2,3,4-tetrahydro-1,4-methanonaphthalen-5-yl]-3-(difluoromethyl)-1-methyl-1H-pyrazole-4-carboxamide</i>
<b>CAS name</b>	<i>N-[9-(dichloromethylene)-1,2,3,4-tetrahydro-1,4-methanonaphthalen-5-yl]-3-(difluoromethyl)-1-methyl-1H-pyrazole-4-carboxamide</i>
<b>Structure</b>	
<b>CAS Reg. No.</b>	1072957-71-1
<b>Formula</b>	C <sub>18</sub> H <sub>15</sub> Cl <sub>2</sub> F <sub>2</sub> N <sub>3</sub> O



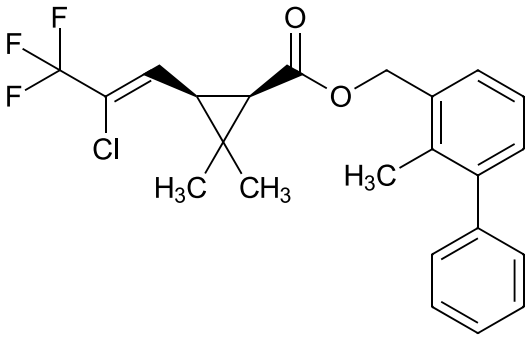
<b>Common name</b>	<b>benzovindiflupyr</b>
Use	fungicide
InChIKey	CCCGEKHKPTUJ-UHFFFAOYSA-N
InChI	InChI=1S/C18H15Cl2F2N3O/c1-25-7-11(15(24-25)17(21)22)18(26)23-12-4-2-3-8-9-5-6-10(13(8)12)14(9)16(19)20/h2-4,7,9-10,17H,5-6H2,1H3,(H,23,26)

<b>Common name</b>	<b>benzpyrimoxan</b>
IUPAC PIN	5-(1,3-dioxan-2-yl)-4-[[4-(trifluoromethyl)phenyl]methoxy]pyrimidine
IUPAC name	5-(1,3-dioxan-2-yl)pyrimidin-4-yl 4-(trifluoromethyl)benzyl ether
CAS name	5-(1,3-dioxan-2-yl)-4-[[4-(trifluoromethyl)phenyl]methoxy]pyrimidine
Structure	
CAS Reg. No.	1449021-97-9
Formula	C <sub>16</sub> H <sub>15</sub> F <sub>3</sub> N <sub>2</sub> O <sub>3</sub>
Use	insecticide
InChIKey	ZYXYTGQFPZEUFX-UHFFFAOYSA-N
InChI	InChI=1S/C16H15F3N2O3/c17-16(18,19)12-4-2-11(3-5-12)9-24-14-13(8-20-10-21-14)15-22-6-1-7-23-15/h2-5,8,10,15H,1,6-7,9H2

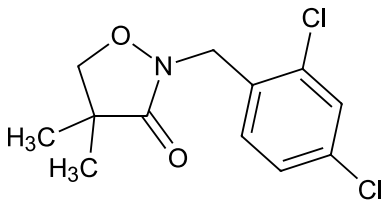
<b>Common name</b>	<b>bicyclopyrone</b>
IUPAC PIN	<i>rac</i> -(1 <i>R</i> ,5 <i>R</i> )-4-hydroxy-3-{2-[(2-methoxyethoxy)methyl]-6-(trifluoromethyl)(pyridine-3-carbonyl)}bicyclo[3.2.1]oct-3-en-2-one
IUPAC name	(1 <i>RS</i> ,5 <i>SR</i> )-4-hydroxy-3-({2-[(2-methoxyethoxy)methyl]-6-(trifluoromethyl)-3-pyridyl}carbonyl)bicyclo[3.2.1]oct-3-en-2-one
CAS name	4-hydroxy-3-[[2-[(2-methoxyethoxy)methyl]-6-(trifluoromethyl)-3-pyridinyl]carbonyl]bicyclo[3.2.1]oct-3-en-2-one

<b>Common name</b>	<b>bicyclopyrone</b>
<b>Structure</b>	
<b>CAS Reg. No.</b>	352010-68-5
<b>Formula</b>	C <sub>19</sub> H <sub>20</sub> F <sub>3</sub> NO <sub>5</sub>
<b>Use</b>	herbicide
<b>InChIKey</b>	HUYBEDCQLAEVPD UHFFFAOYSA-N
<b>InChI</b>	InChI=1S/C19H20F3NO5/c1-27-6-7-28-9-13-12(4-5-14(23-13)19(20,21)22)18(26)15-16(24)10-2-3-11(8-10)17(15)25/h4-5,10-11,24H,2-3,6-9H2,1H3

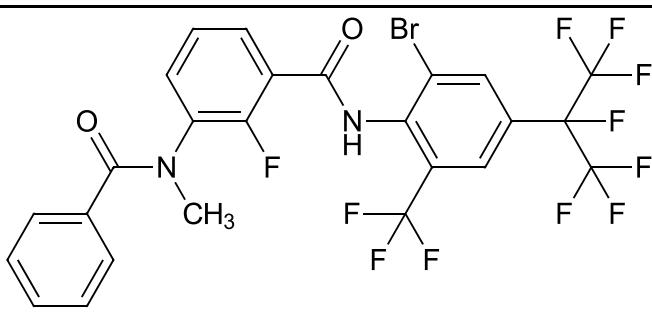
<b>Common name</b>	<b>kappa-bifenthrin</b>
<b>IUPAC PIN</b>	(2-methyl[1,1'-biphenyl]-3-yl)methyl (1 <i>R</i> ,3 <i>R</i> )-3-[(1 <i>Z</i> )-2-chloro-3,3,3-trifluoroprop-1-en-1-yl]-2,2-dimethylcyclopropane-1-carboxylate
<b>IUPAC name</b>	(2-methylbiphenyl-3-yl)methyl (1 <i>R</i> ,3 <i>R</i> )-3-[( <i>Z</i> )-2-chloro-3,3,3-trifluoroprop-1-enyl]-2,2-dimethylcyclopropanecarboxylate Rothamsted-style stereodescriptors: (2-methylbiphenyl-3-yl)methyl (1 <i>R</i> )- <i>cis</i> -3-[( <i>Z</i> )-2-chloro-3,3,3-trifluoroprop-1-enyl]-2,2-dimethylcyclopropanecarboxylate
<b>CAS name</b>	(2-methyl[1,1'-biphenyl]-3-yl)methyl (1 <i>R</i> ,3 <i>R</i> )-3-[(1 <i>Z</i> )-2-chloro-3,3,3-trifluoro-1-propen-1-yl]-2,2-dimethylcyclopropanecarboxylate

<b>Common name</b>	<b>kappa-bifenthrin</b>
<b>Structure</b>	
<b>CAS Reg. No.</b>	439680-76-9
<b>Formula</b>	C <sub>23</sub> H <sub>22</sub> ClF <sub>3</sub> O <sub>2</sub>
<b>Use</b>	insecticide
<b>InChIKey</b>	OMFRMAHOOUJSGP-IRHGGOMRSA-N
<b>InChI</b>	InChI=1S/C23H22ClF3O2/c1-14-16(10-7-11-17(14)15-8-5-4-6-9-15)13-29-21(28)20-18(22(20,2)3)12-19(24)23(25,26)27/h4-12,18,20H,13H2,1-3H3/b19-12-/t18-,20-/m0/s1
<b>Notes</b>	This is one of the 2 isomers of bifenthrin [82657-04-3].

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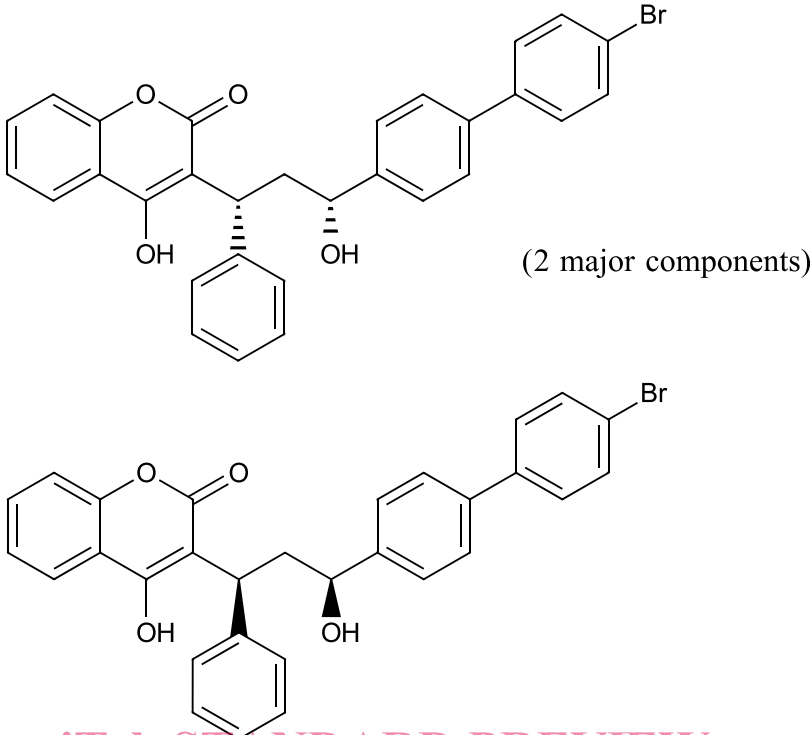
<b>Common name</b>	<b>bixlozone</b> (standards.iteh.ai)
<b>IUPAC PIN</b>	2-[(2,4-dichlorophenyl)methyl]-4,4-dimethyl-1,2-oxazolidin-3-one
<b>IUPAC name</b>	2-[(2,4-dichlorobenzyl)-4,4-dimethylisoxazolidin-3-one
<b>CAS name</b>	2-[(2,4-dichlorophenyl)methyl]-4,4-dimethyl-3-isoxazolidinone
<b>Structure</b>	
<b>CAS Reg. No.</b>	81777-95-9
<b>Formula</b>	C <sub>12</sub> H <sub>13</sub> Cl <sub>2</sub> NO <sub>2</sub>
<b>Use</b>	herbicide
<b>InChIKey</b>	FHUKASKVKWLCY-UHFFFAOYSA-N
<b>InChI</b>	InChI=1S/C12H13Cl2NO2/c1-12(2)7-17-15(11(12)16)6-8-3-4-9(13)5-10(8)14/h3-5H,6-7H2,1-2H3

<b>Common name</b>	<b>broflanilide</b>
<b>IUPAC PIN</b>	N-[2-bromo-4-(1,1,1,2,3,3,3-heptafluoropropan-2-yl)-6-(trifluoromethyl)phenyl]-2-fluoro-3-(N-methylbenzamido)benzamide
<b>IUPAC name</b>	2'-bromo-2-fluoro-3-(N-methylbenzamido)-4'-[1,2,2,2-tetrafluoro-1-(trifluoromethyl)ethyl]-6'-(trifluoromethyl)benzanilide

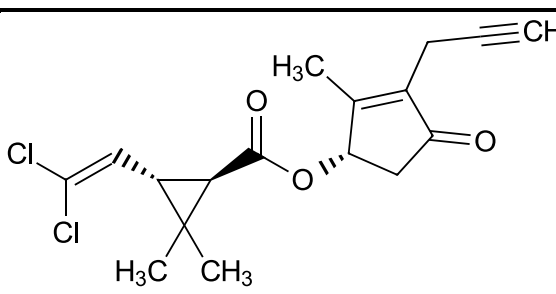
Common name	broflanilide
CAS name	3-(benzoylmethylamino)- <i>N</i> -[2-bromo-4-[1,2,2,2-tetrafluoro-1-(trifluoromethyl)ethyl]-6-(trifluoromethyl)phenyl]-2-fluorobenzamide
Structure	
CAS Reg. No.	1207727-04-5
Formula	C <sub>25</sub> H <sub>14</sub> BrF <sub>11</sub> N <sub>2</sub> O <sub>2</sub>
Use	insecticide
InChIKey	QSLZKWPYTWWEHC-UHFFFAOYSA-N
InChI	InChI=1S/C25H14BrF11N2O2/c1-39(21(41)12-6-3-2-4-7-12)17-9-5-8-14(18(17)27)20(40)38-19-15(23(29,30)31)10-13(11-16(19)26)22(28,24(32,33)34)25(35,36)37/h2-11H,1H3,(H,38,40)

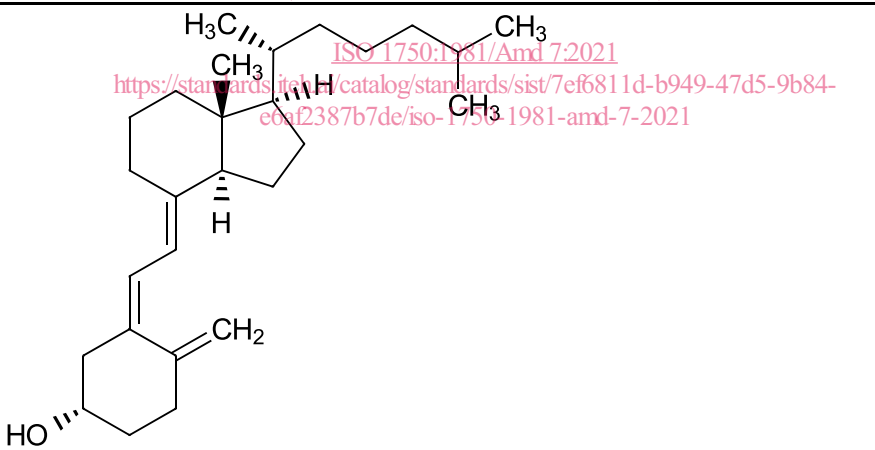
iTech STANDARD PREVIEW

Common name	alpha-bromadiolone (standards.itech.ai)
IUPAC PIN	mixture of 80–100% <i>rac</i> -3-[(1 <i>R</i> ,3 <i>R</i> )-3-(4'-bromo[1,1'-biphenyl]-4-yl)-3-hydroxy-1-phenylpropyl]-4-hydroxy-2 <i>H</i> -1-benzopyran-2-one and 20–0% of the <i>rac</i> -(1 <i>R</i> ,3 <i>S</i> )-isomers
IUPAC name	mixture of 80–100% 3-[(1 <i>RS</i> ,3 <i>RS</i> )-3-(4'-bromobiphenyl-4-yl)-3-hydroxy-1-phenylpropyl]-4-hydroxycoumarin and 20–0% of the (1 <i>RS</i> ,3 <i>SR</i> )-isomers
CAS name	3-[3-(4'-bromo[1,1'-biphenyl]-4-yl)-3-hydroxy-1-phenylpropyl]-4-hydroxy-2 <i>H</i> -1-benzopyran-2-one

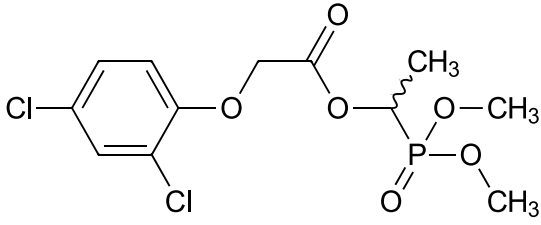
<b>Common name</b>	<b>alpha-bromadiolone</b>
Structure	 <p>(2 major components)</p>
CAS Reg. No.	28772-56-7
Formula	C <sub>30</sub> H <sub>23</sub> BrO <sub>4</sub>
Use	rodenticide <a href="https://standards.iteh.ai/catalog/standards/sist/7cf6811d-b949-47d5-9f84-c6a12387b7de/iso-1750-1981-amd-7-2021">ISO 1750:1981/Amd 7:2021</a>
InChIKey	LQAGAOMQXARIFO-CPBLEMPESA-N (major component)
InChI	InChI=1S/2C30H23BrO4/c2*31-23-16-14-20(15-17-23)19-10-12-22(13-11-19)26(32)18-25(21-6-2-1-3-7-21)28-29(33)24-8-4-5-9-27(24)35-30(28)34/h2*1-17,25-26,32-33H,18H2/t2*25-,26-/m10/s1 (major component)
Notes	The mixture of 80–100% (1 <i>R</i> ,3 <i>S</i> )-isomers and 20–0% (1 <i>R</i> ,3 <i>R</i> )-isomers has the ISO common name <a href="#">bromadiolone</a> .

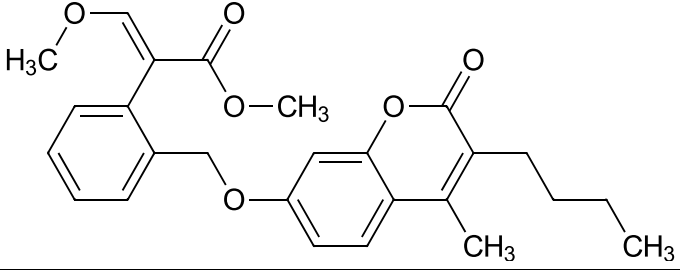
<b>Common name</b>	<b>chloroprallethrin</b>
IUPAC PIN	(1 <i>S</i> )-2-methyl-4-oxo-3-(prop-2-yn-1-yl)cyclopent-2-en-1-yl (1 <i>R</i> ,3 <i>S</i> )-3-(2,2-dichloroethenyl)-2,2-dimethylcyclopropane-1-carboxylate
IUPAC name	(1 <i>S</i> )-2-methyl-4-oxo-3-(prop-2-ynyl)cyclopent-2-enyl (1 <i>R</i> ,3 <i>S</i> )-3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate Rothamsted-style stereodescriptors: (1 <i>S</i> )-2-methyl-4-oxo-3-(prop-2-ynyl)cyclopent-2-enyl (1 <i>R</i> )- <i>trans</i> -3-(2,2-dichlorovinyl)-2,2-dimethylcyclopropanecarboxylate
CAS name	(1 <i>S</i> )-2-methyl-4-oxo-3-(2-propyn-1-yl)-2-cyclopenten-1-yl (1 <i>R</i> ,3 <i>S</i> )-3-(2,2-dichloroethenyl)-2,2-dimethylcyclopropanecarboxylate

<b>Common name</b>	<b>chloroprallethrin</b>
<b>Structure</b>	
<b>CAS Reg. No.</b>	250346-55-5
<b>Formula</b>	C <sub>17</sub> H <sub>18</sub> Cl <sub>2</sub> O <sub>3</sub>
<b>Use</b>	insecticide
<b>InChIKey</b>	RQNZCZHWWQMZER-ZLDLUXBVSA-N
<b>InChI</b>	InChI=1S/C17H18Cl2O3/c1-5-6-10-9(2)13(8-12(10)20)22-16(21)15-11(7-14(18)19)17(15,3)4/h1,7,11,13,15H,6,8H2,2-4H3/t11-,13+,15+/m1/s1

<b>Common name</b>	<b>cholecalciferol</b>
<b>IUPAC name</b>	(5Z,7E)-(3S)-9,10-secocholesta-5,7,10(19)-trien-3-ol
<b>CAS name</b>	(1S,3Z)-3-[(2E)-2-[(1R,3aS,7aR)-1-[(1R)-1,5-dimethylhexyl]octahydro-7a-methyl-4H-inden-4-ylidene]ethylidene]-4-methylenecyclohexanol
<b>Structure</b>	
<b>CAS Reg. No.</b>	67-97-0
<b>Formula</b>	C <sub>27</sub> H <sub>44</sub> O
<b>Use</b>	rodenticide
<b>InChIKey</b>	QYSXJUF SXHHAJI-YRZJJWOYSA-N
<b>InChI</b>	InChI=1S/C27H44O/c1-19(2)8-6-9-21(4)25-15-16-26-22(10-7-17-27(25,26)5)12-13-23-18-24(28)14-11-20(23)3/h12-13,19,21,24-26,28H,3,6-11,14-18H2,1-2,4-5H3/b22-12+,23-13-/t21-,24+,25-,26+,27-/m1/s1
<b>Notes</b>	In China, the name "vitamin D3" (维生素 D3) is used.

<b>Common name</b>	<b>clacyfos</b>
<b>IUPAC PIN</b>	<i>rac</i> -(1R)-1-(dimethoxyphosphoryl)ethyl (2,4-dichlorophenoxy)acetate

<b>Common name</b>	<b>clacyfos</b>
<b>IUPAC name</b>	(1 <i>RS</i> )-1-(dimethoxyphosphinoyl)ethyl (2,4-dichlorophenoxy)acetate
<b>CAS name</b>	1-(dimethoxyphosphinyl)ethyl 2-(2,4-dichlorophenoxy)acetate
<b>Structure</b>	
<b>CAS Reg. No.</b>	215655-76-8
<b>Formula</b>	C <sub>12</sub> H <sub>15</sub> Cl <sub>2</sub> O <sub>6</sub> P
<b>Use</b>	herbicide
<b>InChIKey</b>	UUHXXNQVWVFJLW-UHFFFAOYSA-N
<b>InChI</b>	InChI=1S/C12H15Cl2O6P/c1-8(21(16,17-2)18-3)20-12(15)7-19-11-5-4-9(13)6-10(11)14/h4-6,8H,7H2,1-3H3
<b>Notes</b>	The parent acid has the common name 2,4-D [94-75-7].

<b>Common name</b>	<b>coumoxystrobin</b>
<b>IUPAC PIN</b>	methyl (2 <i>E</i> )-2-(2-[[[(3-butyl-4-methyl-2-oxo-2 <i>H</i> -1-benzopyran-7-yl)oxy]methyl}phenyl]-3-methoxyprop-2-enoate
<b>IUPAC name</b>	methyl (2 <i>E</i> )-2-(2-[[[(3-butyl-4-methyl-2-oxo-2 <i>H</i> -chromen-7-yl)oxy]methyl}phenyl]-3-methoxyacrylate
<b>CAS name</b>	methyl ( $\alpha$ <i>E</i> )-2-[[[(3-butyl-4-methyl-2-oxo-2 <i>H</i> -1-benzopyran-7-yl)oxy]methyl]- $\alpha$ -(methoxymethylene)benzeneacetate
<b>Structure</b>	
<b>CAS Reg. No.</b>	850881-70-8
<b>Formula</b>	C <sub>26</sub> H <sub>28</sub> O <sub>6</sub>
<b>Use</b>	fungicide
<b>InChIKey</b>	CWVRPJSBNHNJSI-XQNSMLJCSA-N
<b>InChI</b>	InChI=1S/C26H28O6/c1-5-6-10-21-17(2)20-13-12-19(14-24(20)32-26(21)28)31-15-18-9-7-8-11-22(18)23(16-29-3)25(27)30-4/h7-9,11-14,16H,5-6,10,15H2,1-4H3/b23-16+

<b>Common name</b>	<b>cyantraniliprole</b>
<b>IUPAC PIN</b>	3-bromo-1-(3-chloropyridin-2-yl)- <i>N</i> -[4-cyano-2-methyl-6-(methylcarbamoyl)phenyl]-1 <i>H</i> -pyrazole-5-carboxamide